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Microscopic calculation of the β^- decays of ^{151}Sm , ^{171}Tm , and ^{210}Pb with implications to detection of the cosmic neutrino background



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ABSTRACT

The electron spectral shapes corresponding to the low- Q β^- -decay transitions $^{151}\text{Sm}(5/2_{g.s.}^-) \rightarrow ^{151}\text{Eu}(5/2_{g.s.}^+)$, $^{151}\text{Sm}(5/2_{g.s.}^-) \rightarrow ^{151}\text{Eu}(7/2_1^+)$, $^{171}\text{Tm}(1/2_{g.s.}^+) \rightarrow ^{171}\text{Yb}(1/2_{g.s.}^-)$, $^{171}\text{Tm}(1/2_{g.s.}^+) \rightarrow ^{171}\text{Yb}(3/2_1^-)$, $^{210}\text{Pb}(0_{g.s.}^+) \rightarrow ^{210}\text{Bi}(1_{g.s.}^-)$, and $^{210}\text{Pb}(0_{g.s.}^+) \rightarrow ^{210}\text{Bi}(0_1^-)$ have been computed using beta-decay theory with several refinements for these first-forbidden nonunique (ff-nu) β^- transitions. These ff-nu β^- transitions have non-trivial electron spectral shapes with transition nuclear matrix elements (NMEs) computed by using the microscopic Interacting Boson-Fermion Model (IBFM-2) for the decays of ^{151}Sm and ^{171}Tm , and the nuclear shell model (NSM) for the decay of ^{210}Pb . Within the respective Q windows, the computed ff-nu electron spectral shapes deviate maximally at sub-percent level from the universal allowed shape, except for the transition $^{210}\text{Pb}(0_{g.s.}^+) \rightarrow ^{210}\text{Bi}(1_{g.s.}^-)$, where the maximal deviation is some 2.7%. This confirms that the so-called ξ approximation is fairly good for most of these low- Q β^- transitions and thus the allowed shape is a rather good first approximation. Our computed spectral shapes could be of interest for experiments aiming to measure the cosmic neutrino background (CvB), like the PTOLEMY experiment. We have also derived CvB cross sections for the ground-state transitions of the considered nuclei at the β endpoint. Our findings indicate that more work on the atomic mismatch correction is needed in the future in order to extract reliable and precise CvB cross sections for any nuclear target.

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Electron spectral shapes of forbidden nonunique β decays can play a prominent role in several contexts of the present-day nuclear and particle physics, e.g., when trying to pin down the effective value g_A^{eff} of the weak axial coupling in the context of g_A -dependent spectral shapes (the Spectrum-Shape Method, SSM, introduced in [1,2] and applied in [3,4]), and when trying to explain the reactor antineutrino anomaly (RAA) [5–7] and the spectral “bump” related to the measured antineutrino flux from nuclear reactors [8–12]. The problem of the effective value of g_A can have serious consequences for the sensitivity of the running and future experiments trying to detect the neutrinoless double beta decay [13–16].

Another important context where the electron spectral shape of a forbidden nonunique β transition plays a decisive role is the detection of the cosmic neutrino background (CvB) [17–19]. The CvB

is a relic of the early Universe and plays an essential role in understanding many key features of the microwave and dark-matter cosmology [20,21]. The wide evidence from cosmological surveys supports indirectly the existence of CvB, but direct evidence on the CvB is still lacking. Detection of CvB in controlled laboratory conditions would thus provide the first proof of the existence of non-relativistic neutrinos.

In the proposed experimental methods the detection of CvB leans on relic neutrino capture on an unstable but long-lived beta emitter with a sizable neutrino-capture cross section. In addition, a small decay energy (Q value) is desirable in order to improve the detection potential of the experiment [18]. The PTOLEMY collaboration [22] considered adsorption of target isotope Tritium on graphene layers [23]. Such a design allows to achieve sufficient event rate without spoiling the performance of energy measurement of the emitted electrons. However, it was shown [24] that localization of Tritium on graphene layer leads to unacceptably large fluctuations of the energy of the emitted β electron due to the Heisenberg uncertainty principle. This problem may be over-

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come by using heavier target nuclei or weakening the bonding potential of graphene [25]. Heavy β emitters as target nuclei for the PTOLEMY experiment should have a sufficiently large $C_{\nu B}$ capture rate (e.g., small Q value) and small energy fluctuations of β electrons. In addition, it is crucial that the daughter nuclei do not experience β decay with larger Q value, creating a fatal background to the electron spectrum. The nuclei ^{151}Sm and ^{171}Tm are proposed as suitable candidates [24] and the capture cross sections are estimated in [26,27]. Another potential candidate is ^{241}Pu [28] if the α decays of the nuclide are proven to not create a background. Also the decays of the nuclei ^{210}Pb and ^{228}Ra are deemed interesting [27] but they are not suitable for PTOLEMY since their daughter isobars β decay with large Q values. However, they could be interesting for other type of experiments. All the mentioned nuclei decay via first-forbidden nonunique (ff-nu) β^- transitions to the ground state and the first excited state. These are the only final states in the β -decay Q windows due to the smallness of the Q values.

For low- Q -value ff-nu β transitions in heavy nuclei the so-called ξ approximation, where the associated electron spectral shape can be well approximated by the allowed shape, is usually valid [29,30]. In particular cases, dictated by the nuclear wave functions of the initial and final states, this approximation can be insufficient and a nuclear-structure calculation has to be done for the involved nuclear matrix elements (NMEs) [29–31]. In [27] it was found indirectly that the ξ approximation should be valid for the decays of ^{151}Sm and ^{171}Tm to their daughter nuclei ^{151}Eu and ^{171}Yb . In this work we want to verify if this indirect method really gives reliable electron spectral shapes for these transition by calculating the NMEs involved in the transitions $^{151}\text{Sm}(5/2_{g.s.}^-) \rightarrow ^{151}\text{Eu}(5/2_{g.s.}^+)$, $^{151}\text{Sm}(5/2_{g.s.}^-) \rightarrow ^{151}\text{Eu}(7/2_1^+)$, $^{171}\text{Tm}(1/2_{g.s.}^-) \rightarrow ^{171}\text{Yb}(1/2_{g.s.}^+)$, and $^{171}\text{Tm}(1/2_{g.s.}^-) \rightarrow ^{171}\text{Yb}(3/2_1^-)$ by using a nuclear-structure model called the microscopic Interacting Boson-Fermion Model (IBFM-2) [32], suitable for extracting wave functions of heavy, possibly deformed, nuclei. For the transitions $^{210}\text{Pb}(0_{g.s.}^+) \rightarrow ^{210}\text{Bi}(1_{g.s.}^-)$ and $^{210}\text{Pb}(0_{g.s.}^+) \rightarrow ^{210}\text{Bi}(0_1^-)$ the involved NMEs have been calculated by using the nuclear shell model (MSM) [33]. This is possible owing to the (near) semi-magicity of the involved nuclei.

In addition to the β spectral shapes we set out to determine the $C_{\nu B}$ scattering cross sections using the endpoint region of the computed β -electron spectra. The various corrective contributions to the cross sections are quantified and analyzed.

The beta-decay transitions discussed in this work are of the β^- type and the corresponding half-life can be cast into the form

$$t_{1/2} = \frac{\kappa}{\tilde{C}}, \quad (1)$$

where $\kappa = 6289\text{s}$ is a universal constant and \tilde{C} is the so-called integrated shape function which is given by

$$\tilde{C} = \int_1^{w_0} F_0(Z, w_e) p w_e (w_0 - w_e)^2 K(Z, w_e) C(w_e) dw_e, \quad (2)$$

where $F_0(Z, w_e)$, is the Fermi function taking into account the final-state Coulomb distortion of the wave function of the emitted electron, Z is the proton number of the final nucleus, and $w_0 = W_0/m_e$, $w_e = W_e/m_e$, and $p = p_e/m_e = \sqrt{w_e^2 - 1}$ are dimensionless kinematic variables. Here p_e and W_e are the momentum and energy of the emitted electron, respectively, and W_0 is the beta endpoint energy. The factor $K(Z, w_e)$ includes all the correction terms. Here we follow the corrections discussed in [34] taking into account finite size, finite mass, radiative corrections,

atomic exchange effects, and screening effects. The quantity of special interest in this work is the factor $C(w_e)$, known as the shape factor [29] and given by

$$C(w_e) = \sum_{k_e, k_\nu, K} \lambda_{k_e} \left[M_K(k_e, k_\nu)^2 + m_K(k_e, k_\nu)^2 - \frac{2\gamma_{k_e} \mu_{k_e}}{k_e w_e} M_K(k_e, k_\nu) m_K(k_e, k_\nu) \right], \quad (3)$$

where k_e and k_ν come from the partial-wave expansion of the lepton wave functions, $\gamma_{k_e} = \sqrt{k_e^2 - (\alpha Z)^2}$, $\mu_{k_e} \approx 1$, and $\lambda_{k_e} = F_{k_e-1}(Z, w_e)/F_0(Z, w_e)$ is the Coulomb function with $F_{k_e-1}(Z, w_e)$ being the generalized Fermi function. The quantities $M_K(k_e, k_\nu)$ and $m_K(k_e, k_\nu)$ have lengthy expressions which can be found from [29], and include both universal kinematic factors as well as model-dependent form factors. The form factors are either of the vector ${}^V F_{KLS}^{(N)}$ or axial-vector ${}^A F_{KLS}^{(N)}$ type, with K, L, s corresponding to the angular momenta of the operators and $N = 0, 1, 2, \dots$ emerge from a power expansion discussed in length in [29].

In order to evaluate these form factors, a common approach is to use the so-called impulse approximation, where the neutron turning into a proton is assumed not to interact with the other nucleons when the decay takes place. In this approximation the form factors can be related to nuclear matrix elements ${}^{V/A} \mathcal{M}_{KLS}^{(N)}$ by

$${}^V F_{KLS}^{(N)} = (-1)^{K-L} R^{-L} g_V {}^V \mathcal{M}_{KLS}^{(N)} \quad (4)$$

$${}^A F_{KLS}^{(N)} = (-1)^{K-L+1} R^{-L} g_A {}^A \mathcal{M}_{KLS}^{(N)}, \quad (5)$$

where R is the nuclear radius, K is the total and L is the orbital angular momentum of the operator, while s is the spin. ${}^V \mathcal{M}_{KLS}^{(N)}$ is a vector-type matrix element and is always multiplied by the vector coupling constant g_V , while ${}^A \mathcal{M}_{KLS}^{(N)}$ is an axial-vector-type matrix element and is always multiplied by the axial-vector coupling g_A . The couplings originate in the formalism when moving from the quark level to the nucleon level as a way to renormalize the hadronic current. The sign convention is chosen as in, e.g., [2]. The nuclear matrix elements can be further broken down to single-particle matrix elements

$${}^{V/A} \mathcal{M}_{KLS}^{(N)} = \frac{\sqrt{4\pi}}{\hat{J}_i} \sum_{pn} {}^{V/A} m_{KLS}^{(N)}(pn) (\Psi_f || [c_p^\dagger \tilde{c}_n]_K || \Psi_i), \quad (6)$$

where $\hat{J}_i = \sqrt{2J_i + 1}$ with J_i being the spin of the initial state in the nucleus, ${}^{V/A} m_{KLS}^{(N)}(pn)$ is the single-particle matrix element corresponding to the proton orbital p and neutron orbital n , and $(\Psi_f || [c_p^\dagger \tilde{c}_n]_K || \Psi_i)$ is the one-body transition density (OBTD), which contains the relevant nuclear-structure information. The choice of nuclear model enters the calculation through the evaluation of the OBTDs.

For the first-forbidden non-unique decays considered in this work, the relevant NMEs are those of the transition operators corresponding to spin-parity changes $0^-, 1^-,$ and 2^- . In the expansion of Behrens and Bühring [29] there are six matrix elements corresponding to the operators

$$\mathcal{O}(0^-) : g_A \in \text{MEC}(\boldsymbol{\sigma} \cdot \mathbf{p}_e), \quad g_A(\boldsymbol{\sigma} \cdot \mathbf{r}) \quad (7)$$

$$\mathcal{O}(1^-) : g_V \mathbf{p}_e, \quad g_A(\boldsymbol{\sigma} \times \mathbf{r}), \quad g_V \mathbf{r} \quad (8)$$

$$\mathcal{O}(2^-) : g_A[\boldsymbol{\sigma} \mathbf{r}]_2, \quad (9)$$

where \mathbf{r} is the radial coordinate and \mathbf{p}_e is the electron momentum. For a given decay the relevant operators are those with a rank between $|J_i - J_f|$ and $J_i + J_f$. For example, for the transition $^{210}\text{Pb}(0_{g.s.}^+) \rightarrow ^{210}\text{Bi}(0_1^-)$ only the two $\mathcal{O}(0^-)$ operators contribute.

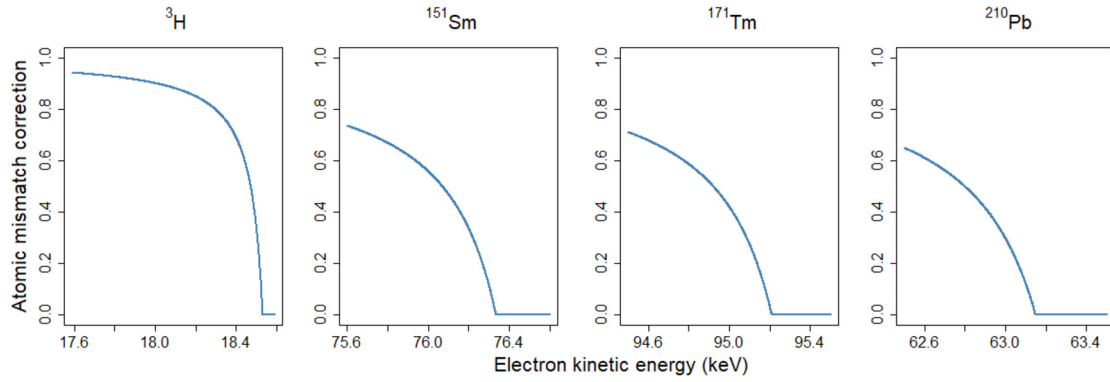


Fig. 1. The mismatch correction of Eq. (11) for ${}^3\text{H}$ and the other three $\text{C}\nu\text{B}$ candidates studied in this work.

In this work the well-known fundamental enhancement (see, e.g., [35]) of the axial-charge NME ($\sigma \cdot \mathbf{p}_e$) is denoted by ϵ_{MEC} .

The Behrens-Bühring formalism is based on expanding the matrix elements in the small quantities $W_e R$, $m_e R$, and $Z\alpha$ which allows for some power-series considerations. In the so-called ξ approximation the complicated shape factor (3) is expressed in powers of $1/\xi$, where $\xi = Z\alpha/((W_e - m_e)R)$. It can then be shown that the shape factor of a forbidden transition is that of an allowed one with corrections of the order $Z\alpha/Q R$, where Q is the decay energy (Q value) of the transition. When $1/\xi$ is small, allowed approximation of the spectrum shape is reasonably accurate. For the transitions considered here $\xi \approx 150$ for the ground-state-to-ground-state decay of ${}^{171}\text{Tm}$ and even larger for the other transitions. Therefore, one would expect at most corrections of the order of 0.67% to the spectra. However, as pointed out in e.g. [27], there can be limitations to the applicability of the ξ approach, in particular if notable cancellations appear among the various terms of the shape factor (3) [31].

In order to see whether such cancellations appear, a proper microscopic calculation of the shape factor must be performed. As the spectrum shape depends on the nuclear matrix elements only through the ratio g_A/g_V (see [2]) the way uncertainties in the nuclear matrix elements propagate to the spectrum shape can be estimated by fixing g_V to its free-nucleon value of unity and varying g_A . Based on the papers [1,2,40,41] reasonable error estimates can be obtained by using an effective value for the axial-vector coupling constant $g_A = g_A^{\text{eff}}$, and varying it between 0.80 and 1.20. However, since this method alone cannot address uncertainties in the ratios of all matrix elements (e.g., the two rank-1 vector matrix elements) and thus we also varied the matrix elements independently by $\pm 20\%$ to get robust error estimates. Based on the studies [35–39] we vary the value of the mesonic enhancement factor ϵ_{MEC} of the axial-charge matrix element between 1.4 and 2.0.

The neutrino capture rate at the beta endpoint can be derived from equations (1) and (2) and can be expressed as [19]

$$\bar{\sigma} = F_0(Z, w_e) p w_e K(Z, w_e) C(w_e) \Big|_{w_e=w_0}. \quad (10)$$

Relevant corrections here are the finite size, atomic screening, radiative, and atomic exchange corrections for which we adopt the expressions given in the comprehensive review [34] on allowed beta spectrum shape. The correction terms were evaluated at $W_e = Q - 0.01$ eV, up to which point they were stable in value. Here Q denotes the Q -value of the transition, i.e., the amount of energy released in the transition. The full correction term is relatively stable up to this point. For forbidden transitions the shape factor is the most important correction. In addition to these corrections, there is one more non-trivial correction related to the shake-up and shake-off effects, where the final atom is either left in an excited state or ionized. For the heavy nuclei considered

here, the shake-off probability is the dominant one, occurring for roughly 20–30% of the decays [34,42]. These effects can be reasonably accounted for with the so-called atomic mismatch correction, which is of the form [42,43]

$$r(Z, W_e) = 1 - \frac{1}{W_0 - W_e} (44.2 Z^{0.41} + 2.3196 \times 10^{-7} Z^{4.45}) \text{eV} + \text{small correction}. \quad (11)$$

The correction (11) is discussed in [34] in the context of spectral shapes, where its wild behavior near the endpoint is greatly mitigated by the kinematic term $(W_e - W_0)^2$ in Eq. (2) and is thus not problematic for the shape function and its integrated form (2). However, when considering the cross section (10), evaluated at the beta endpoint, the correction is problematic, as it is quite small before rapidly going to zero very close to the endpoint. In [27] this difficulty was dealt with by fixing this correction to its value at $0.98Q$ for $W_e > 0.98Q$, but this procedure is arbitrary and leaves room for improvements. In order to see the problem with this correction term, its behavior is presented for ${}^3\text{H}$, a considered candidate for the $\text{C}\nu\text{B}$ detection [19], and the three candidates studied in this work in Fig. 1 for the energy range $W_e = [Q - 1 \text{ keV}, Q]$. The Q value is now shifted downwards by the mean atomic excitation energy. Also, the cross section near the new lower endpoint goes rapidly to zero. Since in reality the shake-up and shake-off processes have some finite probability (smaller than one) of occurring, the Q -value shift would only happen for some fraction of the transitions. Thus, it is clear that for the $\text{C}\nu\text{B}$ detection this effect needs to be considered in a more sophisticated way. Since details of the atomic corrections are out of the scope of the present paper we leave this correction out from the present calculations and concentrate on nuclear structure of the shape factor and the other well-behaved corrections.

The shape factors related to the decay transitions in ${}^{151}\text{Sm}$ and ${}^{171}\text{Tm}$ were treated in the nuclear-structure framework of the microscopic Interacting Boson-Fermion Model (IBFM-2) [32]. IBFM-2 is an extension of the well-known microscopic Interacting Boson Model (IBM-2) [44,45] to odd-mass nuclear systems. The IBM-2 is a phenomenological approach that has been one of the most successful models in reproducing collective features of the low-lying levels of medium-heavy as well as heavy nuclei. The IBM-2 deals with even-even nuclei, where one replaces valence-nucleon pairs with bosons with angular momentum 0 or 2. By coupling an extra fermion to this bosonic system, one is able to extend the IBM-2 to the study of odd- A nuclei. This extension is the IBFM-2.

The mapping of the single-fermion creation operator onto the IBFM-2 space follows the procedure introduced in Refs. [46,47] where relevant terms, using exact values for the fermion matrix elements in the Generalized Seniority scheme, were worked out and thus use of the Number Operator Approximation (NOA) was

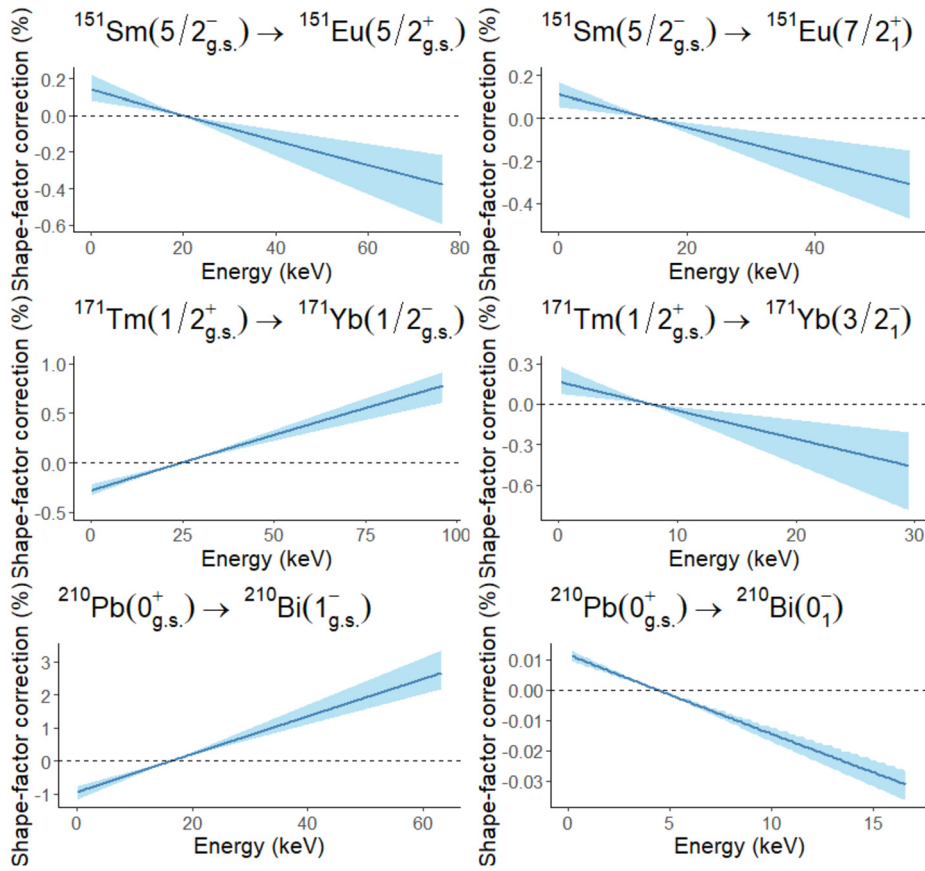


Fig. 2. Shape-factor corrections to the allowed shape of the normalized electron spectrum and the related uncertainties for each transition discussed in this work. The uncertainties have been obtained by varying the effective axial-vector coupling constant g_A^{eff} between 0.80 and 1.20 and the mesonic enhancement ϵ_{MEC} of the axial-charge matrix element between 1.4 and 2.0. In addition, each involved nuclear matrix element has been varied by $\pm 20\%$ in order to capture the possible uncertainties in the nuclear-structure calculations.

Table 1
Boson-fermion interaction parameters (in MeV) used in the IBFM-2 calculations.

	Γ_ρ	Δ_ρ	A_ρ
^{151}Sm	-0.361	0.265	-0.150
^{151}Eu	-0.080	0.090	-0.050
^{171}Tm	-0.050	-0.027	0.580
^{171}Yb	-0.062	-0.021	0.200

avoided. This method has already been applied to allowed beta decays in Ref. [48] and now we extend its use to ff-nu beta decays in the ^{151}Sm and ^{171}Tm nuclei. Since these nuclei are mid-shell nuclei they are best described by IBFM-2. Contrary to this, the nuclei ^{210}Pb and ^{210}Bi , for which the IBFM-2 model is not applicable since they are even- A nuclei, are in the vicinity of the doubly-magic nucleus ^{208}Pb , and thus can be best described using the nuclear shell model.

In the IBFM-2 calculations the even-even ^{150}Sm nucleus was used as a common core for the odd ^{151}Sm and ^{151}Eu nuclei, and ^{170}Yb and ^{172}Yb were adopted as cores for the ^{171}Yb and ^{171}Tm nuclei, respectively. The parameters for the core Sm and Yb nuclei were taken from Refs. [49,50], respectively. The valence space was chosen to span $2p, 1f, 0h_{9/2}, 0i_{13/2}$ neutron and $2s, 1d, 0g, 0h_{11/2}$ proton orbitals with unperturbed single-particle energies taken from [51], where the effect of single-particle energies on occupation probabilities was studied. The used boson-fermion interaction parameters are listed in Table 1.

The wave functions and one-body densities for the decay of ^{210}Pb were calculated in the shell-model framework using the computer program NuShellX@MSU [52]. The calculations

Table 2
Beta-spectrum endpoint corrections and their uncertainties. The uncertainties have been obtained by varying the effective axial-vector coupling constant g_A^{eff} between 0.80 and 1.20 and the mesonic enhancement ϵ_{MEC} of the axial-charge matrix element between 1.4 and 2.0, and by varying the individual nuclear matrix elements by $\pm 20\%$.

Nucleus	J_f^π	Endpoint correction (%)
^{151}Sm	$5/2^+$	$-0.38^{+0.16}_{-0.22}$
	$7/2^+$	$-0.31^{+0.15}_{-0.17}$
^{171}Tm	$1/2^-$	$0.77^{+0.14}_{-0.17}$
	$3/2^-$	$-0.46^{+0.25}_{-0.32}$
^{210}Pb	0^-	$-0.031^{+0.004}_{-0.006}$
	1^-	$2.66^{+0.67}_{-0.50}$

were done in the full model space spanning the proton orbitals $0h_{9/2}, 2p, 1f, 0i_{13/2}$ and the neutron orbitals $0i_{11/2}, 1g, 2d, 3s, 0j_{15/2}$ using the effective Hamiltonian $khpe$ [53].

The calculated shape factors for the decays of ^{151}Sm , ^{171}Tm , and ^{210}Pb are compared with the allowed approximation in Fig. 2. For all six transitions the corrections to the allowed spectral shape are largest at the endpoint, which is the region of interest for detecting the cosmic neutrino background. The endpoint corrections are given in Table 2. While the pure pseudoscalar transition $^{210}\text{Pb}(0_{g.s.}^+) \rightarrow ^{210}\text{Bi}(0_1^-)$ gets only a tiny correction of $-0.031^{+0.004}_{-0.006}\%$ the transition $^{210}\text{Pb}(0_{g.s.}^+) \rightarrow ^{210}\text{Bi}(1_{g.s.}^-)$ gets a non-trivial $2.66^{+0.67}_{-0.50}\%$ correction. The shape-factor corrections for the four other transitions are roughly 0.50%. Looking at the dominating ground-state-to-ground-state transitions the corrections are $0.7/\xi$ for ^{151}Sm , $1.2/\xi$ for ^{171}Tm , and $7.1/\xi$ for ^{210}Pb . In addition, for

Table 3

Endpoint cross sections and their leading errors for the ground-state transitions of the discussed mother nuclei. All relevant corrections except the atomic mismatch effect are included. The half-life and Q -value data have been taken from [55].

Nucleus	$\bar{\sigma}$ (cm ²)	Half-life error	Q -value error	Spectrum-shape error	Total error
¹⁵¹ Sm	4.79×10^{-48}	0.44×10^{-48}	0.01×10^{-48}	0.01×10^{-48}	0.44×10^{-48}
¹⁷¹ Tm	1.14×10^{-46}	0.01×10^{-46}	0.07×10^{-46}	0.01×10^{-46}	0.07×10^{-46}
²¹⁰ Pb	3.27×10^{-48}	0.04×10^{-48}	0.02×10^{-48}	0.01×10^{-48}	0.05×10^{-48}

the three decays to the excited states the corrections are roughly $1/\xi$. While these results are consistent with the $\mathcal{O}(1/\xi)$, the results highlight the fact that one should not assume that the error is necessarily exactly $\approx 1/\xi$. The decay of ²¹⁰Pb to the ground state is a good example of destructive interference, where the ξ approximation does not hold very well.

The cross sections and the impacts of the relevant corrections for the ground-state-to-ground-state transitions of ¹⁵¹Sm, ¹⁷¹Tm, and ²¹⁰Pb are given in Table 3. For ¹⁵¹Sm and ¹⁷¹Tm the obtained values $(4.79 \pm 0.44) \times 10^{-48}$ cm² and $(1.14 \pm 0.07) \times 10^{-46}$ cm² are somewhat larger than the values $(4.77 \pm 0.01) \times 10^{-48}$ cm² and $(1.12 \pm 0.01) \times 10^{-46}$ cm² reported in [27]. The small deviations in the estimates are due to some differences in how the correction terms are evaluated [54] and the larger uncertainties in the present work are due to the inclusion of all the relevant sources of uncertainty. From Table 3 it is clear that the error in the ξ approximation is overwhelmed by the half-life error for ¹⁵¹Sm, by the Q -value error for ¹⁷¹Tm, and by a combination of both for ²¹⁰Pb. But, nevertheless, our computations of the ff-nu spectral shapes quantitatively verify that there are no large accidental cancellations of their various terms thus preserving a sufficient accuracy of the ξ approximation.

The errors in the computed cross sections affect directly the needed target mass in experiments. In [27] it was estimated that in an ideal case, for an experiment running for one year, the needed target mass would be some 6 tonnes for ¹⁵¹Sm and 350 kg for ¹⁷¹Tm (for caveats in the case of ¹⁷¹Tm, see [27]). The presently obtained errors in the cross sections then indicate a variation in the needed target mass of the order of 10%, much more than the sub-1% level predicted in [27].

In this Letter we have calculated the endpoint cross sections for ¹⁵¹Sm, ¹⁷¹Tm, and ²¹⁰Pb using realistic microscopic nuclear models for the involved wave functions. These low- Q -value decays are potential candidates for the detection of cosmic neutrino background. Out of these candidates the most promising one is ¹⁷¹Tm with $\bar{\sigma} = (1.14 \pm 0.07) \times 10^{-46}$ cm². The validity of the ξ approximation for forbidden spectral shapes was investigated and errors up to $7.1/\xi$ were recorded. While this can be considered consistent with an $\mathcal{O}(1/\xi)$ estimate, our results highlight the fact that one should not assume a-priori an error $\approx 1/\xi$, but nuclear structure can alter the situation depending on the details of the initial and final nuclear wave functions.

Declaration of competing interest

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

Data availability

No data was used for the research described in the article.

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