





Measurement of the ^{87}Rb *D*-line vector tune-out wavelengthA. J. Fallon , E. R. Moan , E. A. Larson , and C. A. Sackett ^{*}*Department of Physics, University of Virginia, Charlottesville, Virginia 22904, USA*

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We report a precision measurement of a tune-out wavelength for ^{87}Rb using circularly polarized light. A tune-out wavelength characterizes a zero in the electric polarizability of the atom. For circularly polarized light, the total polarizability depends on both the scalar and vector polarizability components. This shifts the location of the tune-out wavelength and makes it sensitive to different combinations of atomic dipole matrix elements than the scalar polarizability alone. Using σ_- polarized light with a purity of 0.9931(1), we observe a tune-out wavelength of 785.1522(3) nm, which agrees with theoretical expectations when small contributions from the core electrons and off-resonant valence states are taken into account.

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A tune-out wavelength describes a light frequency at which an atom or molecule in a given state experiences a zero-energy shift from an optical field, due to cancellation of the positive and negative shifts from nearby blue- and red-detuned electronic transitions [1,2]. Tune-out wavelengths find applications in experiments involving multiple species, where it can be useful to apply an energy shift to one species without affecting another [3,4]. Tune-out wavelength measurements are also useful in their own right because they provide information about the dipole matrix elements of the target particle that may not otherwise be easily accessible. Knowledge of dipole matrix elements is important for many reasons, including the interpretation of parity violation experiments, accurate estimation of blackbody radiation shifts in atomic clocks [5,6], and as benchmarks for atomic theory calculations. These benefits have prompted a series of precise tune-out-wavelength measurements in alkali-metal and other atoms [3,7–17]. These experiments have mainly focused on zeros of the scalar electric polarizability of the atoms. However, additional information can be obtained from the vector character of the polarizability, which is exhibited through a dependence on the optical polarization of the applied light [12,13,17]. We here explore this polarization dependence through a precise measurement of a vector tune-out wavelength.

Vector tune-out measurements are useful both for trapping applications and for fundamental atomic physics. For applications, they add flexibility by allowing the tune-out value to be adjusted [4]. For instance, the *D*-line scalar tune-out wavelength for Rb is fixed at 790.032 nm, but by adjusting the light polarization, the tune-out wavelength can be set anywhere between 785.112 nm and the *D*₁ line at 794.978 nm. This flexibility can make it easier to satisfy other experimental requirements. It can also be useful that vector fields cause the tune-out wavelength to depend on the magnetic sublevel of the particle [12,13,18].

In regards to fundamental physics, precise measurements of the polarization allows resolution of contributions to the atomic polarizability from different angular momentum states. For instance, interpreting alkali-metal atom parity violation amplitudes in terms of nuclear physics parameters requires knowledge of the $nS_{1/2} \leftrightarrow nP_{1/2}$ dipole matrix elements [19,20]. Vector tune-out measurements can allow the $P_{1/2}$ matrix elements to be constrained separately from the $P_{3/2}$ matrix elements, whereas a purely scalar measurement depends jointly on both $P_{1/2}$ and $P_{3/2}$ elements [21]. Although parity violation experiments do not generally focus on Rb, we expect that precise measurements in Rb could help improve theoretical calculation methods which would be applicable to similar atoms such as Cs or Fr.

The theoretical framework for vector tune-out wavelengths is well understood [22–25]. However, making a precise comparison between theory and measurement requires careful control of both the light polarization and the alignment of the laser beam axis to the quantizing magnetic field. The measurement reported here has a wavelength precision of order 1 pm, and agrees to this level with theoretical expectations. At this precision, the measurement is sensitive to small effects including the polarizability of the ionic core and contributions from far off-resonant valence states. With realistic improvements in precision and by combining tune-out measurements for different states, the technique could provide constraints on important dipole matrix elements and yield accuracies better than the best current theoretical uncertainties.

The energy shift of a particle in an optical field \mathcal{E} can be expressed as

$$U = -\frac{1}{2}\alpha\langle\mathcal{E}^2\rangle = -\frac{1}{2\epsilon_0 c}\alpha I, \quad (1)$$

where α is the electric polarizability, I is the light intensity, c is the speed of light, and ϵ_0 is the electric constant. (The next-order term in the expansion is an estimated 10^7 times smaller for the maximum intensity used here.) The polarizability depends on both the frequency ω and the polarization state $\hat{\epsilon}$ of the light. For an atom in the ground hyperfine state

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$|n, J, F, m\rangle$, a spherical tensor decomposition gives [23,24]

$$\alpha = \alpha^{(0)} - \alpha^{(1)} S_3 \hat{k} \cdot \hat{b} \frac{m}{2F} + \alpha^{(2)} \left(\frac{3|\hat{\epsilon} \cdot \hat{b}|^2 - 1}{2} \right) \frac{3m^2 - F(F+1)}{F(2F-1)}, \quad (2)$$

where the $\alpha^{(i)}$ parameters are the scalar, vector, and tensor polarizability components, for $i = 0, 1$, and 2 , respectively. Here, the light field is taken as a plane wave propagating in direction \hat{k} , with complex polarization vector $\hat{\epsilon}$. The atomic states are defined relative to a magnetic field pointing in direction \hat{b} . The parameter $S_3 = i(\hat{\epsilon}^* \times \hat{\epsilon}) \cdot \hat{k}$ is the fourth Stokes parameter for the light, with $S_3 = \pm 1$ for σ_{\pm} circularly polarized light. Our measurements use the $5S_{1/2}$ ground state of ^{87}Rb , with $F = m = 2$.

In the case of an alkali-metal atom, the polarizability components can be separated into a contribution from the valence electron, a contribution from the core electrons, and a term reflecting interactions between the valence electron and the core. The valence contribution can be calculated using perturbation theory as a sum over excited P states $|n', J', F', m'\rangle$. We measure the tune-out wavelength near the $5P_{1/2}$ and $5P_{3/2}$ states, so it is necessary to account for the hyperfine splittings of these states in order to achieve sufficient precision. For higher-lying states, the hyperfine shifts can be neglected since the polarizability contributions are much smaller. We therefore express the polarizability as

$$\alpha^{(i)} = \alpha_c^{(i)} + \alpha_{cv}^{(i)} + \alpha_{5P}^{(i)} + \alpha_{v'}^{(i)}, \quad (3)$$

where α_c denotes the core contribution, α_{cv} the core-valence interaction, α_{5P} the contribution from the $5P$ states, and $\alpha_{v'}$ the contribution from other valence states. Furthermore, the core contribution has only a scalar component $i = 0$, since the core is spherically symmetric. We also neglect the tensor components $\alpha_{cv}^{(2)}$ and $\alpha_{v'}^{(2)}$ since they are of order 10^{-5} a.u. or smaller. The remaining valence contributions are then [24]

$$\alpha_{5P}^{(0)} = \frac{2}{\hbar} \frac{1}{\sqrt{3(F+1)}} \sum_{J',F'} \frac{|d'|^2 \omega'}{\omega'^2 - \omega^2} \begin{Bmatrix} 1 & 0 & 1 \\ F & F' & F \end{Bmatrix} C', \quad (4)$$

$$\alpha_{5P}^{(1)} = \frac{2}{\hbar} \sqrt{\frac{2F}{(F+1)(2F+1)}} \times \sum_{J',F'} \frac{|d'|^2 \omega'}{\omega'^2 - \omega^2} \begin{Bmatrix} 1 & 1 & 1 \\ F & F' & F \end{Bmatrix} C', \quad (5)$$

$$\alpha_{5P}^{(2)} = \frac{2}{\hbar} \sqrt{\frac{2F(2F-1)}{3(F+1)(2F+1)(2F+3)}} \times \sum_{J',F'} \frac{|d'|^2 \omega'}{\omega'^2 - \omega^2} \begin{Bmatrix} 1 & 2 & 1 \\ F & F' & F \end{Bmatrix} C', \quad (6)$$

$$\alpha_{v'}^{(0)} = \frac{1}{3\hbar} \sum_{n',J'} \frac{|d'|^2 \omega'}{\omega'^2 - \omega^2}, \quad (7)$$

$$\alpha_{v'}^{(1)} = \frac{1}{3\hbar} \sum_{n',J'} \frac{|d'|^2 \omega'}{\omega'^2 - \omega^2} \left(3J' - \frac{7}{2} \right). \quad (8)$$

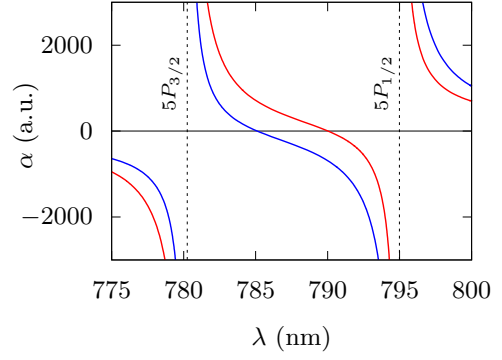


FIG. 1. Electric polarizability of ^{87}Rb in the $F = 2$, $m = 2$ ground state, as a function of optical wavelength. The red curve shows the case of linearly polarized light, and exhibits a tune-out wavelength near 790 nm. The blue curve shows the case of σ_- polarized light, with a tune-out wavelength near 785 nm.

Here, ω is the light frequency, and ω' is the transition frequency from $|5S_{1/2}, F\rangle$ to $|n'P_{J'}, F'\rangle$. We neglect Zeeman shifts since they are small (about 30 MHz) compared to our measurement precision (about 150 MHz). The arrays in braces are Wigner 6- j symbols. The reduced matrix elements are $d' \equiv \langle 5S_{1/2} || d || n'P_{J'} \rangle$, and in the α_{5P} terms we use

$$C' = (-1)^{F+F'+1} (2F+1)(2F'+1) \begin{Bmatrix} F & 1 & F' \\ J' & I & J \end{Bmatrix}^2, \quad (9)$$

with nuclear angular momentum $I = 3/2$.

We take $\alpha_c^{(0)} = 9.116(9)$ from Ref. [26], and $\alpha_{cv}^{(0)} = -0.37(4)$ and $\alpha_{cv}^{(1)} = -0.04(4)$ from Ref. [21]. For the $5P$ states we use $d_{5P_{1/2}} = 4.234(2)$ and the ratio $d_{5P_{3/2}}/d_{5P_{1/2}} = 1.99217(3)$ from Ref. [11]. For higher-lying valence states we use the matrix elements tabulated in Ref. [11]. With these values, we can calculate the net polarizability α for given values of the experimental parameters S_3 , $\hat{k} \cdot \hat{b}$, and $\hat{\epsilon} \cdot \hat{b}$. Figure 1 shows how α varies with wavelength for the cases of linear and σ_- polarized light. The tune-out wavelength is located where $\alpha = 0$. Table I lists the various contributions to α at the tune-out wavelength for σ_- polarized light.

TABLE I. Contributions to the total polarizability α at the tune-out wavelength $\lambda = 785.112$ nm for ideal σ_- polarized light with $S_3 = \hat{k} \cdot \hat{b} = 1$, $\hat{\epsilon} \cdot \hat{b} = 0$. Values in parentheses show the estimated errors. In the case of the α_{5P} contributions, the first parentheses show the uncertainty arising from the uncertainty in the ratio of the $d_{5P_{1/2}}$ to $d_{5P_{3/2}}$ matrix elements. The second parentheses show the uncertainty from the $d_{5P_{1/2}}$ element itself, which is large but correlated among the different components, and therefore has negligible impact on the value of the tune-out wavelength. Uncertainty in the $\alpha_{5P}^{(2)}$ contribution is negligible since the term itself is very small.

Term	Value (a.u.)	Term	Value (a.u.)
$\alpha_{5P}^{(0)}$	12347.7(4)(11.6)	$\alpha_{v'}^{(1)}$	0.2(1)
$\alpha_{5P}^{(1)}$	24716.8(4)(23.3)	$\alpha_c^{(0)}$	9.12(1)
$\alpha_{5P}^{(2)}$	-0.044	$\alpha_{cv}^{(0)}$	-0.37(4)
$\alpha_{v'}^{(0)}$	2.0(1)	$\alpha_{cv}^{(1)}$	-0.04(4)

The experimental apparatus consists of a Bose-Einstein condensate interferometer similar to that of Ref. [11]. We use small condensates of about 10^4 atoms, which provide good phase coherence and well-localized wave packets. The atoms are confined in a weak magnetic trap, with harmonic oscillation frequencies of 5.1, 1.1, and 3.2 Hz along the x , y , and z directions, respectively. The z direction is vertical. The trap uses the time-orbiting potential (TOP) technique, with a bias field of 21.4 G rotating in the xz plane at frequency $\Omega = 2\pi \times 12.8$ kHz. The TOP trap is completed using a linear quadrupole field oscillating in phase with the bias, and also a weaker spherical quadrupole oscillating at 1 kHz.

The interferometer operation is described in Ref. [27]. An off-resonant standing-wave laser along the y axis of the trap applies velocity kicks in units of $v_B = 2\hbar k/m = 11.8$ mm/s via Bragg scattering. The interferometer uses a total of four Bragg pulses. At time $t = 0$, an initial pulse splits the condensate into wave packets moving at $\pm v_B$. At time $t = 10$ ms, the laser is applied again so as to reverse the atoms' motion. The packets then pass through each other with minimal interactions, and at $t = 30$ ms a third laser pulse reverses the motion again. Finally at $t = 40$ ms, the packets are overlapped, the initial splitting pulse is reapplied, and the wave packets are recombined. A fraction N_0/N of the atoms are brought back to rest in the center of the trap, with signal

$$S = \frac{N_0}{N} = \frac{1}{2}[1 + V \cos(\phi + \phi_r)]. \quad (10)$$

Here, the wave packets have developed a phase difference ϕ , the phase of the recombination pulse relative to the initial splitting pulse is ϕ_r , and the visibility is $V = 0.7$. The fraction of atoms at rest is detected by absorption imaging after a short time of flight. We set $\phi_r = \pi/2$ by shifting the frequency of the Bragg laser prior to the final pulse.

A Stark phase shift ϕ is applied by directing a second laser beam, traveling along z , onto one arm of the interferometer. The beam is focused to a waist of about $50 \mu\text{m}$, which is smaller than the maximum wave-packet separation of $240 \mu\text{m}$ and comparable to the wave-packet size of $40 \mu\text{m}$. The Stark beam is derived from an MBR-110 Ti:sapphire laser from Coherent, Inc. This is an improvement over the tapered amplifier used in our previous work [11], since the Ti:sapphire laser is not expected to contain a significant amount of amplified spontaneous emission light at other frequencies. The Stark beam is applied for 20 ms at the start of the interferometer, so that one packet passes through it twice. This leads to a phase $\phi = \int \alpha I dt / (2\epsilon_0 \hbar c)$ which the interferometer detects.

To control the vector portion of the polarizability, we use a pair of acousto-optic modulators to pulse the Stark beam synchronously with the rotating bias field, such that the light is on only when the field points along z . Two modulators are used to provide an extinction ratio better than 60 dB. The Stark beam is aligned to the field by tuning to the D_1 resonance at 795 nm and setting the polarization state to σ_+ . When the laser is optimally aligned to the magnetic field, the $m = 2$ atoms scatter no light since there is no $m = 3$ state in the D_1 hyperfine manifold. Details of this measurement are provided in Ref. [28]. From the residual scattering rate, the alignment error $\delta\theta$ between \hat{k} and \hat{z} is constrained to be less than 16 mrad. In the interferometry experiments, the duration of the pulses is $\tau = 5 \mu\text{s}$, and the angle between \hat{k} and \hat{b} varied

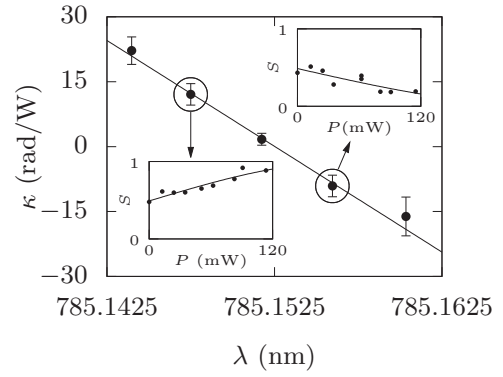


FIG. 2. Tune-out measurement data. For each of the points in the main graph, an interferometric measurement is performed to determine $\kappa = \phi/P$, where ϕ is the interferometer phase and P is the peak power of the Stark beam pulses. The inset graphs show example plots of the interferometer signal vs power, from which κ is determined via a fit to the form of Eq. (10). The scatter of the points illustrates our typical measurement noise, and the error bars for κ are estimated as the offset needed to double the goodness of fit parameter χ^2 . The κ data in the main graph are fit to a line, and the x intercept of 785.1525 nm is taken as the tune-out wavelength value.

during the pulse as the field rotated at frequency Ω . This gives an average value for $\hat{k} \cdot \hat{b}$ of $(2/\Omega\tau) \sin(\Omega\tau/2) \cos\delta\theta = 0.99321(6)$, where the uncertainty reflects the angular misalignment. It is important that the shape and duration of the light pulse are well characterized, since they impact the time average of $\hat{k} \cdot \hat{b}$. We monitor the pulse during the experiment using a fast photodiode, and errors from the nonuniformity and from variations in τ are about an order of magnitude smaller than that from the alignment. The integrated value of $\hat{\epsilon} \cdot \hat{b}$ in the tensor term can also be calculated as $\Omega^2\tau^2/48 = 3.3 \times 10^{-3}$, but this is insignificant because the tensor component $\alpha^{(2)}$ is much smaller than the vector component $\alpha^{(1)}$.

It is also critical to control the light polarization accurately. We set the polarization close to circular using a calcite polarizer, two wave plates, and a Fresnel rhomb, as described in Ref. [28]. The wave plates provide a small correction to the Fresnel rhomb to account for polarization distortions in the vacuum window and other optics. The polarization can be set accurately using again the photon scattering measurements at 795 nm, but the mirrors that direct the beam onto the atoms are slightly chromatic and the polarization is not sufficiently preserved when the Stark laser is scanned to the tune-out wavelength at 785 nm. Instead, we optimize the polarization by setting the laser slightly blue of the tune-out wavelength, and adjusting the wave plates to minimize the interferometer phase ϕ . Since the tune-out wavelength is as blue as possible for σ_- polarized light, this optimizes the polarization at the atoms. The wave-plate angles could be set to an accuracy of about 0.5° , from which we determine $S_3 = 0.99988(12)$.

To perform the measurement, we set the Stark laser to a series of wavelengths near 785 nm, and at each wavelength we vary the pulse power P to scan ϕ and trace out an interference curve. Example data are shown as insets in Fig. 2. We assume $\phi = \kappa P$ and fit the trace data to obtain a value for κ . The main

TABLE II. Sources of error in the tune-out wavelength measurement and calculation. For each contribution, the impact is given both as the uncertainty $\delta\lambda_0$ in the tune-out wavelength and as the uncertainty in the value $\delta\alpha$ of the polarizability at the measured tune-out wavelength. These are related using the calculated derivative $|d\alpha/d\lambda| = 2.527$ a.u./pm. The uncertainty contribution labeled as “atomic parameters” refers to the values in Table I. The division between “Measurement” and “Calculation” is based on the uncertainties affecting the actual tune-out value measured in the experiment, and those affecting what we expect to observe given our experimental setup.

Source	$\delta\lambda_0$ (pm)	$\delta\alpha$ (a.u.)
Measurement total:	0.46	1.2
Statistical	0.43	1.1
Wave meter	0.15	0.4
Calculation total:	0.88	2.2
Atomic parameters	0.35	0.9
Polarization	0.70	1.8
Alignment	0.35	0.9
Pulse length	0.06	0.2
Pulse symmetry	0.06	0.2

graph of Fig. 2 shows that near 785 nm, κ is a linear function of wavelength which crosses zero at 785.1525(5) nm. Here, the uncertainty is primarily from noise in the linear fit, but also includes the 0.15 nm uncertainty in our wave-meter calibration. A second independent measurement yielded a consistent value of 785.1519(4) nm, so we report the average of these results as $\lambda_0 = 785.1522(3)$ nm.

In comparison, using the the experimental estimates for $\hat{k} \cdot \hat{b}$ and S_3 in conjunction with the theoretical values from Table I, we calculate an expected tune-out value of 785.1538(9) nm, which is about 2σ different from our measurement. The optical polarization is the largest source of uncertainty in the calculated value, with the alignment error and atomic parameters contributing about half as much. Table II summarizes the main contributions to the uncertainties of the measurement and calculation. Our result is consistent with that obtained by Wen *et al.* [17], who found $\lambda_0 = 785.146(12)$ nm for σ_- polarized light.

Although the discrepancy between our measurement and calculation is not large enough to be significant, the sign is interesting, since larger-than-estimated errors in the polarization or alignment would result in a measurement redder than expected, whereas our result is bluer. If we assume that the Stark beam parameters are perfect and account only for the rotation of the bias field during the Stark pulses, we would expect a tune-out wavelength of 785.1530(4) nm, still about 2σ redder than observed. This could suggest an inaccuracy in the atomic parameters of Table I, and motivates further investigations.

For atom-trapping applications, the level of precision demonstrated here shows how accurately a vector tune-out application can be implemented. For instance, in the conditions of our experiment, a rubidium atom in the $F = 2$, $m = 1$ state would experience a total polarizability α of 6222 a.u. This can be compared to a residual polarizability $\delta\alpha$ of 2 a.u. for an $F = 2$, $m = 2$ atom. The ratio $\alpha/\delta\alpha = 3 \times 10^3$ indicates

how strongly the $m = 1$ atom can be manipulated before the $m = 2$ atom is affected.

In terms of atomic physics, we see that the precision demonstrated here is already sufficient to distinguish the larger non- $5P$ contributions to the net polarizability. If, for instance, the core contribution $\alpha_c^{(0)} + \alpha_{cv}^{(0)}$ were excluded from the calculation, the expected tune-out wavelength would shift by about 4σ . Our measurement thus tests the theory in a nontrivial way, but with reasonable increases in precision, it could provide a more meaningful comparison. For instance, with a factor of 5 improvement the experiment would be sensitive to the $\alpha_{cv}^{(0)}$ core-valence interaction, which has not previously been experimentally observed. With a factor of 50 improvement, the experimental precision would exceed the theoretical precision in most cases. This would be particularly interesting for the $\alpha_{v'}$ terms, where the theoretical uncertainty is dominated by the contribution from the high n' Rydberg tail. This same contribution is the largest source of uncertainty in the relationship between measured atomic parity violation amplitudes and the weak mixing angle of the standard model [20], so providing a precise benchmark via the polarizability can be expected to help improve the parity violation interpretation.

Improvement by a factor of 50 is experimentally feasible. We have previously demonstrated a scalar tune-out measurement with an uncertainty of 0.035 pm, which was limited primarily by statistics [11]. Improvement to 0.01 pm should involve no new challenges. The vector measurement is more difficult due to the requirement for polarization control, but many of the limitations encountered here could be resolved for atoms confined in an optical trap rather than a TOP trap, since it would then be possible to use a static bias field and a continuous-wave Stark beam. This would allow significantly higher average power to be applied to the atoms, so that the polarization and alignment optimizations could be made more precise. Recent experiments at Los Alamos National Laboratory have demonstrated an optically trapped atom interferometer with performance comparable to that used here [29]. An optical trap may also benefit from being able to use different Zeeman levels. We therefore argue that reaching experimental precision comparable to the theoretical precision is likely achievable.

Looking forward to such experiments, it will be necessary to distinguish the various contributions to α so that, for instance, the Rydberg tail contribution can be isolated from the core-valence interaction. This can be achieved by comparing tune-out measurements near different states, such as the $6P$ states near $\lambda = 420$ nm for Rb. The core and Rydberg contributions have different frequency dependencies, allowing their impact to be resolved [21]. Further, since the $J = 1/2$ and $J = 3/2$ states contribute differently to the scalar and vector components of $\alpha_{v'}$ in Eqs. (7) and (8), these two contributions can be distinguished as well. The parity violation interpretation depends only on the $J = 1/2$ matrix elements. We therefore expect vector measurements to be an important component of this approach.

In summary, we have carried out a precise measurement of a vector tune-out wavelength, for near-circularly-polarized light. We show that the polarization and alignment factors can be controlled with 10 ppm precision, even in the rotating

magnetic field of a TOP trap. The 1 ppm precision that we obtain in the wavelength is similar to that of many scalar tune-out measurements, but the vector character provides both more utility and more information. We believe that this work illustrates the feasibility and utility of precise vector tune-out measurements, and we hope that our results stimulate further improvements to the point that the method becomes useful for

interpreting parity violation and other experiments that rely on atomic dipole matrix elements.

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