# First calculation of the $\gamma \gamma$-decay width of a nuclear $2_{1}^{+}$state: The case of ${ }^{48} \mathbf{C a}$ 

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#### Abstract

The competitive double $-\gamma$ decay of the $2_{1}+$ state of an even-even spherical nucleus is studied for the first time. The coupling of one-, two-, and three-phonon terms in the wave functions of excited states is taken into account within the microscopic model based on the Skyrme energy density functional. The approach enables one to perform the calculations in very large configurational spaces. We estimate the generalized electric dipole polarizabilities involved in the $\gamma \gamma / \gamma$ decay process and make a prediction for the branching ratio of the competitive $\gamma \gamma$-decay relative to its single $\gamma$-decay calculated to be $3 \times 10^{-8}$ for the case of ${ }^{48} \mathrm{Ca}$.


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## I. INTRODUCTION

In Ref. [1], the $\gamma \gamma$ decay of a nuclear transition in competition with an allowed $\gamma$ decay has been discovered. This is the observation of the $\gamma \gamma$ decay of the first excited $J^{\pi}=11 / 2^{-}$ state of ${ }^{137} \mathrm{Ba}$ directly competing with an allowed $\gamma$ decay to the $J^{\pi}=3 / 2^{+}$ground state. The branching ratio of the competitive $\gamma \gamma$ decay of the $11 / 2^{-}$isomer of the odd-even nucleus ${ }^{137} \mathrm{Ba}$ to the ground state relative to its single $\gamma$ decay was determined to be $(2.05 \pm 0.37) \times 10^{-6}$. This discovery has very recently been confirmed and the data were made more precise, in particular with respect to the contributing multipolarities [2].

The $\gamma \gamma$-decay reactions are formally analogous to neutrinoless double- $\beta$ decay ( $0 \nu \beta \beta$-decay) processes where in the latter two $\beta$ particles and in the former two $\gamma$-quanta appear in the final state and share the total transition energy. Indeed, $\gamma \gamma$-decay processes have first theoretically been postulated and studied by Maria Göppert in her Ph.D. thesis [3] with Max Born in Göttingen, even before discussing $\beta \beta$-decay processes. While various predictions for the $0 \nu \beta \beta$-nuclear matrix elements are on the market, e.g., from Refs. [4-6], there is no way to test the accuracy of the theoretical calculations without a firm measurement of the $0 \nu \beta \beta$-decay rate. The $2 \nu \beta \beta$-decay rate is a useful tool for testing the accuracy of calculations that try to predict the $0 \nu \beta \beta$ rate, but it is certainly not enough for making predictions for the $0 \nu \beta \beta$ rates [5]. Although $\beta \beta$ decay processes are nuclear reactions of second order in the electroweak interaction, it is surprising to find that even less data exist for nuclear decay reactions that proceed in second order in the electromagnetic interaction where two $\gamma$ quanta are simultaneously emitted in a single quantum transition from one quantum state to another. Up to recently, $\gamma \gamma$-decay reactions in the even-even nuclei were known only in three particular cases, ${ }^{16} \mathrm{O}[7,8]$ and ${ }^{40} \mathrm{Ca}$ and ${ }^{90} \mathrm{Zr}[9,10]$, where the first excited states of these even-even nuclei have spin and
parity quantum numbers $0^{+}$and a single- $\gamma$ decay is strictly forbidden by helicity conservation.

The $E 1 E 1$ transition from the $2 s$ to $1 s$ level in the H atom and low- Z H -like ions is a proverbial example for atomic systems. It is the most probable decay mode and therefore defines the lifetime of the $2 s$ level. The first estimate was obtained by Breit and Teller in Ref. [11]. However, the nuclear $\gamma \gamma$ decay dominated by $E 1 E 1$ contribution has not yet been found.

In this paper, we report on the more general situation, in which the $\gamma \gamma$ decay of the low-energy quadrupole state occurs in a nuclear transition which could proceed by a single- $\gamma$ decay in competition.

Using $\hbar=c=1$, the $\gamma$ decay width of the $2_{1}{ }^{+}$state of even-even nuclei is related to its reduced electric quadrupole transition strength, $B\left(E 2 ; 2^{+}{ }_{1} \rightarrow 0^{+} \mathrm{gs}\right)$, via

$$
\begin{equation*}
\Gamma_{\gamma}=\frac{4 \pi}{75}\left(E_{2^{+}}\right)^{5} B\left(E 2 ; 2_{1}^{+} \rightarrow 0^{+}{ }_{\mathrm{gs}}\right) \tag{1}
\end{equation*}
$$

To describe the $\gamma \gamma$ decay between the $2_{1}{ }^{+}$and $0_{\mathrm{gs}}{ }^{+}$states, we use a formalism that explicitly relates the electromagnetic interaction up to second order in the electromagnetic operators and two-quantum processes in atomic nuclei [12]. Thus, the $\gamma \gamma$-decay width can be estimated as

$$
\begin{equation*}
\Gamma_{\gamma \gamma^{\prime}}=\frac{64 \pi}{42525}\left(E_{2^{+}, 1}\right)^{7}\left(\alpha_{E 1 E 1}\right)^{2}(1+\delta) \tag{2}
\end{equation*}
$$

with

$$
\begin{align*}
& \alpha_{E 1 E 1}=\sum_{i} \frac{\left\langle 0_{\mathrm{gs}}^{+}\right||M(E 1)|\left|1_{i}^{-}\right\rangle\left\langle 1_{i}^{-}\|M(E 1)\| 2^{+}{ }_{1}\right\rangle}{E_{1_{i}^{-}}-0.5 E_{2^{+}}}  \tag{3}\\
& \delta=\left(\frac{\alpha_{M 1 M 1}}{\alpha_{E 1 E 1}}\right)^{2}+\frac{3}{11} 10^{-4}\left(\frac{\alpha_{E 2 E 2}}{\alpha_{E 1 E 1}}\right)^{2}\left(E_{2^{+}}{ }_{1}\right)^{4}+\cdots \tag{4}
\end{align*}
$$

The $\gamma \gamma$-decay width is dominated by the $E 1 E 1$ contribution, i.e., $\delta \ll 1$. One may easily confirm this statement in
a two-state scenario by using the $M 1$ sum rule. The $M 1 M 1$ component to the $\gamma \gamma / \gamma$ decay of the $2_{1}{ }^{+}$state of any nucleus is much smaller than the $\alpha_{E 1 E 1}$ component discussed here in the case of ${ }^{48} \mathrm{Ca}$, taken as an example.

The doubly magic nucleus ${ }^{48} \mathrm{Ca}$ has provided a crucial testing ground for nuclear theories (see, e.g., Refs. [13,14]). In this connection, the electric dipole polarizability,

$$
\begin{equation*}
\alpha_{D}=\frac{8 \pi}{9} \sum_{i} \frac{\left\langle 0_{\mathrm{gs}}^{+}\|M(E 1)\| 1_{i}^{-}\right\rangle\left\langle 1_{i}^{-}\|M(E 1)\| 0_{\mathrm{gs}}^{+}\right\rangle}{E_{1_{i}^{-}}} \tag{5}
\end{equation*}
$$

is playing an important role; in particular, its value has strong implications in constraining the symmetry energy $J$ including its density dependence and slope parameter $L$ of the nuclear equation of state [15]. The symmetry energy also plays an important role in nuclei, where it contributes to the formation of neutron skins in the presence of a neutron excess. Calculations based on energy density functionals (EDFs) pointed out that $J$ and $L$ can be correlated with isovector collective excitations of the nucleus, such as pygmy dipole resonances (PDRs) [16] and giant dipole resonances (GDRs) [17], thus suggesting that the neutron skin thickness [18], the difference of the neutron and proton root-mean-square radii, could be constrained by studying properties of collective isovector observables at low energy. Since $\alpha_{D}$ and $\alpha_{E 1 E 1}$ are challenging but in principle accessible observables, it is useful to compare their values. We provide here first information on their relation.

## II. FORMALISM

Our tool is the random phase approximation (RPA) with Skyrme EDF. The residual particle-hole interaction is obtained as the second derivative of the energy density functional with respect to the particle density. By means of the standard procedure [19], we obtain the familiar equations of the RPA in the one particle-one hole ( $1 \mathrm{p}-1 \mathrm{~h}$ ) configuration space. The eigenvalues of the RPA equations are found numerically as the roots of a relatively simple secular equation within the finite-rank separable approximation (FRSA) [20,21] which allows one to perform the calculations in large configurational spaces. In particular, the cutoff of the discretized continuous part of the single-particle spectra is at the energy of 100 MeV . This is sufficient to exhaust practically all the energyweighted sum rule within the RPA [21]. Further technical details of the calculations are provided in Appendix A.

Being a linear combination of $1 \mathrm{p}-1 \mathrm{~h}$ states, the RPA solutions are treated as quasibosons with quantum numbers $\lambda^{\pi}: Q_{\lambda \mu i}^{+}|0\rangle$. The value $\mu$ denotes the $z$ projection of the total angular momentum in the laboratory system. Among these solutions, there are one-phonon states corresponding to collective GDRs and pure $1 \mathrm{p}-1 \mathrm{~h}$ states. The configurations with various degrees of complexity can be built by combining different one-phonon configurations of fixed quantum number $J^{\pi}=1^{-}$. Taking into account the basic ideas of the quasiparticle-phonon model (QPM) [22], the Hamiltonian is then diagonalized in a space spanned by states composed of one, two, and three RPA phonons. This implies that the rank of the set of linear equations is equal to the number of one-, two-, and three-phonon configurations included in the wave
function,

$$
\begin{align*}
\Psi_{v}(J M)= & \left(\sum_{i} R_{i}(J v) Q_{J M i}^{+}+\sum_{\substack{\lambda_{1} i_{1} \\
\lambda_{2} i_{2}}} P_{\lambda_{2} i_{2}}^{\lambda_{1} i_{1}}(J v)\left[Q_{\lambda_{1} i_{1}}^{+} Q_{\lambda_{2} i_{2}}^{+}\right]_{J M}\right. \\
& \left.+\sum_{\substack{\lambda_{1} i_{1} \\
\lambda_{1} i_{2} \\
\lambda_{3} i_{3} J^{\prime}}} T_{J^{\prime} \lambda_{3} i_{3}}^{\lambda_{1} i_{1} \lambda_{2} i_{2}}(J v)\left[\left[Q_{\lambda_{1} i_{1}}^{+} Q_{\lambda_{2} i_{2}}^{+}\right]_{J^{\prime}} Q_{\lambda_{3} i_{3}}^{+}\right]_{J M}\right)|0\rangle
\end{align*}
$$

Its solution requires computation of the matrix elements of the quasiparticle-phonon interaction [23], $\langle 0| Q_{J i} H\left[Q_{\lambda_{1} i_{1}}^{+} Q_{\lambda_{2} i_{2}}^{+}\right]_{J}|0\rangle$. The equations of the phonon-phonon coupling (PPC) have the same form as the QPM equations [24,25], but the single-particle spectrum and the parameters of the residual interaction are calculated with the chosen Skyrme EDF without any further adjustments [23]. We consider widely used SLy5 EDF [26], which is adjusted to reproduce the enhancement factor of the Thomas-Reiche-Kuhn (TRK) sum rule $\kappa=0.25$ and the nuclear matter properties, as well as nuclear charge radii and binding energies of doubly magic nuclei [27]. Also the excitation energy $E_{x}=3.832 \mathrm{MeV}$ and the decay transition strength $B\left(E 2 ; 2^{+}{ }_{1} \rightarrow 0_{\mathrm{gs}}^{+}\right)=1.71 \pm 0.09$ W.u. of the $2_{1}{ }^{+}$state of ${ }^{48} \mathrm{Ca}$ [28] are reasonably well described in the RPA with the same set of parameters, yielding $E_{x}=3.19 \mathrm{MeV}$ and $B(E 2)=1.3$ W.u. [29]. The inclusion of the phonon-phonon coupling plays a minor role in the $2_{1}{ }^{+}$state's description. The crucial contribution to the wave function comes from the neutron configuration $\left\{1 f_{7 / 2}^{-1}, 2 p_{3 / 2}\right\}$.

In the actual calculations (PPC3), we have included in our model space different multipoles $\lambda^{\pi}=1^{-}, 2^{+}, 3^{-}$, and $4^{+}$. Tentative estimates for the position of the resonance centroid $E_{c}$ and the spreading width $\Gamma$ are defined by means of the energy-weighted moments $m_{k}=\sum B(E 1) E^{k}$ : (i) $E_{\mathrm{c}}=$ $m_{1} / m_{0}$ and (ii) $\Gamma=\sqrt{m_{2} / m_{0}-\left(m_{1} / m_{0}\right)^{2}}$, taking $92 \%$ of the TRK sum rule, i.e., $(1+\kappa) 14.8 N Z / A e^{2} \mathrm{fm}^{2} \mathrm{MeV}$. The $E 1$ transitions are corrected for the center-of-mass motion; see Appendix B. All one-, two-, and three-phonon configurations with energies up to $E=27 \mathrm{MeV}$ are included. The inclusion of high-energy configurations plays a minor role in our calculations. At the same time, the extension of the configurational space to the two- and three-phonon configurations has a strong effect on the low-energy $1^{-}$spectrum, shown in Fig. 1. If we omit the three-phonon configurations, then this calculation is hereafter called PPC2. Further, the PPC2 calculation taking into account the coupling with $\left[2_{1}{ }^{+}\right]_{\text {RPA }}$ phonon only is named PPC2-2 ${ }_{1}{ }^{+}$.

## III. RESULTS

The photoabsorption cross section in ${ }^{48} \mathrm{Ca}\left(p, p^{\prime}\right)$ is measured in the range from 10 to 25 MeV [32]. The general shape of the $E 1$ strength distribution obtained in the PPC3 calculation is rather close to that observed in experiment; see Fig. 2(a). The calculated integral characteristics of the $\operatorname{GDR}\left(E_{c}=19.1 \mathrm{MeV}\right.$ and $\left.\Gamma=3.1 \mathrm{MeV}\right)$ are in


FIG. 1. The phonon-phonon coupling effect on the $B(E 1)$ strength distribution. Experimental data are taken from Ref. [30].
satisfactory agreement with the experimental data on ${ }^{48} \mathrm{Ca}$ [32]. The inclusion of the PPC effects yields a noticeable redistribution of the GDR strength in comparison with the RPA results. In Ref. [29], the role of the coupling of the one- and


FIG. 2. (a) The $B(E 1)$ strength distribution. The smoothing parameter 1 MeV is used for the strength distribution described by the Lorentzian function. Experimental data are taken from Refs. [31,32]. The dashed, dotted, and solid lines correspond to the PPC2-2 ${ }^{+}$, PPC2, and PPC3 calculations, respectively. (b) Running sum of the polarizability $\alpha_{D}$. The shaded area indicates experimental upper and lower limits [32]. (c) Running sum of the generalized polarizability $\alpha_{E 1 E 1}$ and the value $M_{E 1 E 1} / E_{s c}$ (the dash-dotted line).
two-phonon configurations on the $E 1$ strength-distribution calculations with the SLy5 EDF has been examined.

The electric dipole polarizability of $\alpha_{D}=1.73 \pm 0.18 \mathrm{fm}^{3}$ has been measured in the range from 10 to 25 MeV in Ref. [32]. Running sums of the $\alpha_{D}$ value for ${ }^{48} \mathrm{Ca}$ in the energy region below 27 MeV are given in Fig. 2(b). The steep rise in the theoretical band around 19 MeV indicates that the position of the GDR peak is consistent with the experimental centroid. The fact that the slope is well reproduced indicates that the spreading is well described. Inclusion of PPC does not change the value of $\alpha_{D}$ obtained by integrating the $E 1$ strength up to $25 \mathrm{MeV}: \alpha_{D}=2.16 \mathrm{fm}^{3}$ in the case of the PPC2 and $\alpha_{D}=2.17 \mathrm{fm}^{3}$ in the PPC3. The contribution [0.0101 $\pm$ $0.0006 \mathrm{fm}^{3}$ ] found experimentally below 10 MeV is negligible. Our calculations predict, below $10 \mathrm{MeV}, \alpha_{D}=0.015 \mathrm{fm}^{3}$ and $0.017 \mathrm{fm}^{3}$ within PPC2 and PPC3, respectively. These values are indeed negligible, as for the experimental case. A similar conclusion was made in Ref. [33], where the EDFs SGII and SLy4 were used. Thus, the main contribution to the $\alpha_{D}$ polarizability comes from the one-phonon configurations.

The calculated value for the $\alpha_{E 1 E 1}$ matrix element for the $\gamma \gamma$ decay is $0.018 \mathrm{fm}^{3}$ within the PPC2 and $0.019 \mathrm{fm}^{3}$ in the case of the PPC3 at $1^{-}$energies below 27 MeV . The running sum of $\alpha_{E 1 E 1}$ is very instructive with respect to various energy regions of the $1^{-}$spectrum; see Fig. 2(c). We obtain two regimes: a rapid rise in the energy region $16<E_{x}<20 \mathrm{MeV}$ and a steep decline in the energy region $22<E_{x}<25 \mathrm{MeV}$, respectively. In the first energy region, the crucial component to the $\alpha_{E 1 E 1}$ value comes from a coherent contribution of $1^{-}$ states in the sum of product of $E 1$ matrix elements

$$
\begin{equation*}
M_{E 1 E 1}=\sum_{i}\left\langle 0_{\mathrm{gs}}^{+}\|M(E 1)\| 1_{i}^{-}\right\rangle\left\langle 1_{i}^{-}\|M(E 1)\| 2^{+}{ }_{1}\right\rangle \tag{7}
\end{equation*}
$$

belonging to the GDR; see Fig. 3(a). The running sum for $\alpha_{E 1 E 1}$ decreases at energies above 22 MeV due to a coherent contribution with a negative sign in this energy range; see Fig. 3(b). This implies that the calculated $E 1$ matrix elements represent two dominant substructures of the dipole strength distribution; see the dash-dotted line of Fig. 2(c). We study them first in a simple two-state mixing scenario taking into account the GDR and the initial, $2_{1}{ }^{+}$, and final, $0^{+}{ }_{\mathrm{gs}}$, states of the generalized polarizability.

We construct a two-state mixing scheme by considering the GDR state, the coupled GDR $\otimes 2_{1}{ }^{+}$state, and the interaction $V$ between them. The relative phases of amplitudes are opposite in the perturbed states I and II, i.e.,

$$
\begin{align*}
\left|1_{\mathrm{I}}^{-}\right\rangle & =\alpha|\mathrm{GDR}\rangle+\beta\left|\mathrm{GDR} \otimes 2_{1}^{+}\right\rangle  \tag{8}\\
\left|1_{\mathrm{II}}^{-}\right\rangle & =-\beta|\mathrm{GDR}\rangle+\alpha\left|\mathrm{GDR} \otimes 2_{1}{ }^{+}\right\rangle \tag{9}
\end{align*}
$$

The two-state mixing scheme provides the simple expression

$$
\begin{equation*}
\frac{\alpha_{E 1 E 1}}{\alpha_{D}}=\frac{9}{8 \pi} \frac{\alpha \beta\left(E_{\mathrm{II}}-E_{\mathrm{I}}\right)}{\alpha^{2} E_{\mathrm{II}}+\beta^{2} E_{\mathrm{I}}}, \tag{10}
\end{equation*}
$$

in the typical situation where the energies $E_{\text {II }}$ and $E_{\mathrm{I}}$ are much large than $E_{x}\left(2_{1}{ }^{+}\right) / 2$. Notice that $\alpha_{E 1 E 1} / \alpha_{D}=0$ if the basis configurations do not mix. The amplitudes, $\alpha$ and $\beta$, can be estimated in the two limiting cases of weak or strong mixing, where either the interaction V is much smaller than


FIG. 3. Reduced matrix elements of $E 1$ transitions in the following energy interval: from 16 until 20 MeV (a) and from 22 until 25 MeV (b).
the difference of the unperturbed energies $\left[\Delta E_{u}=E_{x}\left(2_{1}{ }^{+}\right)\right]$, or the two unperturbed states are degenerate $\left(\Delta E_{u}=0\right)$. For these cases, we obtain

$$
\begin{equation*}
\frac{\alpha_{E 1 E 1}}{\alpha_{D}}=\frac{9}{8 \pi} \frac{V}{E_{\mathrm{GDR}}+0.5 \Delta E_{u}} \tag{11}
\end{equation*}
$$

Figure 2(c) shows that the running sum of $\alpha_{E 1 E 1}$ remains more or less unchanged when comparing the full result with the calculation taking into account the PPC with the $\left[2_{1}{ }^{+}\right]_{\text {RPA }}$ phonon, only. Therefore, we can simulate the full diagonalization reasonably well by using a two-state mixing [34]; see Fig. 4. The GDR state built on the most collective RPA states, the two-phonon GDR $\otimes 2_{1}{ }^{+}$state, and the interaction

$$
\begin{equation*}
V=\frac{1}{\sqrt{N_{1 \mathrm{ph}} N_{2 \mathrm{ph}}}} \sum_{i=1}^{N_{\mathrm{lph}}} \sum_{k=1}^{N_{2 \mathrm{ph}}}\langle 0| Q_{1 i} H\left[Q_{21}^{+} Q_{1 k}^{+}\right]_{1}|0\rangle, \tag{12}
\end{equation*}
$$

between them, have been calculated from the SLy5 EDF.
The two-state model can also help to estimate the value $\alpha_{M 1 M 1}$ of $0.05 \mu_{N}^{2} \mathrm{MeV}^{-1}$; see Appendix C. The $M 1$ transition matrix elements calculated without spin-gyromagnetic quenching factor leads to an upper limit for $\alpha_{M 1 M 1}$. The M1M1 component dominates in the correction $\delta$ of Eq. (4): $\delta(M 1)=1.7 \times 10^{-3}$ and $\delta(E 2)=1.9 \times 10^{-7}$. Thus, keeping


FIG. 4. The ratio between polarizabilities $\alpha_{D}$ and $\alpha_{E 1 E 1}$ as a function of the mean value $V$ of the matrix elements coupling one- and two-phonon configurations. The solid line corresponds to the PPC2-2 ${ }^{+}$calculation taking into account scaling of the quasiparticle-phonon interaction. The results of the two-state model with $\Delta E_{u}=0$ and 3.2 MeV are denoted by the dashed line and the dotted line, respectively. The PPC3 calculation of SLy5 is denoted by the circle.
only $1^{-}$virtual states is a satisfactory approximation for the $\gamma \gamma$ decay width. It is worth mentioning that in ${ }^{48} \mathrm{Ca}$ the spin-flip $M 1$ resonance having a noncollective structure being dominated by the neutron configuration $\left\{1 f_{7 / 2}^{-1}, 1 f_{5 / 2}\right\}$ is a key reference for an interpretation of the quenching phenomenon; see, e.g., Ref. [35].

In total, our PPC3 analysis predicts the $\gamma \gamma$-decay width of $1.0 \times 10^{-10} \mathrm{eV}$, almost seven orders smaller than the width corresponding to the single $\gamma$ decay. This may be considered as an upper limit of the $\gamma \gamma / \gamma$ branching ratio since the calculated width of the $\gamma$ decay, $3.5 \times 10^{-3} \mathrm{eV}$, is smaller than the experimental data [28].

## IV. SUMMARY AND CONCLUSION

Starting from Skyrme mean-field calculations we have studied for the first time the $\gamma \gamma / \gamma$ decay of the $2{ }_{1}{ }^{+}$state of an even-even nucleus. As our test case we considered ${ }^{48} \mathrm{Ca}$ for which its dipole polarizability has recently been measured. It can be compared also to the generalized dipole polarizability. We use the Skyrme EDF SLy5 to create a single-particle spectrum and to analyze excited states of ${ }^{48} \mathrm{Ca}$. Our calculations take into account the coupling between one-, two- and three-phonon terms in the wave functions. It is shown that the $\gamma \gamma$ decay width is sensitive to the energy difference of the one-phonon doorway mode and the two-phonon structure which arises from the coupling of this doorway to the lowenergy phonons. It is further correlated to the PPC strength. The maximal branching ratio of the competitive $\gamma \gamma$-decay relative to its single $\gamma$-decay is predicted for ${ }^{48} \mathrm{Ca}$ as $3 \times 10^{-8}$. It is desirable to experimentally establish the $\gamma \gamma$ decay of a first $2^{+}$state of an even-even nucleus.

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FIG. 5. The transition probability, the $\gamma$-decay width, and $\gamma \gamma / \gamma$ branching ratio of the $2_{1}{ }^{+}$state are given in panels (a), (b), and (c), respectively. PPC3 calculations are performed with different cutoffs of single-particle energies ( $E_{\mathrm{sp}}$ ). The solid lines are drawn as guide to the eye.
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## APPENDIX A

The FRSA model enables one to exhaust the Thomas-Reiche-Kuhn (TRK) sum rule with the enhancement factor $\kappa$ for the Skyrme EDF, i.e., $(1+\kappa) 14.8 N Z / A \mathrm{e}^{2} \mathrm{fm}^{2} \mathrm{MeV}$. One of the basic ingredients for the fitting protocol of the SLy5 EDF is the enhancement factor of the TRK sum rule $\kappa=0.25$ [26], i.e., $216 \mathrm{e}^{2} \mathrm{fm}^{2} \mathrm{MeV}$ in the case of ${ }^{48} \mathrm{Ca}$. The energy-weighted sum rule of the $E 1$ distribution calculated within the FRSA amounts to $210 \mathrm{e}^{2} \mathrm{fm}^{2} \mathrm{MeV}$. The model configuration space depends on the cutoff of the discretized continuous part of the single-particle spectra. To confirm the two-state scenario of the $\alpha_{E 1 E 1}$ calculation, one needs a maximum full configurational space. The phonon-phonon coupling calculations (PPC3) take into account the different values of single-particle energy cutoff; see Fig. 5. For the $\gamma \gamma / \gamma$-decay branching, satisfactory convergence is reached from $E_{\mathrm{sp}} \approx 50$ MeV . Thus, the cutoff at 100 MeV is sufficient to describe correctly the $\gamma \gamma / \gamma$-decay branching ratio.

## APPENDIX B

To calculate the dipole strength distributions, the spurious isoscalar dipole mode appears at zero excitation energy. Because of small numerical inaccuracies, the $E 1$ transitions are corrected for the center-of-mass motion. The effective charges $-Z / A$ for neutrons and $N / A$ for protons eliminate contaminations associated with the operator,

$$
\begin{equation*}
\hat{S}=\sum_{i=1}^{A} r_{i} Y_{1 \mu}\left(\hat{r}_{i}\right) \tag{B1}
\end{equation*}
$$

To crosscheck the reliability of the prediction of $\alpha_{E 1 E 1}$, we use the orthogonalization of the spurious state to all physical states, as proposed in Ref. [36]. Starting from the wave functions (6) defined as the set $|\nu\rangle$, we construct a new set of normalized states $|\tilde{v}\rangle$,

$$
\begin{equation*}
|\tilde{v}\rangle=\mathcal{N}_{\tilde{v}}\left(|\nu\rangle-\eta_{\tilde{v}}|S\rangle\right), \tag{B2}
\end{equation*}
$$

where the state $|S\rangle$ is defined as $|S\rangle \equiv \hat{S}|0\rangle$, and $|0\rangle$ being the RPA vacuum. The set $|\tilde{v}\rangle$ is required to satisfy the following


FIG. 6. (a) The $B(E 1)$ strength distribution. (b) Running sum of the polarizability $\alpha_{D}$. (c) Running sum of the generalized polarizability $\alpha_{E 1 E 1}$. The dashed line corresponds to the calculation taking into account the orthogonalization of the spurious center-of-mass state to all physical states, the dotted line is the calculation without the orthogonalization, and the solid line is the calculation with the effective charges, $e_{\text {eff }}^{n}=-\frac{Z}{A} e$ and $e_{\text {eff }}^{p}=\frac{N}{A} e$.
conditions:

$$
\begin{equation*}
\langle\tilde{v}| \hat{S}|0\rangle=0 . \tag{B3}
\end{equation*}
$$

As shown in Fig. 6, we found a remarkable agreement between the running sum of the generalized polarizability $\alpha_{E 1 E 1}$, obtained with the orthogonalization, and the ones, generated by means of the effective charges. Moreover, the calculation without the orthogonalization leads to very similar results for the value of $\alpha_{E 1 E 1}$. This means that the spurious center-of-mass state is well separated and it plays a minor role in our calculation of $\alpha_{E 1 E 1}$. Nevertheless, the $E 1$ transitions are corrected for the center-of-mass motion.

## APPENDIX C

For estimating the value of the generalized polarizability $\alpha_{M 1 M 1}$, we take the spin-flip $M 1$ resonance $(G R)$ and the initial
$\left(2_{1}{ }^{+}\right)$and final $\left(0_{\mathrm{gs}}^{+}\right)$states into account. The two-state mixing scheme then consists of the GR state, the coupled GR $\otimes$ $2_{1}{ }^{+}$state, and the interaction $V$ between them. The model parameters are defined by means of the energy-weighted moments $m_{k}=\sum B(M 1) E^{k}$ : the centroid, $E_{\mathrm{GR}}=m_{1} / m_{0}$, and the $M 1$-transition probability, $B(M 1)_{\mathrm{GR}}=m_{0}$. The mixed wave functions are

$$
\begin{align*}
\left|1_{\mathrm{I}}^{+}\right\rangle & =\alpha|\mathrm{GR}\rangle+\beta\left|\mathrm{GR} \otimes 2_{1}{ }^{+}\right\rangle  \tag{C1}\\
\left|1_{\mathrm{II}}{ }^{+}\right\rangle & =-\beta|\mathrm{GR}\rangle+\alpha\left|\mathrm{GR} \otimes 2_{1}{ }^{+}\right\rangle \tag{C2}
\end{align*}
$$

The two-state model gives the following expression:

$$
\begin{equation*}
\alpha_{M 1 M 1}=B(M 1)_{\mathrm{GR}} \frac{\alpha \beta \sqrt{E_{2_{1}+}^{2}+4 V^{2}}}{E_{\mathrm{GR}}^{2}-0.25 E_{2_{1}+}^{2}-V^{2}} \tag{C3}
\end{equation*}
$$

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