Ab initio studies of the double–Gamow-Teller transition and its correlation with neutrinoless double- β decay

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We use chiral interactions and several *ab initio* methods to compute the nuclear matrix elements (NMEs) for ground-state-to-ground-state double Gamow-Teller transitions in a range of isotopes and explore the correlation of these NMEs with those for neutrinoless double beta decay produced by the exchange of a light Majorana neutrino. When all the NMEs of both isospin-conserving and isospin-changing transitions from the *ab initio* calculations are considered, the correlation is strong. For the experimentally relevant isospin-changing transitions by themselves, however, the correlation is weaker and may not be helpful for reducing the uncertainty in the NMEs for neutrinoless double beta decay.

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I. INTRODUCTION

Neutrinoless double- β ($0\nu\beta\beta$) decay is a hypothetical lepton-number-violating process [1] in which two neutrons in a parent nucleus decay into two protons in its daughter nucleus via the emission of only two electrons. The hunt for $0\nu\beta\beta$ decay is of particular importance as its observation would demonstrate the Majorana nature of neutrinos and provide a key ingredient for generating the matter-antimatter asymmetry in the Universe. If $0\nu\beta\beta$ decay is driven by the standard mechanism of exchanging light Majorana neutrinos, then its half-life can also be used to determine the effective neutrino mass $\langle m_{\beta\beta} \rangle = \sum_i U_{ei}^2 m_i$, where m_i are the masses of light neutrinos and U_{ei} are elements of the neutrino-mixing matrix [2]. The $0\nu\beta\beta$ decay rate is governed by a nuclear matrix element (NME) that must be computed. The precise determination of the NMEs for candidate nuclei, which are vital for interpreting and planning the current- [3-6] and nextgeneration ton-scale [7-9] experiments, is thus being pursued energetically by theoretical physicists [10].

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applied to compute the NMEs in nuclei of interest to experiment [11-30], under the assumption that light-neutrino exchange is the dominant decay mechanism. The discrepancy between these predictions is as large as a factor of about three, causing uncertainty at the level of an order of magnitude in the half-life for a given value of the neutrino mass. Resolving this discrepancy has been one of the most significant objectives in the nuclear community [31,32]; see for instance the recent reviews in Refs. [33-36]. Unfortunately, the systematic uncertainty turns out to be difficult to reduce because each model has its phenomenological assumptions and uncontrolled approximations. In recent years, remarkable progress has been achieved in first-principles calculations of nuclear structure and reactions [37]. It has enabled the first wave of multimethod results for $0\nu\beta\beta$ -decay NMEs in light nuclei [38-41], the lightest experimental candidate ⁴⁸Ca [42,43], and in one case even in the heavier candidates ⁷⁶Ge and ⁸²Se [44]. These calculations start with realistic two-nucleon-plus-three-nucleon (NN+3N) interactions from either phenomenological parametrization or chiral effective field theory (EFT). In heavier candidate nuclei, such calculations that include fully controllable uncertainties are still challenging, however [35]. Under these circumstances, it is worthwhile to explore correlations between the $0\nu\beta\beta$ -decay NMEs and other observables. Such correlations may provide model-independent constraints on the NMEs.

A wide range of conventional nuclear models have been

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Recently, Shimizu et al. [45] found a linear correlation between the NMEs of $0\nu\beta\beta$ decay and those that govern the ground-state-to-ground-state double Gamow-Teller (DGT) transition, a double spin-isospin flip excitation mode accessible in high-energy heavy-ion double-charge-exchange (HIDCE) processes [46–49]. The correlation, which appeared in both medium-mass and heavy nuclei in calculations based on the large-scale nuclear shell-model, was attributed to the mainly short-range character of both transitions [50,51]. These studies provide strong support to experimental programs to measure HIDCE reactions, through examples such as $^{12}C(^{18}O, ^{18}Ne)^{12}Be$ [52] and $(^{11}B, ^{11}Li)$ [53] at the Research Center of Nuclear Physics, Osaka University, and others in the NUMEN project at the Laboratori Nazionali del Sud, Istituto Nazionale di Fisica Nucleare [54]. The expectation is that the cross section in HIDCE reactions can place a constraint on the NMEs for $0\nu\beta\beta$ decay if the correlation exists and is universal. Santopinto et al. [55] argued that it is possible to factorize the HIDCE cross section into reaction and nuclear parts. The latter can be further written as a product of the DGT NMEs for projectile and target nuclei. This study showed that the DGT NME is linearly correlated with the total NME for $0\nu\beta\beta$ decay predicted by the interacting boson model. Recently, Brase et al. [56] exploited this correlation to determine the NMEs for $0\nu\beta\beta$ decay in heavier candidate nuclei with an EFT.

Reference [45] shows, however, that the correlation does not appear in the results of the quasiparticle random-phase approximation (QRPA) [57], which has been extensively used to calculate $0\nu\beta\beta$ -decay NMEs [14,24,58–61]. The contradiction between this method and others needs to be resolved to clarify the significance of HIDCE experiments [53,54,62]. In this paper, we use *ab initio* methods to address the issue. So that we can compare our results with those obtained previously, we do not include the recently discovered contact transition operator [63,64], even though it might affect the $M^{0\nu\beta\beta}$ significantly within an *ab initio* framework [65].

The paper is organized as follows. In Sec. II, the formulas for the NMEs of both DGT transition and $0\nu\beta\beta$ decay are presented. The feature of neutrino potentials regularized with dipole form factors in coordinate space is discussed. In Sec. III, we briefly introduce the many-body methods and nuclear Hamiltonians that are employed in the *ab initio* calculations. The results are also discussed in comparison with a scale-separation analysis in Sec. IV. The in-medium renormalization effect is discussed in comparison with conventional shell-model results in Sec. V. Our conclusions are summarized in Sec. VI.

II. THE DGT AND $0\nu\beta\beta$ TRANSITIONS

The spin-parity J^{π} of the ground state of an even-even nucleus is 0⁺. A DGT transition connects this state to the final states of its neighboring even-even nucleus with spin-parity of 0⁺ or 2⁺. Here we only consider the ground-state-to-groundstate DGT transition as its NME is expected to be closest to that of the $0\nu\beta\beta$ decay, even though it might be just a small fraction (about 10⁻⁴) of the total DGT strength for the isospin-changing transitions [45,46,66]. The NME of the DGT transition is defined as

$$M^{\rm DGT} = \langle 0_f^+ | \sum_{1,2} [\sigma_1 \otimes \sigma_2]^0 \tau_1^+ \tau_2^+ | 0_i^+ \rangle, \tag{1}$$

where σ and τ are the spin and isospin operators, respectively. The nonzero matrix element of the isospin-raising operator is $\langle pp | \tau_1^+ \tau_2^+ | nn \rangle = 1$. The ground-state wave functions of initial and final nuclei $|0_{i/f}^+\rangle$ also enter into the expression for the NME $M^{0\nu\beta\beta}$ of $0\nu\beta\beta$ decay, which in the standard mechanism has the following form:

$$M^{0\nu\beta\beta} = \sum_{\alpha} \langle 0_f^+ | \sum_{1,2} h_{\alpha,K}(r_{12}) C_{\alpha}^K \cdot S_{\alpha}^K \tau_1^+ \tau_2^+ | 0_i^+ \rangle.$$
(2)

The symbol α runs over Fermi, GT, and tensor terms. The spin-spatial part $C_{\alpha}^{K} \cdot S_{\alpha}^{K}$ in (2) is the scalar product of two tensors with their expressions given by

$$C_F^0 = 1, \quad S_F^0 = 1,$$
 (3a)

$$C_{\rm GT}^0 = 1, \quad S_{\rm GT}^0 = \boldsymbol{\sigma}_1 \cdot \boldsymbol{\sigma}_2, \tag{3b}$$

$$C_T^2 = \sqrt{\frac{24\pi}{5}} Y_2(\hat{\boldsymbol{r}}_{12}), \quad S_T^2 = [\boldsymbol{\sigma}_1 \otimes \boldsymbol{\sigma}_2]^2.$$
(3c)

The coordinate-space neutrino potential is given by

$$h_{\alpha,K}(r_{12}) = \frac{2R_A}{\pi g_A^2} \int_0^\infty \mathrm{d}q \, q^2 \frac{h_{\alpha,K}(\boldsymbol{q}^2)}{q(q+E_d)} j_K(qr_{12}), \quad (4)$$

where $R_A = 1.2 A^{1/3}$ fm is introduced to make the matrix element $M^{0\nu\beta\beta}$ dimensionless. The relative coordinate between the two decaying neutrons is defined as $\mathbf{r}_{12} = \mathbf{r}_1 - \mathbf{r}_2$ and its magnitude $r_{12} = |\mathbf{r}_{12}|$ and direction vector $\hat{\mathbf{r}}_{12} =$ $\mathbf{r}_{12}/|\mathbf{r}_{12}|$. The average excitation energy E_d is chosen as $E_d =$ $1.12 A^{1/2}$ MeV [66]. We make this choice to facilitate comparison with prior work. In an EFT framework, E_d corresponds to a subleading correction [67]. The function $j_K(qr_{12})$ is the spherical Bessel function of rank K, where K = 0 for the Fermi and GT terms and K = 2 for the tensor term. The functions $h_{\alpha,K}(\mathbf{q}^2)$ are defined in terms of the vector (g_V) , axial-vector (g_A) , induced pseudoscalar (g_P) and weakmagnetism (g_M) coupling constants,

$$h_{F,0}(\boldsymbol{q}^2) = -g_V^2(\boldsymbol{q}^2),$$
 (5a)

$$h_{\text{GT},0}(\boldsymbol{q}^2) = g_A^2(\boldsymbol{q}^2) - \frac{2}{3} \frac{\boldsymbol{q}^2}{2m_p} g_A(\boldsymbol{q}^2) g_P(\boldsymbol{q}^2) + \frac{1}{3} \frac{\boldsymbol{q}^4}{4m_p^2} g_P^2(\boldsymbol{q}^2) + \frac{2}{3} \frac{\boldsymbol{q}^2}{4m_p^2} g_M^2(\boldsymbol{q}^2), \qquad (5b)$$

$$h_{T,2}(\boldsymbol{q}^2) = \frac{2}{3} \frac{\boldsymbol{q}^2}{2m_p} g_A(\boldsymbol{q}^2) g_P(\boldsymbol{q}^2) - \frac{1}{3} \frac{\boldsymbol{q}^4}{4m_p^2} g_P^2(\boldsymbol{q}^2) + \frac{1}{3} \frac{\boldsymbol{q}^2}{4m_p^2} g_M^2(\boldsymbol{q}^2),$$
(5c)

where the coupling constants are regularized by the following dipole form factors:

$$g_V(\boldsymbol{q}^2) = g_V(0) (1 + \boldsymbol{q}^2 / \Lambda_V^2)^{-2},$$
 (6a)

$$g_A(\boldsymbol{q}^2) = g_A(0) \left(1 + \boldsymbol{q}^2 / \Lambda_A^2 \right)^{-2},$$
 (6b)

$$g_M(q^2) = g_V(q^2)(1+\kappa_1),$$
 (6c)

$$g_P(q^2) = g_A(q^2) \left(\frac{2m_p}{q^2 + m_\pi^2}\right).$$
 (6d)

If not mentioned explicitly, then we choose the vector and axial-vector coupling constants $g_V(0) = 1$, $g_A(0) = 1.27$, the anomalous nucleon isovector magnetic moment $\kappa_1 = \mu_n^{(a)} - \mu_p^{(a)} = 3.7$, and the cutoff values $\Lambda_V = 0.85$ GeV and $\Lambda_A = 1.09$ GeV, following Refs. [68,69].

The dipole form factor regularizes the short-range behavior of the neutrino potentials. In the simplest case of $h_{F,0}(q^2) = -g_V^2(0)$, and $h_{\text{GT},0}(q^2) = g_A^2(0)$, i.e., without dipole form factors and induced higher-order currents, the Fermi- and GT-type neutrino potentials in coordinate space can be derived analytically [70]:

$$h_{F,0}(r_{12}) \to \tilde{h}_{F,0}(r_{12}) = -\left(\frac{g_V^2}{g_A^2}\right) R_A \frac{\phi(E_d r_{12})}{r_{12}},$$
 (7a)

$$h_{\text{GT},0}(r_{12}) \to \tilde{h}_{\text{GT},0}(r_{12}) = R_A \frac{\phi(E_d r_{12})}{r_{12}},$$
 (7b)

where the function $\phi(x)$ is defined as [66,70]

$$\phi(x) = \frac{2}{\pi} \sin(x) \operatorname{Ci}(x) + \cos(x) \left[1 - \frac{2}{\pi} \operatorname{Si}(x) \right], \quad (8)$$

with the sine and cosine integrals

$$\operatorname{Si}(x) = \int_0^x \frac{\sin t}{t} dt, \quad \operatorname{Ci}(x) = -\int_x^\infty \frac{\cos t}{t} dt. \quad (9)$$

To see how the short-range behavior of the neutrino potential $h_{\alpha,0}$ is regularized by the dipole form factors, we choose different values for the cutoffs Λ_V and Λ_A in $g_V(q^2)$ and $g_A(q^2)$, respectively. The corresponding Fermi and GT neutrino potentials are displayed in Figs. 1(a) and 1(b), respectively. It is shown that the use of a smaller cutoff value $\Lambda_{V/A}$ leads to a larger modification on the neutrino potentials in the short-distance region. As expected, the neutrino potentials approach a Coulomb-like potential in the limit $\Lambda_{V/A} \rightarrow \infty$ and $E_d = 0$, i.e., $\tilde{h}_{GT,0}(r_{12}) \rightarrow R_A/r_{12}$ as $\phi(0) = 1$. Besides, we show in Fig. 1(c) how a nonzero value of E_d reduces the entire neutrino potential. For the $0\nu\beta\beta$ -decay candidate nuclei with mass number $48 \leq A \leq 150$, the empirical value of E_d is $E_d \in [7.76, 13.71]$ MeV [66], and $r_{12} \simeq 1.0$ fm, which gives $\phi(E_d r_{12}) \in [0.86, 0.91]$, as shown in Fig. 2.

The NMEs of both DGT and $0\nu\beta\beta$ transitions can be conveniently rewritten as a function of the relative coordinate r_{12} between decaying nucleons [71],

$$M_A^{\kappa} = \int dr_{12} C_A^{\kappa}(r_{12}), \tag{10}$$

where κ stands for either $0\nu\beta\beta$ or DGT. It was pointed out in Ref. [56] that in conventional nuclear models, nuclear wave functions and neutrino potentials are represented in a harmonic oscillator basis with the oscillator length given by $b = \sqrt{\hbar/(M_N\omega)}$, where M_N is nucleon mass and the frequency ω scales as $A^{-1/3}$. Thus, the NME of $M_{GT}^{0\nu\beta\beta}$ (proportional to R_A/b) is expected to scale as $A^{1/6}$ and the DGT matrix element M_{GT}^{DGT} is expected to be correlated with $M_{GT}^{0\nu\beta\beta}A^{-1/6}$ for all isotopes. From another point of view, if the $0\nu\beta\beta$



FIG. 1. The (a) Fermi and [(b) and (c)] GT-type neutrino potentials (4) as a function of r_{12} . The radius parameter R_A is excluded in the neutrino potentials to facilitate comparison with the Coulomblike potential $1/r_{12}$. Different choices of cutoff values Λ_V and Λ_A are employed in the dipole form factors. In (c), a different value of the average excitation energy E_d is used. All the Fermi-type neutrino potentials are multiplied with a minus sign to facilitate comparison. In panel (b), the potential $h_{GT,0}^{AA}(q^2) = g_A^2(q^2)$ is compared to that of the full GT operator (5b) with the cutoff value change to $\Lambda_A = 1.09$ GeV.

decay is dominated by the short-range contribution, namely the long-range Coulomb-like decaying behavior is regularized by a faster decaying two-nucleon wave function, then one may expect that the DGT matrix element M^{DGT} is correlated with $M_{\text{GT}}^{0\nu\beta\beta}A^{-1/3}$, where the factor $A^{-1/3}$ is from the radius R_A introduced to make the NME dimensionless, cf. (4). These two correlation relations will be discussed using the results from the calculations of both conventional nuclear models and *ab initio* methods in the next section.



FIG. 2. The function $\phi(x)$ defined in Eq. (8) with the limits of $\phi(0) = 1$ and $\phi(\infty) = 0$. The shaded area indicates the interval where the average excitation energy takes $E_d \in [7.76, 13.71]$ MeV and the relative distance between the two decaying nucleons $r_{12} =$ 1.0 fm.

III. AB INITIO MANY-BODY CALCULATIONS

In this section, we carry out ab initio nuclear manybody calculations with the importance-truncated no-core shell model (IT-NCSM) [72], valence-space in-medium similarity renormalization group (VS-IMSRG) [73], and in-medium generator coordinate method (IM-GCM) [42,74] starting from chiral NN+3N interactions. The latter two are different variants of IMSRG [75], which introduces a flow equation to gradually decouple the off-diagonal elements of the Hamiltonian that are connecting the valence space and the excluded spaces or to decouple a preselected reference state from all other states. In the VS-IMSRG, an effective Hamiltonian in a specific valence space is obtained, while in the IM-GCM [42,74], the reference state becomes a reasonable approximation to the ground state of the evolved Hamiltonian. The close-to-exact ground state is obtained with GCM by admixing other states that differ only in their collective parameters. A unitary transformation is defined via the flow equation and this transformation is consistently applied to all operators of interest. We employ the chiral nuclear interaction (up to $N^{3}LO$) by Entem and Machleidt [76], which we indicate by the label "EM." We use the free-space SRG [77] to evolve the EM interaction to a resolution scale of $\lambda = 1.8 \text{ fm}^{-1}$. Following Refs. [78,79], we construct the 3N interaction directly, with a chiral cutoff of $\Lambda = 2.0 \text{ fm}^{-1}$. We refer to the resulting NN+3N Hamiltonian as $EM\lambda/\Lambda$, i.e., EM1.8/2.0—see Refs. [78,79] for details. For comparison, we also employ the recently proposed chiral force $\Delta N^2 LO_{GO}(394)$ [80], a low-cutoff NN+3N interaction whose construction accounts for Δ isobars and whose parameters are constrained by $A \leq 4$ few-body data as well as nuclear matter properties. For the 3N interaction, we discard all matrix elements involving states with $e_1 + e_2 + e_3 > 14$, where $e_i = 2n_i + \ell_i$ is the number of oscillator quanta in state *i*. The frequency of the harmonic oscillator basis is chosen as $\hbar \omega = 16$ MeV.

Figures 3 and 4 display the results of three *ab initio* calculations for both isospin-conserving and isospin-changing DGT transitions and GT- $0\nu\beta\beta$ decay in a set of *p*-, *sd*-, and *fp*shell nuclei, respectively, where the same value of $e_{\text{Max}} = 8$



FIG. 3. The transition densities $\tilde{C}^{\alpha}(r_{12})$ for the isospinconserving transitions of DGT [(a)–(c)] and GT-0 $\nu\beta\beta$ decay [(d)–(f)] from three *ab initio* calculations for ⁶He, ¹⁰Be, and ¹⁴C using the EM1.8/2.0 chiral force. The tilde means that the transition densities are normalized to unity at the first peak position.

is used in the three methods for comparison. As discussed in the previous paper [41], $N_{\text{Max}} = 8$ is usually employed in the IT-NCSM calculations, except for the transitions of ⁸He and ²²O where the $N_{\text{Max}} = 6$ and 4 is used, respectively. In the two variants of IMSRG calculations, the transition density $C^{\kappa}(r_{12})$ is evaluated using the corresponding IMSRG-evolved transition operator at each value of the relative coordinate r_{12} and the nuclear wave functions by the evolved Hamiltonian. One can see that the short-range parts of both transition densities by all the three *ab initio* methods are on top of each other. The predictions for the long-range part of the isospin-changing DGT transitions differ because the long-range part of the transition densities is more sensitive to the way each method models many-body correlations. Due to the presence of the neutrino potential which decreases with r_{12} approximately as $1/r_{12}$ (cf. Fig. 1), the discrepancy in the long-range part of $\tilde{C}^{\text{GT}}(r_{12})$ among the three methods is strongly suppressed. As a result, the short- and long-range parts of $0\nu\beta\beta$ decay of both isospin-conserving and isospin-changing types are consistently described in the three calculations. It is also seen that the transition densities of isospin-conserving transitions do not change sign as a function of r_{12} . In contrast, those of isospin-changing transitions oscillate with r_{12} , and thus the contributions of long- and short-range regions compensate for



FIG. 4. Same as Fig. 3 but for isospin-changing transitions of 8 He, 22 O, and 48 Ca.

each other. Figures 4(b) and 4(c) show that the long-range contribution in the VS-IMSRG is generally larger than that in the IM-GCM. In the VS-IMSRG calculation, the long-range contribution to the DGT transition can be even larger than the short-range part, resulting in an inverted sign for the DGT NME. We note that varying the e_{Max} around the selected value does not change the shape of the transition densities but modifies slightly the height of the peaks.

As discussed in Ref. [45] and in the next section, if the processes of both DGT transition and $0\nu\beta\beta$ decay are dominated by the short-range contribution, then these two types of matrix elements are expected to be correlated, irrespective of if the isospin is changing or conserving in the process. To verify this finding, we show all NMEs of DGT transitions and $0\nu\beta\beta$ decay from the three *ab initio* calculations for the isotopes in different mass regions in Fig. 5. In these calculations, the value of e_{Max} is chosen to ensure the convergence of the NMEs, as discussed in Ref. [41]. In the VS-IMSRG calculation for heavier isotopes, $e_{\text{Max}} = 12$ is used. The NMEs M^{DGT} are plotted against those of $0\nu\beta\beta$ decay scaled as $M^{0\nu\beta\beta}A^{-1/3}$ and $M^{0\nu\beta\beta}A^{-1/6}$ in Figs. 5(a) and 5(b) and Figs. 5(c) and 5(d), respectively. The results are fitted to the following relation:

$$M^{\rm DGT} = \alpha M^{0\nu\beta\beta} A^{\gamma} + \beta, \qquad (11)$$

where the power parameter is fixed to be either $\gamma = -1/3$ or -1/6. The coefficient of correlation *r* between the two quantities,

$$r = \frac{\sum_{i} (x_{i} - \bar{x})(y_{i} - \bar{y})}{\sqrt{\sum_{i} (x_{i} - \bar{x})^{2} \sum_{i} (y_{i} - \bar{y})^{2}}},$$
(12)

is introduced as a statistical measure of the strength of the correlation relationship, where x_i and y_i are the samples on the *x* and *y* axis, respectively, and \bar{x} and \bar{y} are the average of the samples on each axis. The value r = 1 corresponds to a perfect positive linear correlation and r = 0 indicates no linear relationship. The correlation relationships that we obtain are as follows:

$$M^{\text{DGT}} = 1.21(3)M^{0\nu\beta\beta}_{\text{Tot}}A^{-1/3} - 0.32(2), \quad r = 0.99,$$
(13a)

$$M^{\rm DGT} = 1.57(4) M_{\rm GT}^{0\nu\beta\beta} A^{-1/3} - 0.37(3), \quad r = 0.99,$$
(13b)

$$M^{\text{DGT}} = 0.85(3)M^{0\nu\beta\beta}_{\text{Tot}}A^{-1/6} - 0.42(4), \quad r = 0.98,$$
(13c)

$$M^{\text{DGT}} = 1.12(3)M^{0\nu\beta\beta}_{\text{GT}}A^{-1/6} - 0.50(4), \quad r = 0.98.$$
 (13d)

The finite-sample-size error on the coefficient r is evaluated using the formula [81] $\sigma_r \approx (1 - r^2)/\sqrt{N - 2}$, which gives $\sigma_r = 0.003$ for r = 0.99 and $\sigma_r = 0.006$ for r = 0.98. One can see that the M^{DGT} is correlated slightly stronger with the quantity $M^{0\nu\beta\beta}A^{-1/3}$ than with the quantity $M^{0\nu\beta\beta}A^{-1/6}$, irrespective of if only the GT part or the total NME is considered for $M^{0\nu\beta\beta}$. The Fermi part of $M^{0\nu\beta\beta}$ is approximately one-third of the GT part, which explains why we still see a correlation when considering the total NME $M^{0\nu\beta\beta}$. Besides, the residuals exhibit some clear pattern at low values of $M^{0\nu\beta\beta}$ in all cases, more predominantly when scaling $M^{0\nu\beta\beta}$ by $A^{-1/6}$. This finding not only further validates that using $A^{-1/3}$ is a better choice for the results of *ab initio* calculations, but it also suggests that the good correlation we find when looking at the whole data set is not representative of the whole situation. We note that the results from ab initio calculations using the $\Delta N^2 LO_{GO}(394)$ interaction are consistent with the above correlation relation (13a).

It is worth emphasizing that the matrix elements for isospin-changing transitions and isospin-conserving transitions are significantly different from each other. The relevant $0\nu\beta\beta$ decay is an isospin-changing transition, and thus its matrix element is generally smaller than those of isospinconserving transitions for the same mass number A. Therefore, one may expect the correlation obtained from these two types of transitions to be different. With this consideration, we plot the NMEs of isospin-changing and isospin-conserving transitions separately in Figs. 6(a) and 6(b), respectively. Two different linear regressions are carried out for these two types of transitions with the parameter $\gamma = -1/3$. We find the following correlation relationships:

$$\Delta T = 2: M^{\text{DGT}} = 1.21(20) M_{\text{GT}}^{0\nu\beta\beta} A^{-1/3} - 0.30(5), \quad (14a)$$

$$\Delta T = 0: M^{\text{DGT}} = 1.43(14)M^{0\nu\beta\beta}_{\text{GT}}A^{-1/3} - 0.17(20).$$
(14b)

It is shown that the correlation coefficient (r = 0.72, $\sigma_r = 0.08$) of isospin-changing transitions ($\Delta T = 2$) is much smaller than that (r = 0.95, $\sigma_r = 0.03$) of isospin-conserving ($\Delta T = 0$) transitions. In other words, the correlation relation is much weaker for the NMEs of isospin-changing transitions.



FIG. 5. Correlation between the NMEs of DGT transitions (M^{DGT}) with those of $GT-0\nu\beta\beta$ decay $(M_{GT}^{0\nu\beta\beta})$ or the full NMEs of $0\nu\beta\beta$ decay $(M_{Tot}^{0\nu\beta\beta})$, scaled by a mass-number-dependent factor of either $A^{-1/6}$ or $A^{-1/3}$, respectively. The NMEs are obtained from *ab initio* calculations with the chiral NN+3N interactions EM1.8/2.0 or $\Delta N^2 LO_{go}$ for isotopes ranging from A = 6 to A = 76. Only the results obtained with EM1.8/2.0 are used in the linear regression with residuals given in the bottom row. The shaded area indicates the confidence interval with a 95% confidence level while the dashed line indicates the prediction interval at a 95% confidence level. See text for details.



FIG. 6. Correlation between M^{DGT} and $M^{0\nu\beta\beta}_{\text{GT}}A^{-1/3}$ when considering only isospin-changing transitions (left column) or only isospin-conserving transitions (right column). The green bands show the 95% confidence interval while the dashed lines show the 95% prediction interval. The best fit lines from both cases agree with each other within 1σ .

Figures 7 and 8 summarize the results from the calculations of both *ab initio* and conventional nuclear models (shell models [45,50,51], energy-density-functional (EDF) [82] and QRPA [57]). Treating the matrix elements of isospin-conserving and isospin-changing processes, separately, we



FIG. 7. Correlation between M^{DGT} and $M^{0\nu\beta\beta}_{\text{Tot}}A^{-1/3}$ derived from the results of both *ab initio* (red) and conventional (blue) nuclear models [45,50,51,57,82], except for the results of QRPA calculations [57] for isospin-conserving transitions in (a) and isospin-changing in (b).



FIG. 8. Same as Fig. 7, but for the correlation between M^{DGT} and $M^{0\nu\beta\beta}_{\text{Tot}}A^{-1/6}$.

find the following relation for the nuclear matrix elements from the *ab initio* calculations,

$$\Delta T = 2: M^{\text{DGT}} = 1.08(15)M^{0\nu\beta\beta}_{\text{Tot}}A^{-1/3} - 0.26(4), \quad (15a)$$

$$\Delta T = 0: M^{\text{DGT}} = 1.20(9) M_{\text{Tot}}^{0\nu\beta\beta} A^{-1/3} - 0.30(15), \quad (15b)$$

$$\Delta T = 2: M^{\text{DGT}} = 0.51(7) M^{0\nu\beta\beta}_{\text{Tot}} A^{-1/6} - 0.23(4), \quad (15c)$$

$$\Delta T = 0: M^{\text{DGT}} = 1.09(10) M^{0\nu\beta\beta}_{\text{Tot}} A^{-1/6} - 1.06(26), \ (15d)$$

and the following relation for the nuclear matrix elements from the conventional nuclear-model calculations:

$$\Delta T = 2: M^{\text{DGT}} = 1.32(3)M^{0\nu\beta\beta}_{\text{Tot}}A^{-1/3} - 0.19(2), \quad (16a)$$

$$\Delta T = 0: M^{\text{DGT}} = 1.87(20) M^{0\nu\beta\beta}_{\text{Tot}} A^{-1/3} - 0.68(22), \quad (16b)$$

$$\Delta T = 2: M^{\text{DGT}} = 0.58(2)M^{0\nu\beta\beta}_{\text{Tot}}A^{-1/6} - 0.14(2), \quad (16c)$$

$$\Delta T = 0: M^{\text{DGT}} = 1.10(9)M^{0\nu\beta\beta}_{\text{Tot}}A^{-1/6} - 0.93(18).$$
(16d)

It is seen that the correlation relations are significantly different for the matrix elements of $\Delta T = 2$ and $\Delta T = 0$ transitions. The slope for the $\Delta T = 0$ transitions is generally larger than that for the $\Delta T = 2$ transitions. Besides, the slope by the *ab initio* calculation is generally smaller than that by the conventional nuclear model calculations because of a stronger cancellation between the long-range and short-range contributions in the DGT NMEs by *ab initio* methods. Among all the methods concerned, the QRPA predicts the smallest value for the DGT NMEs, indicating the occurrence of the strongest cancellations [57].

To compare with the correlation relation in Refs. [45,56], we follow their way to derive the correlation relation based on the matrix elements of both isospin-conserving and isospin-changing processes from the calculations of conventional nuclear models (excluding the results of QRPA), which reads

$$M^{\text{DGT}} = 1.41(3)M^{0\nu\beta\beta}_{\text{Tot}} \cdot A^{-1/3} - 0.23(2), \quad r = 0.96, \quad (17a)$$

$$M^{\text{DGT}} = 0.65(2)M^{0\nu\beta\beta}_{\text{Tot}} \cdot A^{-1/6} - 0.20(2), \quad r = 0.94,$$
(17b)

the latter of which is consistent with the interval of parameters $\alpha \in [0.447, 0.699]$, $\beta \in [-0.18, -0.056]$, with the power parameter $\gamma = -1/6$ found in Ref. [56]. The value of the coefficient *r* indicates that the DGT matrix elements M^{DGT} are slightly stronger correlated with $M_{\text{Tot}}^{0\nu\beta\beta}A^{-1/3}$ than with $M_{\text{Tot}}^{0\nu\beta\beta}A^{-1/6}$, as discussed in Fig. 5.

It is worth mentioning that the NMEs of the groundstate-to-ground-state DGT transition of isospin-changing transitions would be exactly zero if the spin-isospin SU(4) symmetry were conserved in atomic nuclei as in this case the initial and final states would belong to different irreducible representations of the SU(4) group. Therefore, different values of the DGT NMEs predicted by different nuclear models indicate that the SU(4) symmetry is broken to different extents in the ground-state wave functions. The studies on the breaking of the SU(4) symmetry in different nuclear models and its impact on the correlation relation might be able to provide us with a more profound understanding, but these studies are beyond the scope of this work.

IV. A SCALE-SEPARATION ANALYSIS

In this section, we carry out a scale-separation analysis of the correlation between the NMEs of DGT transitions and $0\nu\beta\beta$ decay based on the argument [83–85] that the longand short-distance physics in atomic nuclei can be rather well separated. In this case, the nuclear many-body wave functions $\Psi_{i/f}(1, 2, ..., A)$ of initial and final nuclei in the transitions can be approximately factorized into the product of a universal short-distance two-body wave function $\phi_c(\mathbf{r}_{12})$ and a statedependent long-range *A*-body wave function $\chi_c^{(i/f)}$ [86]

$$\Psi_{i/f} \xrightarrow{r_{12} \to 0} \sum_{c} \phi_c(\mathbf{r}_{12}) \chi_c^{(i/f)}(\mathbf{R}_{12}; 3, \dots, A), \quad (18)$$

where $\mathbf{R}_{12} = (\mathbf{r}_1 + \mathbf{r}_2)/2$, and *c* distinguishes channels (spinisospin) with different quantum numbers for the pair of nucleons. For the NMEs of DGT and $0\nu\beta\beta$ transitions, the pair of neutrons in the *s*-wave state with total spin S = 0, angular momentum J = 0, and isospin T = 1 converting into a pair of protons with the same quantum numbers provides the predominate contribution. If only this channel (labeled with c = 0) is considered and the two-body wave function $\phi_c(r_{12})$ is assumed to be isospin independent, then one has $\phi_{nn}(r_{12}) = \phi_{pp}(r_{12}) = \phi_0(r_{12})$. Thus, it is reasonable to parametrize the transition densities defined in (10) into the following forms:

$$\tilde{C}_{A}^{\text{GT}}(r_{12}) = -3r_{12}^{2}h_{\text{GT},0}(r_{12})\rho_{nn}(r_{12})C_{ppnn}^{0}(A_{f}, A_{i}, r_{12}),$$
(19a)

$$\tilde{C}_{A}^{\text{DGT}}(r_{12}) = \sqrt{3}r_{12}^2\rho_{nn}(r_{12})C_{ppnn}^0(A_f, A_i, r_{12}),$$
(19b)

where the factor of (-3) in the GT transition is from the spin operator. The DGT defined in (1) brings an additional factor of $(-1/\sqrt{3})$. The two-nucleon density is defined as $\rho_{nn}(r_{12}) = |\phi_0(r_{12})|^2$. The overlap function $C_{ppnn}^0(A_f, A_i, r_{12})$ is determined by the *A*-body wave functions $\chi_0^{(i/f)}$ of initial and final nuclei multiplied by the number of pairs [86]. In the limit of $r_{12} \rightarrow 0$, the two-nucleon density $\rho_{nn}(r_{12})$ is expected to be universal for different nuclei but may depend

on the employed nuclear force. The overlap function $C_{ppnn}^{0}(A_{f}, A_{i}, r_{12})$ in the short-distance region $r_{12} < 1.0$ fm can be approximated by a constant $\mathcal{N}_{ppnn}^{A}(f, i)$ depending only on the nucleus and on the details of employed nuclear force [85,87,88]. The ratios $\mathcal{N}_{ppnn}^{A}(f, i)/\mathcal{N}_{ppnn}^{A'}(f, i)$ for any two nuclei with mass numbers A and A' are less model dependent. Recently, this property has been exploited in the calculation of the NMEs of $0\nu\beta\beta$ candidate nuclei by combining quantum Monte Carlo and the nuclear shell model [86].

If the processes of both DGT transition and $0\nu\beta\beta$ decay are dominated by the short-range contribution, as pointed out in Ref. [45], then one would have the ratio of these two matrix elements from Eq. (19),

$$\frac{M_{\rm GT}^{0\nu\beta\beta}}{M^{\rm DGT}} \approx \lim_{\epsilon \to 0} \frac{-3\int_0^{\epsilon} dr_{12}r_{12}^2\rho_{nn}(r_{12})h_{\rm GT,0}(r_{12})}{\sqrt{3}\int_0^{\epsilon} dr_{12}r_{12}^2\rho_{nn}(r_{12})} \\
= \lim_{\epsilon \to 0} -\sqrt{3}R_A \frac{\int_0^{\epsilon} dr_{12}r_{12}\rho_{nn}(r_{12})}{\int_0^{\epsilon} dr_{12}r_{12}^2\rho_{nn}(r_{12})} = \lim_{\epsilon \to 0} C(\epsilon)A^{1/3},$$
(20)

where $C(\epsilon)$ is a constant depending on the short-range parameter ϵ . This expression holds for both isospin-changing and isospin-conserving transitions, because the isospin effects are encoded in the overlap functions, which cancel in the ratio. Of course, in the realistic case, the validity of the approximations employed to derive the above relation varies with atomic nuclei, and it will be studied in combination with the results of *ab initio* calculations.

Let us first parametrize the two-nucleon density $\rho_{nn}(r)$ with the following function form:

$$\rho_{nn}(r) = F_{\text{SRC}}(r)e^{-r^2/d}, \qquad (21)$$

where the parameter *d* controls the long-range decay behavior. The function $F_{\text{SRC}}(r)$ is introduced to mimic the effect of short-range correlation (SRC) [89],

$$F_{\rm SRC}(r) = 1 - ce^{-ar^2} \left(\gamma + \sum_{i=1}^3 b_i r^{i+1}\right).$$
(22)

The parameters $a = 3.17 \text{ fm}^{-2}$, $\gamma = 0.995$, c = 1, $b_1 = 1.81 \text{ fm}^{-2}$, $b_2 = 5.90 \text{ fm}^{-3}$, and $b_3 = -9.87 \text{ fm}^{-4}$ were obtained in Ref. [89] by fitting to the results from cluster variational Monte Carlo calculations for the proton-proton/neutron-neutron correlation functions. Here we vary the values of the parameters *c* and *d* within an interval producing reasonable transition densities to simulate the impacts from the nuclear-force dependence and nucleus dependence [85] on the correlation relation between the NMEs of DGT and GT- $0\nu\beta\beta$ transitions. Specifically, we vary the parameter *c* between 0 and 1 and the *d* between 0.5 and 5 to examine the sensitivity of the correlation relation to these two parameters.

As shown in Figs. 3 and 4, the transition densities $\tilde{C}^{\kappa}(r_{12})$ of isospin-conserving and isospin-changing transitions have different dependence on r_{12} , so we treat these two transitions separately. Besides, it has been found in previous many-body calculations [11,42,71] that the transition densities $C^{\kappa}(r_{12})$ may possess node structures varying in detail with nuclear models and the mass regions of isotopes. We note that this structure is generally similar for the same type of transitions

in the isotopes of the same mass region for a given nuclear model. To reproduce this node structure, we approximate the overlap function $C_{ppnn}^{0}(A_f, A_i, r_{12})$ with the following simple form:

$$C_{ppnn}^{0}(A_f, A_i, r_{12}) \simeq p(r_{12})q(r_{12})\mathcal{N}_{ppnn}^A(f, i),$$
 (23)

where p(r) and q(r) are polynomial functions $c_0[1 +$ $\sum_{i=1}^{N} c_{2i-1} (r - c_{2i})^{2i}$ with different values of c_i for isospin-changing and isospin-conserving transitions. The parameter N is determined by the node structure of the transition densities and the parameters c_i need to be fitted to transition densities of each nuclear model. In our case, as shown later, we fit to the results of a few transitions in each valence space and take the average values given for the c_i 's from those fits. We also imposed p(r) = q(r) for the case of isospin-conserving transitions to ensure isospin symmetry. Figure 9 displays the transition densities of isotopes in p, sd, and fp shells from the VS-IMSRG calculation with the EM1.8/2.0 interaction. One can see that the distribution of the transition density could have a complex structure with more than two peaks. In particular, long-range contribution to the DGT transition could be even more significant than the short-range contribution, and it could enhance (quench) the isospin-conserving (isospin-changing) transitions.

To examine the correlation between the NMEs of DGT and $0\nu\beta\beta$ decay in a more general way, we use the transition densities in Fig. 9 to optimize the parameters *c* and *d* in the two-nucleon density and c_i s in the function $p(r_{12})$ and $q(r_{12})$. Once these parameters are determined, we vary the parameters *c* and *d* around their optimal values which allows to generate more transition densities to look at the correlation relation. The sampled transition densities $\tilde{C}_A^{\kappa}(r_{12})$ are displayed in Fig. 10. One can see that the function form (19) for the transition density, together with the twonucleon density $\rho_{nn}(r)$ in (21) and the polynomial function for the overlap $C_{ppnn}^0(A_f, A_i, r_{12})$ in (23) can reproduce nicely the main structure of both isospin-conserving and isospinchanging transition densities in each mass regions from the VS-IMSRG calculations.

The sampled transitions are then integrated over the coordinate r_{12} which leads to the NME \tilde{M}_{DGT} and $\tilde{M}_{\text{GT}}^{0\nu\beta\beta}$,

$$\tilde{M}^{\kappa} = \int dr_{12} \tilde{C}^{\kappa}_A(r_{12}), \qquad (24)$$

where $\tilde{C}_{A}^{\kappa}(r_{12})$ has been defined in (19) with the overlap function (23). The NME \tilde{M}^{κ} might differ from the actual value of the NME M^{κ} because of the unknown constant $N_{ppnn}^{A}(f, i)$, but it does not impact the analysis of the correlation relation between them, except for the intercept parameter.

between them, except for the intercept parameter. The NMEs \tilde{M}^{DGT} and $\tilde{M}_{\text{GT}}^{0\nu\beta\beta}$ for isospin-conserving transitions are displayed in Fig. 11. It is shown that these two types of NMEs are correlated in some way. With the transition densities generated by varying the parameters c and d around those values from the VS-IMSRG calculations for the p- and sd-shell isotopes, the resultant M^{DGT} is increasing with $M_{\text{GT}}^{0\nu\beta\beta}$ in a parabolic form approximately, while that derived from the f p-shell isotopes is in a linear form. These correlation relations are summarized in Fig. 12, where the NMEs from





FIG. 9. The transition densities [cf. Eq. (10)] of isotopes in *p*, *sd*, and *fp* shells from the VS-IMSRG calculation with the EM1.8/2.0 interaction. Asterisks indicate isospin conserving transitions.

FIG. 10. Sampled transition densities $\tilde{C}_A^{\kappa}(r_{12})$ with the parameters fitted to the results of VS-IMSRG calculations for the isospin-conserving transitions in *p*-, *sd*-, and *f p*-shell nuclei, except for the parameters *c* and *d*, which vary from 0 to 1 and from 0.5 to 4.5 fm², respectively.



FIG. 11. The NMEs from the integral of the sampled transition densities in Fig. 10. Arrows show the direction of increasing the d value.



FIG. 12. The correlation relations are derived from the sampled transition densities for isospin-conserving transitions. The area is originated from the variants of the parameters c and d which control the short-range and long-range behavior, respectively. The NMEs from the VS-IMSRG calculations are added for comparison.



FIG. 13. Sampled transition densities with parameters fitted to the VS-IMSRG results for isospin-changing transitions in the (a) *sd*-and (b) f p-shell nuclei, with the parameters c and d increasing from 0 to 1 and from 0.5 to 4.5 fm², respectively.

the VS-IMSRG calculations are added for comparison. It is shown that the location of the VS-IMSRG calculations is generally within the area of the correlation relation for the isotopes in each mass region. However, the correlation relations for the isotopes in different mass regions are offset from each other. It implies that there are probably different correlation relations for the isospin-conserving transitions of isotopes in different mass regions.

The correlation relationship between the NMEs of isospinchanging transitions is more complicated. The corresponding sampled transition densities are displayed in Fig. 13 and the obtained NMEs are shown in Fig. 14. Again, one can see that the main structure of the transition densities exhibited in those by the VS-IMSRG calculation is reproduced in the sampled ones. Due to the strong cancellation between long-range and short-range contributions, the final DGT NME is significantly quenched. As a result, the value of the DGT NME varies from a small negative value to a small positive value with the



FIG. 14. The NMEs from the integral of the sampled transition densities in Fig. 13. (a) For *sd*-shell nuclei and (b) for fp-shell nuclei. Arrows show the direction of increasing the *d* value.

parameters c and d. There is a kind of weak correlation between the NMEs of DGT transitions and $0\nu\beta\beta$ decay, depending much on the mass region of the isotopes and the values of c and d.

For comparison, we perform a similar analysis based on the transition densities from conventional shell-model calculations for isotopes in sd and fp shells. The results are shown in Fig. 15. In contrast to the results from the VS-IMSRG calculations, the linear correlation relation between the NMEs is very robust. Varying the parameters c and d seems only change the intercept parameter of the linear correlation relation. The main difference between the sampled densities in Fig. 13 and Fig. 15 is the contributions from the intermediate- and longrange regions to the NME. A strong cancellation is shown in transition densities derived from the VS-IMSRG calculation but not in those from the conventional shell-model calculations. The previous analysis starts from (but is not limited to) the assumption that the transition matrix elements are dominated by the short-ranged part of the operator. As can be seen in Fig. 4, the DGT operator satisfies this requirement rather poorly. It is worth considering whether limiting this operator to shorter distances would enhance the scale separation and thus improve the correlation with the $0\nu\beta\beta$ amplitude. Such a restriction to short distances could conceivably be motivated by the light-ion induced double charge exchange reaction mechanism being surface peaked and therefore requiring both exchanged nucleons to be relatively localized. Our aim here is not to model the reaction process realistically but to test whether requiring the DGT operator to be short range improves the correlation with the NME of $0\nu\beta\beta$ decay. To this



FIG. 15. (a) The transition densities from conventional shellmodel calculations; (b) sampled transition densities with parameters fitted to the isospin-changing transition densities in (a). (c) The NMEs from the integral of the sampled transition densities in (b).

end, we define the NME of the surface localized DGT as

$$M_{\rm surf}^{\rm DGT} = \langle 0_f^+ | \sum_{1,2} f(r_{12}) g(r_{\rm CM}) [\boldsymbol{\sigma}_1 \otimes \boldsymbol{\sigma}_2]^0 \tau_1^+ \tau_2^+ | 0_i^+ \rangle, \quad (25)$$



FIG. 16. Correlation between the NMEs M^{DGT} of DGT transitions and the $M_{\text{GT}}^{0\nu\beta\beta}$ scaled by $A^{-1/3}$ using a subset of isospin-changing transitions for the *sd*- and *fp*-shell nuclei from the conventional shell-model calculation as well as the results obtained with the VS-IMSRG wave functions and the transition operators in the harmonic oscillator (HO) basis, Hartree-Fock basis (HF), or fully (full) evolved basis. See text for details.

where

$$f(r_{12}) = \exp\left(-r_{12}^2/2\right),$$
 (26)

$$g(R_{12}) = 1 - \exp\left(-\frac{R_{12}^2}{2R_A^2}\right),$$
 (27)

with $R_{12} = |(\mathbf{r}_1 + \mathbf{r}_2)/2|$ representing the position of the center of mass of the two particles. The functions $f(r_{12})$ and $g(R_{12})$ ensure that the two particles are close to each other and on the surface of the nucleus, respectively. We find that the correlation of the NME by this operator with $M^{0\nu\beta\beta}A^{-1/3}$ is worse than the standard DGT operator in (1).

V. THE IN-MEDIUM RENORMALIZATION EFFECT

In this section, we try to understand the origin of the discrepancy between the results of the conventional nuclear shell model and VS-IMSRG as these two methods are comparable. To further understand why a correlation is found in the nuclear shell-model calculation for isospin-changing transitions, we select a subset of transitions that shows a very strong correlation. This subset consists of the transitions ${}^{22}\text{O} \rightarrow {}^{22}\text{Ne}$, ${}^{28}\text{Mg} \rightarrow {}^{28}\text{Si}$, ${}^{44}\text{Ca} \rightarrow {}^{44}\text{Ti}$, ${}^{48}\text{Ca} \rightarrow {}^{48}\text{Ti}$, ${}^{56}\text{Ca} \rightarrow {}^{56}\text{Ti}$, and ${}^{52}\text{Ti} \rightarrow {}^{52}\text{Cr}$. In the conventional shell-model calculations, the USDB interaction [90] is used for the *sd*-shell nuclei and the GXPF1A interaction [91] for the *f p*-shell nuclei. The results are shown in Fig. 16. Based on the results within this subset, one finds the following correlation relationship:

$$M^{\rm DGT} = 1.86(20) M_{\rm GT}^{0\nu\beta\beta} A^{-1/3} - 0.37(8), \quad r = 0.98.$$
 (28)

The correlation coefficient r = 0.98 indicates that the two NMEs are strongly correlated, as shown in Fig. 16(a). This result is consistent with the previous shell-model study [45], as expected. In contrast, Fig. 16(d) shows that the correlation relation from the full VS-IMSRG calculation is weakened with the correlation coefficient r = 0.91. We note that the $0\nu\beta\beta$ NMEs of these nuclei by the the VS-IMSRG are generally smaller than those by the conventional shell-model calculations, while the DGT NMEs are much smaller and even

with opposite sign. This is mainly due to the configuration mixing in nuclear wave functions predicted differently in the calculations using the conventional shell-model interactions and those derived from VS-IMSRG.

To better understand how the in-medium renormalization effect from the VS-IMSRG evolution on the transition operator affects the correlation, we provide two intermediate results in Figs. 16(b) and 16(c), where the nuclear wave functions are from the VS-IMSRG calculation, while the transition operator in the harmonic oscillator (HO) basis or in the Hartree-Fock (HF) basis is used, respectively. In these two intermediate results, the transition operator is not consistently evolved. One can see that the correlation in the results of calculations with either the HO transition operator or the HF operator is even stronger than that of the conventional shell model in which the operator is represented in the harmonic oscillator basis.

Taking ⁴⁸Ca as an example, we illustrate how the transition density distribution looks in the four types of calculations shown in Fig. 16. The transition densities are displayed in Fig. 17. One can see that the use of the transition operator from the HO one to the IMSRG evolved one modifies the transition densities slightly in both short-range ($\simeq 0.6$ fm) and long-range ($\simeq 2.0-3.0$ fm) regions. Quantitatively, this modification is slightly different for different isotopes. Figure 18 shows the transition densities for the isospin-conserving transition ${}^{6}\text{He} \rightarrow {}^{6}\text{Be}$ and for the isospin-changing transition ⁸He \rightarrow ⁸Be obtained with and without VS-IMSRG evolution. The evolution enhances the short-range part of the transition densities for the isospin-conserving transitions, leading to an overall enhancement of the NME. This behavior is found in all the isospin-conserving cases. For the isospin-changing transition of ⁸He, the evolution increases the magnitude of the peaks of the distributions but, due to cancellations, does not necessarily increase the final NMEs. This behavior is found in all isospin-changing cases. In particular, one finds that the renormalization effect on the short-range part of the transition density is more significant in the isospin-conserving transitions than in the isospin-changing transitions. In short, the effect of VS-IMSRG evolution on isospin-changing



FIG. 17. The transition densities of both DGT and the GT part of the $0\nu\beta\beta$ decay from ⁴⁸Ca to ⁴⁸Ti from the conventional nuclear shell-model calculation with the GXPF1A interaction and the VS-IMSRG calculation using the transition operators in three types of basis: HO basis, HF basis, and the fully evolved basis.



FIG. 18. Evolution of the transition distribution of the GT part of the $0\nu\beta\beta$ decay and DGT transitions from the VS-ISMSRG in (a) ⁶He and (b) ⁸He.



FIG. 19. The transition densities of the DGT and GT part of the $0\nu\beta\beta$ decay in ⁴⁸Ca from the IM-GCM calculation using either the bare or evolved transition operator, where the same nuclear wave functions but different transition operators are used, respectively.

transitions is generally small, but the details of the cancellation between short- and long-range components depend on the nucleus and the operator, which finally degrades the correlation.

To help assess the method dependence of these conclusions, Fig. 19 displays the transition densities of the DGT and GT part of the $0\nu\beta\beta$ decay in ⁴⁸Ca from the IM-GCM calculation using either the bare or evolved transition operator, where the renormalization effect enlarges significantly the short-range contribution for both transitions. It has been discussed in Ref. [42] that the multireference IMSRG flow incorporates the effects of pairing in high-energy orbitals, greatly enhancing the contribution of the J = 0 pair of nucleons to the NMEs. We note that one should not make a direct comparison of the operator renormalization effects in the VS-IMSRG and IM-GCM calculations, as the operator renormalization and many-body correlations are partitioned in a different way. It is, however, meaningful to compare the final matrix elements obtained with renormalized operators; in this case, the discrepancy between the results of these two calculations reflects the error due to the missing of higherbody operators in both methods. As shown in Ref. [41] the inclusion of an induced three-body transition operator helps reduce the discrepancy between them. Therefore, these two variants of IMSRG provide a complementary description of NMEs of neutrinoless double- β decay.

VI. CONCLUSION

In this work, we have explored the possible correlation between the NMEs $M^{0\nu\beta\beta}$ of ground-state-to-ground-state $0\nu\beta\beta$ decay and those M^{DGT} of DGT transitions in a set of nuclei in different mass regions with three *ab initio* methods starting from the same NN+3N chiral interactions. We have found that the obtained M^{DGT} is correlated with the quantity $M^{0\nu\beta\beta}A^{-1/3}$ slightly stronger than with $M^{0\nu\beta\beta}A^{-1/6}$ for isospin-conserving transitions, where the long-range and short-range contributions add coherently, leading to large values of both transition matrix elements. However, the correlation relation turns out to be much weaker in isospin-changing transitions where the long-range and short-range contributions compensate each other, leading to small values of M^{DGT} . This conclusion has been confirmed with a scale-separation analysis in which we have sampled a set of transition densities for both isospinconserving and isospin-changing transitions by changing the short-range and long-range behavior around the results from the VS-IMSRG calculation.

We have also explored the origin of the discrepancy between the NMEs from conventional shell-model calculations and VS-IMSRG calculations. Our studies have shown that apart from the discrepancy mainly due to the configuration mixing predicted differently in the calculations using the conventional shell-model interactions and those derived from VS-IMSRG, the in-medium renormalization effect from the VS-IMSRG evolution on the transition operator varies with isotopes, which spoils somewhat the correlation relation.

Combining the NMEs of both isospin-conserving and isospin-changing transitions from the three ab initio calculations, we observe a strong correlation. However, the correlation is considerably weaker for the experimentally relevant isospin-changing transitions. In other words, a large uncertainty will likely still exist in the $M^{0\nu\beta\beta}$ even if the ground-state-to-ground-state DGT transition of the candidate nucleus is precisely measured. It is worth mentioning that our current analysis is mainly based on the results of calculations with the chiral interaction EM1.8/2.0. A comprehensive way to examine the correlation relation can be carried out by computing the NMEs with a set of chiral nuclear forces with the low-energy constants varying within acceptable regions [92]. Any experimental data on the DGT transition NME may provide a constraint on the chiral interaction, and through that reduce the uncertainty of the predicted $M^{0\nu\beta\beta}$. Besides, we note that recently Romeo et al. [93] found a good linear correlation between the $0\nu\beta\beta$ decays and the double- γ transitions, suggesting another potential way to constrain the NMEs of $0\nu\beta\beta$ decay. The present study can be extended straightforwardly to examine that linear correlation relation as well.

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