# **Reliability of Calculated Nuclear Matrix Element of Two-Neutrino Double-ß decay**



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The goal is to determine the effective mass of the neutrino using the neutrinoless double- $\beta$  (0v $\beta\beta$ ) decay of the nuclei. A problem was reported of a significant discrepancy in the running sum for the nuclear matrix element (NME) of the two-neutrino double- $\beta$  (2v $\beta\beta$ ) decay between the shell model and QRPA calculations. The reliability of the calculation of the NME of the 2vßß decay is important because of the similarity of the  $0v\beta\beta$  and  $2v\beta\beta$  decays. This problem is solved in this study.

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Status:

#### Investigation of cause



Possible change of two neutrons to two protons in a nucleus emitting two electrons with neutrino exchange (neutrinoless double- $\beta$  (0v $\beta\beta$ ) decay). This decay occurs, if the neutrino (v) is a Majorana particle ( $\nu = \overline{\nu}$ ), and the effective neutrino mass can be determined, see the equations below. Determination of the effective neutrino mass is one of the most important subjects in modern physics.

#### Why nuclei?

Because E(final state) < E(initial state) is necessary.

List of nuclei used in the experiments

<sup>76</sup> Ge→ <sup>76</sup> Se	$^{130}\text{Te} \rightarrow ^{130}\text{Xe}$	<sup>136</sup> Xe→ <sup>136</sup> Ba
<sup>150</sup> Nd→ <sup>150</sup> Sm	<sup>48</sup> Ca→ <sup>48</sup> Ti	<sup>82</sup> Se→ <sup>82</sup> Kr
<sup>96</sup> Zr→ <sup>96</sup> Mo	<sup>100</sup> Mo→ <sup>100</sup> Ru	$^{110}\text{Pd} \rightarrow ^{110}\text{Cd}$
<sup>116</sup> Cd→ <sup>116</sup> Sn	$^{124}Sn \rightarrow ^{124}Te$	and more

The calculated NMEs by various approximation methods and groups are distributed in a range of the min.-max. ratio of 2–3. The NME cannot be obtained by experiment. Thus, examination and improvement of the calculation are essential.

#### The problem

The large discrepancy in the running sum for  $2\nu\beta\beta$  NME  $\times (g_A)^2$  of  ${}^{136}Xe \rightarrow {}^{136}Ba$  in two calculations. ×0.01



The intermediate nucleus is <sup>136</sup>Cs. In calculation, effective value  $g_A^{eff}$  is used for  $g_A$ . SM: shell model; GCN and MC: interaction names. QRPA: quasiparticle random-phase approximation. The calculated  $M_{GT}^{(2v)}$  (the right edge) are fitted to the

• It turned out that a small decrease is obtained at the energy of the GT<sup>-</sup> giant resonance (GR<sup>-</sup>) at 12.5 MeV by increasing  $G_{\text{pn}}^{\text{IS}}$ .



• The decrease in the running sum implies that the component of the  $2\nu\beta\beta$  NME changes its sign at that energy. This behavior can be explained by the large strength of the interaction analytically under the separable approximation to the RPA.





 $C_{\mu i}$ : matrix element of the one-body operator used for the interaction of the separable approximation. The particle and hole states are denoted by  $\mu$  and *i*.

 $\langle m_{\nu} \rangle = \left| \sum_{i=1,2,3} U_{ei}^2 m_i \right|$ 

*U*: Pontecorvo-Maki-Nakagawa-Sakata matrix  $m_i$ : eigen mass (*i*=1,2,3)



exp. value.

GT: Gamow-Teller; the Fermi NME is negligible for the  $2\nu\beta\beta$  decay.



### My QRPA calculation

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Skyrme + Coulomb + contact isovector (IV) and isoscalar (IS) pairing (*pp*, *nn*, and *pn*) interactions. *p*: proton; *n*: neutron  $G_{pn}^{IS}$ : strength of the IS *pn* pairing interaction.

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	Menéndez	Horoi	Šimkovic	Terasaki	
Method	SM		QRPA		

 $\varepsilon_{\mu i}$ : particle-hole energy of main component of GR<sup>-</sup>.

 $E_B$ : energy of intermediate state  $|B\rangle$ .  $\overline{M}$ : mean value of initial and final nuclear masses.

 $N_B$ : normalization factor > 0.

When the attractive interaction strength is enhanced, all  $E_B$  are lowered and the one nearest to the  $\varepsilon_{\mu i}$  of the GR<sup>-</sup> switches from  $E_B < \varepsilon_{\mu i}$  to  $E_{B'} > \varepsilon_{\mu i}$ . This causes the sign change of  $\langle B(GR^{-}) | \sigma \tau^{-} | I \rangle$ .  $O_{B_{CR^{-}}}^{F^{\dagger}} | F \rangle$  does not have this component in the single-Slater approximation; see the figure below. Thus, the possibility of the sign change of  $\langle F | \boldsymbol{\sigma} \tau^{-} | B(GR^{-}) \rangle$  is low.



# GT<sup>-</sup>: GT transition from *n* to *p*.; GT<sup>+</sup>: that from *p* to *n*.

The significant decrease in the running sum at the GT<sup>-</sup> GR implies that the interaction is stronger than that of calculations with less decrease.

The cause of the problem is the difference in the interaction strength. **Conclusion 203** 



The transition operator used in my calculation is  $V(r_{12}, E_B) \cong h_+(r_{12})\{-\boldsymbol{\sigma}(1) \cdot \boldsymbol{\sigma}(2) + g_V^2/g_A^2\} \tau^-(1)\tau^-(2)$ Double-Gamow-Teller + Double-Fermi  $g_{V}$ : vector current coupling Neutrino potential = 1

Variation of				
comp. of	Small	Large	Large	Small

The calculations of Menéndez et al. (SM) and Simkovic et al. (QRPA) are those in the above figure (Gando et al.). The calculation of Horoi et al. is in

M. Horoi and A. Brown, PRL **110**, 222502 (2013)

The cause of the discrepancy problem is not the theoretical differences between SM and QRPA.

## Check of my interaction strength

Exp. data and calculation of  $g_A^2 \times GT^-$  strength



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The energy dependence of the strength shows the validity of my interaction.