

α decay in the complex-energy shell modelR. Id Betan^{1,2,3} and W. Nazarewicz^{1,2,4}¹*Department of Physics and Astronomy, University of Tennessee, Knoxville, Tennessee 37996, USA*²*Physics Division, Oak Ridge National Laboratory, P.O. Box 2008, Oak Ridge, Tennessee 37831, USA*³*Departamento de Química y Física, FCEIA(UNR), Instituto de Física Rosario (CONICET), Av. Pellegrini 250, 2000 Rosario, Argentina*⁴*Institute of Theoretical Physics, University of Warsaw, ul. Hoża 69, 00-681 Warsaw, Poland*

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Background: Alpha emission from a nucleus is a fundamental decay process in which the alpha particle formed inside the nucleus tunnels out through the potential barrier.**Purpose:** We describe alpha decay of ²¹²Po and ¹⁰⁴Te by means of the configuration interaction approach.**Method:** To compute the preformation factor and penetrability, we use the complex-energy shell model with a separable $T = 1$ interaction. The single-particle space is expanded in a Woods-Saxon basis that consists of bound and unbound resonant states. Special attention is paid to the treatment of the norm kernel appearing in the definition of the formation amplitude that guarantees the normalization of the channel function.**Results:** Without explicitly considering the alpha-cluster component in the wave function of the parent nucleus, we reproduce the experimental alpha-decay width of ²¹²Po and predict an upper limit of $T_{1/2} = 5.5 \times 10^{-7}$ sec for the half-life of ¹⁰⁴Te.**Conclusions:** The complex-energy shell model in a large valence configuration space is capable of providing a microscopic description of the alpha decay of heavy nuclei having two valence protons and two valence neutrons outside the doubly magic core. The inclusion of proton-neutron interaction between the valence nucleons is likely to shorten the predicted half-life of ¹⁰⁴Te.DOI: [10.1103/PhysRevC.86.034338](https://doi.org/10.1103/PhysRevC.86.034338)

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I. INTRODUCTION

According to Gamow theory of alpha decay [1,2], this fundamental radioactive decay can be considered as a two-step process [3–5]. In the first step, an alpha cluster is formed inside the parent nucleus. The resulting alpha particle resides in a metastable state of an average potential of the daughter system. In the second step, the particle tunnels through the potential barrier. Each step requires different theoretical treatment. To compute the preformation factor that describes the alpha formation probability, one needs to evaluate the overlap integral involving wave functions of the parent and daughter nuclei, and that of the alpha particle. The estimate of the penetration probability requires a careful treatment of the resonance state.

The commonly used formulation of the alpha-decay problem employs the R -matrix expression [6,7]

$$\Gamma_L = 2P_L\gamma_L^2 \quad (1)$$

for the absolute width. In this formalism, the first stage (formation of alpha particle with angular momentum L) is given by the reduced width γ_L^2 , while the second stage (decay) is expressed by means of the penetrability P_L . Alternatively, the absolute width can be obtained from the general reaction-theory expression [8–11]

$$\Gamma_L = S_L\Gamma_L^{\text{sp}}, \quad (2)$$

where S_L is the alpha-spectroscopic factor and Γ_L^{sp} is the single-particle (s.p.) decay width.

Historically, expression (1) was derived in 1954 by Thomas [6] using the time-independent R -matrix theory of nuclear reactions. In 1957, Mang [7] developed the alpha-decay

formalism based on the time-dependent perturbation theory. He made the connection with the shell model and succeeded in expressing the alpha-decay formation amplitude in a basis of s.p. states. As shown in Refs. [12,13], formulations of Thomas and Mang are formally equivalent; there are, however, many differences when it comes to practical implementations.

The reduced width calculated in a shell-model configuration expressed in the harmonic oscillator (h.o.) basis is too small. This can be partly cured by means of configuration mixing involving extended shell-model spaces [14,15] as each admixed configuration contributes coherently to γ_L^2 . To improve asymptotic properties of s.p. wave functions, the particle continuum was taken into account [16] by considering h.o. expansion [17] or within a Woods-Saxon (WS) basis consisting of bound and outgoing single-particle resonant (Gamow) states [18,19]. The configuration mixing calculations of Refs. [15,18] in the valence space of ²¹²Po assumed the seniority-zero (pairing vibrational) wave functions obtained by considering the monopole pairing interaction between like nucleons. However, all these improvements were not sufficient to reproduce the experimental alpha decay in ²¹²Po. It was only after the valence proton-neutron interaction had been considered together with a generalized wave function expressed as a combination of cluster and shell model components [20] that theoretical and experimental widths could be reconciled [21].

The R -matrix expression for the width (1) depends on the channel radius R . This radius should be chosen large enough so that the alpha-daughter interaction in the external region is given by the Coulomb force alone [22]. The infinite range of the Coulomb force implies, however, that the asymptotic behavior of the R -matrix expression is reached only at large values of R ,

at which the asymptotic behavior of the shell-model s.p. basis (h.o. basis in most applications) used to calculate $\gamma_L^2(R)$ does matter. Due to the mismatch between the internal part of the s.p. wave function (well described in the h.o. basis) and the asymptotic part (poorly or not described in the h.o. basis), rather small changes in R may produce appreciable variations in penetrability. Physically, the reason for this sensitivity is the fact that the alpha cluster is formed in the surface region of the nucleus in which the coupling to the alpha continuum that impacts the radial behavior of the formation amplitude is important [16]. Consequently, the absolute R -matrix width depends in general on the channel radius [5,23], and this is an obvious drawback of the method [22].

Our renewed interest in the alpha-decay problem is stimulated by the recent experimental data above the doubly-magic ^{100}Sn [24,25] that demonstrate the presence of very fast alpha decays. Indeed, the observed enhancement of the reduced widths of $^{105,106}\text{Te}$ relative to $^{213,212}\text{Po}$ is 2–3, thus confirming earlier expectations [26] of “superallowed” alpha decays in this region due to the large overlaps of valence s.p. shell-model proton and neutron wave functions. Our long-term goal is to estimate alpha preformation factors in nuclei above ^{208}Pb and ^{100}Sn by using large valence s.p. spaces, including positive-energy Gamow states of a finite-depth WS potential [18,19]. In this study, we focus on ^{212}Po and ^{104}Te nuclei having two valence protons and two valence neutrons outside doubly-magic cores.

Our paper is organized as follows. Section II briefly describes the alpha-decay formalism used in this work, with special emphasis on approximations used to describe wave functions of parent and daughter nuclei. Section III deals with the approximations employed and parameters used. In particular, we discuss the sensitivity of the calculated spectroscopic factor to the parameters defining the shifted Gaussian basis that is used to compute the normalization of the channel function. In Sec. IV we study the sensitivity of the reduced alpha width in ^{212}Po on the choice of s.p. basis used. In Sec. V we discuss the absolute alpha-decay width of ^{212}Po and in Sec. VI we compare it with the absolute width of the superallowed alpha emitter ^{104}Te . Finally, the main conclusions of this work are summarized in Sec. VII.

II. FORMALISM

In this section, we discuss the R -matrix (1) and spectroscopic factor (2) expressions for the decay width. The connection between the two formulations is given in Ref. [8]. We also discuss the so-called δ -approximation for the formation amplitude.

A. R -matrix expression for the decay width

Within the R -matrix theory [6,7,12], the absolute width is given by Eq. (1) with $P_L(R)$ being the barrier penetrability and $\gamma_L(R)$ the reduced width amplitude [27]. While both quantities strongly depend on the value of the channel radius R , the absolute width should be R independent.

For $P_L(R)$ we use the standard expression [6]

$$P_L(R) = \frac{kR}{|H_L^+(\eta, kR)|^2}, \quad (3)$$

where k is given by the alpha energy $E_\alpha = \frac{\hbar^2 k^2}{2\mu}$, obtained from the experimental Q_α value by correcting for electron screening; $\mu = \frac{m_d m_\alpha}{m_d + m_\alpha}$ is the reduced mass of alpha particle with m_d being the mass of the daughter nucleus; $H_L^+(\eta, kR)$ is the outgoing spherical Coulomb-Hankel function; and $\eta = \frac{2Z_d \mu e^2}{\hbar^2 k}$ is the Sommerfeld Coulomb parameter.

The reduced width amplitude $\gamma_L(R)$ may be written in terms of the formation amplitude $g_L(R)$ [5,14]:

$$\gamma_L = \sqrt{\frac{\hbar^2 R}{2\mu}} g_L(R), \quad (4)$$

with

$$g_L(R) = \int d\Omega_R \int d\xi_\alpha \int d\xi_D \times \Phi_{JM}^P \mathcal{A}[\phi_\alpha(\xi_\alpha) \Psi_j^D(\xi_D) Y_L(\hat{R})]_{JM}^*, \quad (5)$$

where ϕ_α is the normalized wave function of the alpha particle with zero angular momentum, Y_{LM} is the angular part of the center-of-mass (c.o.m.) motion of the alpha particle, $\Psi_{j m_j}^D$ is the wave function of the daughter nucleus, and Φ_{JM}^P is the wave function of the parent nucleus. The coordinates ξ_α and ξ_D are the intrinsic coordinates of the alpha particle and daughter nucleus, respectively. All wave functions are normalized in terms of the internal and c.o.m. coordinates [27]. By construction, the parent and daughter wave functions are antisymmetric. The antisymmetrization with respect to interfragment nucleons is done by means of the operator \mathcal{A} . Its action can be approximated by means of a factor $[(\binom{N_v}{2})(\binom{Z_v}{2})]^{1/2}$ [14,27,28], with N_v and Z_v being, respectively, the numbers of valence neutrons and protons in the parent nucleus.

For the internal alpha-particle wave function we take the standard Gaussian ansatz [18,29]:

$$\begin{aligned} \phi_\alpha(\rho_1 \rho_2 \rho_3, \sigma_1 \sigma_2 \sigma_3 \sigma_4) &= \phi(\rho_1 \rho_2 \rho_3) \chi_{00}(\sigma_1 \sigma_2) \chi_{00}(\sigma_3 \sigma_4), \\ \chi_{00}(\sigma_1 \sigma_2) &= [\chi_{1/2}(\sigma_1) \chi_{1/2}(\sigma_2)]_{00}, \\ \phi(\rho_1 \rho_2 \rho_3) &= \left(\frac{8\beta}{\pi}\right)^{9/4} e^{-4\beta(\rho_1^2 + \rho_2^2 + \rho_3^2)}. \end{aligned} \quad (6)$$

The parameter $\beta = \frac{9}{64r_\alpha^2} = 0.057 \text{ fm}^{-2}$ depends on the root-mean-square alpha radius $r_\alpha = 1.57 \text{ fm}$ [29].

The transformation between the intrinsic $\xi_\alpha = \{\boldsymbol{\rho}_1, \boldsymbol{\rho}_2, \boldsymbol{\rho}_3\}$ and nucleonic $\{\mathbf{r}_i\}$ ($i = 1, 2, 3, 4$) coordinates reads

$$\begin{aligned} \boldsymbol{\rho}_1 &= \frac{\mathbf{r}_1 - \mathbf{r}_2}{\sqrt{2}}, & \boldsymbol{\rho}_2 &= \frac{\mathbf{r}_3 - \mathbf{r}_4}{\sqrt{2}}, \\ \boldsymbol{\rho}_3 &= \frac{(\mathbf{r}_1 + \mathbf{r}_2) - (\mathbf{r}_3 + \mathbf{r}_4)}{2}, \end{aligned} \quad (7)$$

and

$$\mathbf{R} = \frac{\mathbf{r}_1 + \mathbf{r}_2 + \mathbf{r}_3 + \mathbf{r}_4}{4} \quad (8)$$

is the c.o.m. coordinate of alpha particle. Let us denote the spherical components of intrinsic coordinates by

$\rho_i = (\rho_i, \tilde{\theta}_i, \tilde{\varphi}_i)$. Assuming $\theta_R = \varphi_R = 0$, the nucleonic coordinates can be written as

$$\begin{aligned} 4r_{1,2}^2 &= 4R^2 + \rho_3^2 + 2\rho_1^2 \pm 2\sqrt{2}\rho_3\rho_1 \cos \tilde{\theta}_{31} \\ &\quad + 4R(\rho_3 \cos \tilde{\theta}_3 \pm \sqrt{2}\rho_1 \cos \tilde{\theta}_1), \\ 4r_{3,4}^2 &= 4R^2 + \rho_3^2 + 2\rho_2^2 \mp 2\sqrt{2}\rho_3\rho_2 \cos \tilde{\theta}_{32} \\ &\quad - 4R(\rho_3 \cos \tilde{\theta}_3 \pm \sqrt{2}\rho_2 \cos \tilde{\theta}_2), \end{aligned} \quad (9)$$

where $\tilde{\theta}_{ij} = \tilde{\theta}_j - \tilde{\theta}_i$, and

$$\begin{aligned} \cos \theta_{1,2} &= \frac{2R + \rho_3 \cos \tilde{\theta}_3 \pm \sqrt{2}\rho_1 \cos \tilde{\theta}_1}{2r_{1,2}}, \\ \cos \theta_{3,4} &= \frac{2R - \rho_3 \cos \tilde{\theta}_3 \pm \sqrt{2}\rho_2 \cos \tilde{\theta}_2}{2r_{3,4}}. \end{aligned} \quad (10)$$

This paper deals with g.s. \rightarrow g.s. alpha decays to the magic daughter nucleus. Assuming the seniority-zero wave function, the corresponding formation amplitude is [13,14]

$$F_0(R) = \frac{\sqrt{8}}{16\pi^{3/2}} \sum_{v_n, v_p} (-)^{l_n+l_p} b_{v_n, v_p} \hat{J}_n \hat{J}_p I_{v_n, v_p}(R), \quad (11)$$

where

$$\begin{aligned} I_{v_n, v_p}(R) &= \int d\rho_1 d\rho_2 d\rho_3 \phi(\rho_1 \rho_2 \rho_3) \\ &\quad \times \frac{u_{v_n}(r_1) u_{v_n}(r_2)}{r_1 r_2} P_{l_n}(\cos \theta_{12}) \\ &\quad \times \frac{u_{v_p}(r_3) u_{v_p}(r_4)}{r_3 r_4} P_{l_p}(\cos \theta_{34}), \end{aligned} \quad (12)$$

with $\theta_{ij} = \theta_j - \theta_i$, $v = \{n, l, j\}$, and $u_v(r)$ being s.p. radial wave functions. The factor $\sqrt{8}$ comes from the Jacobian of the transformation between the nucleonic coordinates $\{\mathbf{r}_i\}$ and the internal and c.o.m. coordinates [5,30]. In Eq. (12) and in the following, the s.p. indices 1, 2 refer to neutrons while 3, 4 refer to protons. The coefficients b_{v_n, v_p} are the shell-model four-particle wave function amplitudes.

B. δ -function approximation

In the calculation of alpha-decay rates based on h.o. wave functions, it was noticed [28] that the relative rates change little with the oscillator length $b_{\text{h.o.}}$ of the basis. Using this argument, Mang proposed to take $\beta \gg 1/b_{\text{h.o.}}^2$. In this limit, the expression for the formation amplitude can be simplified (see also Ref. [31]). In the literature, this is known as δ -function approximation [32].

In practice, one assumes that the alpha-particle wave function is constant inside a small volume of radius $s_\alpha = 2.34$ fm [32] and zero outside. Within this approximation $\rho_i = 0$; hence, it immediately follows from Eqs. (9) that $\mathbf{r}_1 = \mathbf{r}_2 = \mathbf{r}_3 = \mathbf{r}_4 = \mathbf{R}$ [32,33], and the formation amplitude reduces to

$$F_0^\delta(R) = \frac{\sqrt{8}}{16\pi^{3/2}} \left(\frac{4\pi s_\alpha^3}{3} \right)^{3/2} \left(\sum_{v_n} I_{v_n}^n \right) \left(\sum_{v_p} I_{v_p}^p \right), \quad (13)$$

with

$$I_v^\tau = (-)^{l_v} b_v^\tau \hat{J}_v B_v \frac{u_{v\tau}^2(R)}{R^2}, \quad (14)$$

where $\tau = n, p$. The correction factor B_v depends on the relative angular momentum [32]:

$$B_v = 1 - 0.013l_v(l_v + 1). \quad (15)$$

C. Four-particle amplitudes

For the g.s. alpha decay of ^{212}Po and ^{104}Te , we are going to assume that the four valence nucleons move around the rigid, doubly-magic core. The parent-nucleus wave function is approximated by a product of two-neutron and two-proton seniority-zero states:

$$|\Phi_{J=0, M=0}^P\rangle = |\Psi_{2n,00}\rangle \otimes |\Psi_{2p,00}\rangle, \quad (16)$$

where

$$|\Psi_{2\tau,0}\rangle = \sum_v X_v^\tau |v\nu, 00\rangle, \quad (17)$$

$|v\nu, 00\rangle = \frac{[a_v^\dagger a_v^\dagger]_{00}}{\sqrt{2}} |0_\tau\rangle$, and $|0\rangle = |0_n\rangle \otimes |0_p\rangle$ is the shell-model vacuum representing the ^{208}Pb or ^{100}Sn g.s. wave function. The four-particle amplitudes b_{v_n, v_p} in Eq. (11) can thus be written in a separable form:

$$b_{v_n, v_p} = X_{v_n}^n X_{v_p}^p. \quad (18)$$

D. Alpha-decay spectroscopic factor

Based on the general theoretical arguments [8–11], the absolute width can be expressed as a product of the alpha-particle spectroscopic factor and the single-particle width; see Eq. (2). The spectroscopic factor S_L contains information about the probability of forming an alpha cluster in the parent system. Since the alpha particle, when formed, occupies the resonant state, the s.p. width can be obtained from the so-called current expression [5,34,35]:

$$\Gamma_L^{\text{sp}} = i \frac{\hbar^2}{2\mu} \frac{u_L^*(R) u_L(R) - u_L'(R) u_L'^*(R)}{\int |u_L(R)|^2 dR}, \quad (19)$$

where the Gamow function $u_L(R)$ is obtained as a solution of the Schrödinger equation with outgoing boundary condition. When the imaginary part of the complex energy eigenvalue $\mathcal{E}_\alpha = \frac{\hbar^2 k^2}{2\mu}$ is small, which is always the case for the considered g.s. alpha emitters, one can approximate (19) with [36]

$$\Gamma_L^{\text{sp}} = \frac{\hbar^2 \Im(k)}{\mu} \frac{|u_L(R)|^2}{|H_L^+(\eta, kR)|^2}. \quad (20)$$

The s.p. width obtained in this way should be identical to the value $-2\text{Im}(\mathcal{E}_\alpha)$ given by the imaginary part of the Gamow resonance energy, if the latter is computed with a sufficient precision.

The conventional alpha spectroscopic factor as introduced in Ref. [8] is defined by

$$S_L = |\langle \mathcal{A}[\phi_\alpha(\xi_\alpha) \Psi_j^D(\xi_D) \psi_L(\mathbf{R})]_{JM} | \Phi_{JM}^P \rangle|^2, \quad (21)$$

where $\psi_{LM}(\mathbf{R}) = \frac{u_L(R)}{R} Y_{LM}(\hat{R})$ represents the relative motion alpha particle with respect to the daughter. In terms of the formation amplitude, S_L reads [3,5,37]

$$S_L = \int_0^\infty g_L^2(R) R^2 dR. \quad (22)$$

E. Modified spectroscopic factor

Since the formation amplitude Eq. (5) represents the overlap of the parent wave function with the daughter-alpha product state, one would be tempted to associate it with the probability amplitude that in the parent wave function Φ_{JM}^P an alpha particle ϕ_α and a daughter nucleus $\Psi_{jm_j}^D$ are at a distance R . The value of S_L would then be associated with the total probability of formation of an alpha particle. However, the fundamental problem with this interpretation is that the channel function $\mathcal{A}[\phi_\alpha(\xi_\alpha) \Psi_{jm_j}^D(\xi_D) \psi_L(\mathbf{R})]_{JM}$ is not properly normalized [3,10,11,38–41].

The properly defined spectroscopic factor (sometimes referred to as “the amount of clustering”) [21,38,42–45] is given by

$$S_L = \int_0^\infty G_L^2(R) R^2 dR, \quad (23)$$

where

$$G_L(R) = \int \mathcal{N}_L^{-1/2}(R, R') g_L(R') R^2 dR' \quad (24)$$

is the modified formation amplitude. The norm kernel \mathcal{N}_L appearing in Eq. (24) is [43]

$$\begin{aligned} \mathcal{N}_L(R, R') &= \left\langle \mathcal{A} \frac{\delta(R_\alpha - R)}{R^2} \phi_\alpha [Y_L \Psi_j^D]_J \left| \mathcal{A} \frac{\delta(R_\alpha - R')}{R'^2} \phi_\alpha [Y_L \Psi_j^D]_J \right. \right\rangle. \end{aligned} \quad (25)$$

The presence of the norm kernel \mathcal{N} effectively enhances the spectroscopic factor by 1–2 orders of magnitude [21,40,41,45,46].

To compute $\mathcal{N}_L^{-1/2}(R, R')$, we expand the eigenfunctions of the norm kernel in an orthonormalized shifted Gaussian basis (SGB) [43],

$$\tilde{F}_L(R, R_k) = \sum_{k'} (N_F^{-1/2})_{kk'} F_L(R, R_{k'}), \quad (26)$$

with R_k equidistant mesh points in the interval $(0, R_{\max})$ and $k = 1, \dots, M$, where M is the dimension of the basis. The SGB is given by

$$F_L(R, R_k) = 4\pi \left(\frac{8\beta'}{\pi} \right)^{3/4} e^{-4\beta'(R^2 + R_k^2)} i^L j_L(-i8\beta' R R_k), \quad (27)$$

while the SGB overlap $(N_F)_{kk'}$ is given by

$$\begin{aligned} (N_F)_{kk'} &= \int F_L^*(R, R_k) F_L(R, R_{k'}) R^2 dR \\ &= 4\pi e^{-2\beta'(R_k^2 + R_{k'}^2)} i^L j_L(-i4\beta' R_k R_{k'}). \end{aligned} \quad (28)$$

Using the SGB overlaps, the eigenvalue equation for the norm matrix can be expressed in the form

$$\sum_{k'}^M \mathcal{N}_{kk'}^{\tilde{F}} c_{k'}^v = n_v c_k^v, \quad (29)$$

where

$$\mathcal{N}_{kk'}^{\tilde{F}} = \sum_{nn'}^M (N_F^{-1/2})_{kn} \mathcal{N}_{nn'}^F (N_F^{-1/2})_{n'k'} \quad (30)$$

For $\beta' = 4\beta$, the core-projected norm \mathcal{N}^F in Eq. (30) reduces to a simple expression [21,43,47]:

$$\mathcal{N}_{kk'}^F = \left(\langle \psi_k^{(v),L} | \psi_{k'}^{(v),L} \rangle \right)^2 \left(\langle \psi_k^{(\pi),L} | \psi_{k'}^{(\pi),L} \rangle \right)^2, \quad (31)$$

where

$$\begin{aligned} \langle \psi_k^{(\mu),L} | \psi_{k'}^{(\mu),L} \rangle &= \langle \phi_k^L | \phi_{k'}^L \rangle - \sum_{nlj\tau \in \text{core}} \delta_{lL} \langle \phi_k^l | R_{nlj} \rangle \langle R_{nlj} | \phi_{k'}^l \rangle, \end{aligned} \quad (32)$$

with $\phi_k^L(R) = F_L(R, R_k)(\beta' \rightarrow \beta)$ and $R_{nlj}(R) = u_{nlj}/R$ are the radial s.p. wave functions of the core.

In terms of eigenstates c_k^v of (29), the spectral representation of the norm kernel can be written as

$$\mathcal{N}_L^{-1/2}(R, R') = \sum_{\substack{v \\ (n_v > n_{\min})}} n_v^{-1/2} u_v^L(R) u_v^L(R'), \quad (33)$$

where the eigenfunctions $u_v^L(R)$ of the norm kernel are

$$u_v^L(R) = \sum_k^M c_k^v \tilde{F}_L(R, R_k), \quad (34)$$

and n_{\min} represents the usual cutoff on the eigenvalue of the norm kernel. The final expression for the modified formation amplitude in the normalized SGB becomes [43]

$$G_L(R) = \sum_{\substack{v \\ (n_v > n_{\min})}} n_v^{-1/2} u_v^L(R) g_v^L \quad (35)$$

with

$$g_v^L = \int u_v^L(R) g_L(R) R^2 dR. \quad (36)$$

III. THE MODEL

A. Single-particle space

The s.p. space is spanned on resonant states of a WS + Coulomb average potential. The parameters of the s.p. Hamiltonian, namely the WS potential depth V_0 , spin-orbit potential depth V_{so} , diffuseness $a(=a_{\text{so}})$, radius $r_0(=r_{0,\text{so}})$, and the radius of the uniform charge distribution r_c defining the Coulomb potential are listed in Table I. The resulting neutron and proton s.p. energies for ^{208}Pb and ^{100}Sn are given in Tables II and III, respectively. The nucleus ^{101}Sb is proton unbound; the values in Table III are generally consistent with systematics [48]. In particular, we predict a very small splitting

TABLE I. Parameters of the average WS Hamiltonian used in this work to compute s.p. neutron and proton states of ^{208}Pb and ^{100}Sn cores.

Core	τ	V_0 (MeV)	V_{so} (MeV)	a (fm)	r_0 (fm)	r_c (fm)
^{208}Pb	n	44.40	16.5	0.70	1.27	
	p	66.04	19.0	0.75	1.19	1.27
^{100}Sn	n	51.60	11.3	0.70	1.27	
	p	52.20	10.5	0.70	1.27	1.27

between the $0g_{7/2}$ and $1d_{5/2}$ neutron shells outside ^{100}Sn , and a $0g_{7/2}$ g.s. in ^{101}Sn as suggested by recent experiment [49].

B. Two-particle interaction

The correlated two-particle wave functions $|\Psi_{2\tau,0}\rangle$ (17) have been obtained using a separable two-body $T = 1$ pairing interaction [51]:

$$\langle \nu\nu, 00 | V | \nu'\nu', 00 \rangle = -G_\tau f(\nu, \tau) f(\nu', \tau), \quad (37)$$

where

$$f(\nu, \tau) = \frac{(-)^{l_\nu}}{\sqrt{2}} \langle j_\nu || Y_0 || j_\nu \rangle I(\nu, \tau). \quad (38)$$

In Eq. (38) we used the Condon-Shortley phase convention for $\langle j_\nu || Y_0 || j_\nu \rangle$ and

$$I(\nu, \tau) = \int u_{\nu\tau}^2(r) f_\tau(r) dr. \quad (39)$$

TABLE II. The eigenstates (in MeV) of the s.p. Hamiltonian of Table I for ^{208}Pb calculated with the Gamow solver ANTI [50]. The positive-energy eigenvalues represent Gamow resonances; their imaginary energies reflect nonzero particle width.

Orbit	Neutrons	Orbit	Protons
$1g_{9/2}$	-3.926	$0h_{9/2}$	-3.784
$0i_{11/2}$	-2.797	$1f_{7/2}$	-3.542
$2d_{5/2}$	-2.072	$0i_{13/2}$	-1.844
$0j_{15/2}$	-1.883	$2p_{3/2}$	-0.690
$3s_{1/2}$	-1.438	$1f_{5/2}$	-0.518
$2d_{3/2}$	-0.781	$2p_{1/2}$	$0.491 - i0.200 \times 10^{-11}$
$1g_{7/2}$	-0.768	$1g_{9/2}$	$4.028 - i0.130 \times 10^{-7}$
$1h_{11/2}$	$2.251 - i0.026$	$0i_{11/2}$	$5.434 - i0.992 \times 10^{-8}$
$0j_{13/2}$	$5.411 - i0.009$	$0j_{15/2}$	$5.960 - i0.115 \times 10^{-7}$
		$2d_{5/2}$	$6.748 - i0.184 \times 10^{-2}$
		$3s_{1/2}$	$7.843 - i0.367 \times 10^{-1}$
		$1g_{7/2}$	$8.087 - i0.898 \times 10^{-3}$
		$2d_{3/2}$	$8.530 - i0.284 \times 10^{-1}$
		$1h_{11/2}$	$11.390 - i0.215 \times 10^{-1}$
		$0j_{13/2}$	$15.086 - i0.493 \times 10^{-2}$
		$1h_{9/2}$	$15.964 - i0.393$

TABLE III. Similar to Table II except for ^{100}Sn .

Orbit	Neutrons	Protons
$0g_{7/2}$	-10.830	$2.669 - i0.207 \times 10^{-7}$
$1d_{5/2}$	-10.674	$2.869 - i0.963 \times 10^{-5}$
$2s_{1/2}$	-9.074	$4.150 - i0.595 \times 10^{-2}$
$1d_{3/2}$	-8.927	$4.393 - i0.166 \times 10^{-2}$
$0h_{11/2}$	-5.793	$7.280 - i0.110 \times 10^{-2}$
$1f_{7/2}$	-2.346	$9.649 - i0.452$
$2p_{3/2}$	-1.531	
$2p_{1/2}$	-0.912	
$0h_{9/2}$	-0.641	$12.012 - i0.0736$
$1f_{5/2}$	-0.171	
$0i_{13/2}$	$3.254 - i0.132 \times 10^{-2}$	$15.572 - i0.185$

For the radial form factor $f_\tau(r)$ we took the derivative of the WS potential multiplied by r :

$$f_\tau(r) = \frac{r}{a_{\nu\tau}} \frac{e^{\frac{r-R_{\nu\tau}}{a_{\nu\tau}}}}{\left(1 + e^{\frac{r-R_{\nu\tau}}{a_{\nu\tau}}}\right)^2}. \quad (40)$$

In the case of ^{212}Po and ^{104}Te the two-particle amplitudes of Eq. (17) were obtained exactly in the Tamm-Dancoff approximation [52,53]:

$$X_\nu^\tau = N_0 \frac{f(\nu, \tau)}{2\epsilon_\nu^\tau - E_0^\tau}, \quad (41)$$

where ϵ_ν^τ are s.p. energies, E_0^τ is the correlated two-particle energy, and N_0 is the normalization constant fixed by the condition $\sum_\nu (X_\nu^\tau)^2 = 1$.

The parameters $R_{\nu\tau}$ and $a_{\nu\tau}$ defining the radial form factor (40) for ^{210}Pb and ^{210}Po were chosen to reproduce the wave functions used by Harada [14]. Since such data are not available for ^{102}Sn and ^{102}Te , in this case we adopted the values of the WS potential for ^{100}Sn shown in Table I. The pairing strength G_τ was adjusted to fit the experimental two-nucleon separation energies $S_{2\tau}$ through the dispersion relation

$$\frac{1}{G_\tau} = \sum_\nu \frac{f^2(\nu, \tau)}{2\epsilon_\nu^\tau - E_0^\tau}. \quad (42)$$

Since the proton-unbound nucleus ^{102}Te is not known experimentally, for this system we adopted the value of $S_{2p} = -2.14$ MeV obtained by extrapolating down from the heavier Te isotopes [54]. This value is in reasonable agreement with recent phenomenological estimates [48]. Table IV lists the parameters of the residual interaction used in our study.

TABLE IV. Parameters $R_{\nu\tau}$ and $a_{\nu\tau}$ of the residual interaction (37). The last column lists the value of $S_{2\tau}$ that has been used to constrain the pairing strength G_τ for various configuration spaces considered.

Nucleus	R_ν (fm)	a_ν (fm)	$S_{2\tau}$ (MeV)
^{210}Pb	7.525	0.70	9.123
^{210}Po	5.451	0.75	8.783
^{102}Sn	5.895	0.70	24.3
^{102}Te	5.895	0.70	-2.14

TABLE V. Model spaces used in this work to describe ^{212}Po alpha decay.

Model	Neutron States	Proton States
M0	$1g_{9/2}$	$0h_{9/2}$
M1	$1g_{9/2}, 0i_{11/2}, 2d_{5/2}, 0j_{15/2}$ $3s_{1/2}, 2d_{3/2}, 1g_{7/2}$	$0h_{9/2}, 1f_{7/2}, 0i_{13/2}, 2p_{3/2}$ $1f_{5/2}, 2p_{1/2}$
M2	$1g_{9/2}, 0i_{11/2}, 2d_{5/2}$	$0h_{9/2}, 1f_{7/2}, 0i_{13/2}$
M3	$1g_{9/2}, 0i_{11/2}, 2d_{5/2}, 0j_{15/2}$	$0h_{9/2}, 1f_{7/2}, 0i_{13/2}$
M4	$1g_{9/2}, 0i_{11/2}, 2d_{5/2}, 0j_{15/2}$ $3s_{1/2}, 2d_{3/2}, 1g_{7/2}, 1h_{11/2}$ $0j_{13/2}$	$0h_{9/2}, 1f_{7/2}, 0i_{13/2}, 2p_{3/2}$ $1f_{5/2}, 2p_{1/2}, 1g_{9/2}, 0i_{11/2}$ $0j_{15/2}, 2d_{5/2}, 3s_{1/2}, 1g_{7/2}$ $2d_{3/2}, 1h_{11/2}, 0j_{13/2}, 1h_{9/2}$

C. Configuration space

To study the dependence of the formation amplitude on the size of valence space, and to compare with previous work, we considered several model spaces. Those used in the description of the alpha decay of ^{212}Po are given in Table V. The model space M0 contains only one valence shell. The space M1 contains one major shell, including the unusual-parity intruder orbit. The model space M2 is that used by Harada [14]. The model space M3 is that of Glendenning and Harada [55]. Finally, M4 is the extended shell model space employed by Tonzuka and Arima. The model spaces used to describe ^{104}Te alpha decay are shown in Table VI; M1 consists of one major shell, including the unusual-parity intruder orbit, while M4 consists of states with width less than 1 MeV.

D. Wave functions

For the alpha formation amplitude in ^{212}Po discussed in Sec. IV we considered the model spaces M2, M3, and M4. The wave function amplitudes in M2 were taken from Refs. [14,56]. For calculations in M3, we took the $T = 1$ seniority-zero amplitudes of Ref. [55] and renormalized them accordingly. For calculations in the extended space M4, we used the renormalized amplitudes of Ref. [15]; here we retained only configurations having width smaller than 1 MeV. The comparison between ^{212}Po and ^{104}Te discussed in Sec. VI was carried out in the model spaces M1 and M4. The corresponding wave functions were calculated in the two-particle approximation described in Sec. III B, except for ^{212}Po in the M4 model space, where Ref. [55] was used instead.

TABLE VI. Model spaces used in this work to describe ^{104}Te alpha decay.

Model	Neutron States	Proton States
M1	$0g_{7/2}, 1d_{5/2}, 2s_{1/2}, 1d_{3/2}$ $0h_{11/2}$	$0g_{7/2}, 1d_{5/2}, 2s_{1/2}, 1d_{3/2}$ $0h_{11/2}$
M4	$0g_{7/2}, 1d_{5/2}, 2s_{1/2}, 1d_{3/2}$ $0h_{11/2}, 1f_{7/2}, 2p_{3/2}, 2p_{1/2}$ $0h_{9/2}, 1f_{5/2}, 0i_{13/2}$	$0g_{7/2}, 1d_{5/2}, 2s_{1/2}, 1d_{3/2}$ $0h_{11/2}, 1f_{7/2}, 0h_{9/2}, 0i_{13/2}$

E. Penetration factor

The s.p. alpha width Γ_0^{sp} has been obtained from the current expression (20). The alpha-core potential was assumed to be of a WS + Coulomb form with the parameters of Ref. [57]: $r_0 = R_c = 1.315$ fm, $a = 0.65$ fm. The strength of the WS potential has been adjusted to reproduce the measured Q_α value corrected by the electron screening term [6,13,58–60]:

$$Q_\alpha = E_\alpha \frac{A_P}{A_D} + \Delta E_{sc}, \quad (43)$$

where

$$\Delta E_{sc} = 65.3Z_P^{1.4} - 80Z_P^{0.4} \text{ (eV)}. \quad (44)$$

For ^{212}Po , $E_\alpha = 8.785$ MeV [54] and $\Delta E_{sc} = 31.8$ keV; hence, $Q_\alpha = 8.986$ MeV. The g.s. alpha decay of ^{104}Te has not been observed. For that reason, we took the value $Q_\alpha^{\text{BE}} = 5.135$ MeV extrapolated down from the binding energy differences in ^{108}Te (3.445 MeV) and ^{106}Te (4.290 MeV) [61]. By adding the screening correction $\Delta E_{sc} = 16.1$ keV, we arrived at $Q_\alpha = 5.151$ MeV. The resulting WS potential strength is $V_0 = 143.49$ MeV for ^{212}Po and 149.64 MeV for ^{104}Te .

The Gamow wave functions, obtained by means of the code ANTI [50], were normalized using the complex rotation method proposed in Ref. [62]. The complex energy of the metastable alpha state is $\mathcal{E}_\alpha = (8.986 - i0.632 \times 10^{-13})$ MeV for ^{212}Po and $\mathcal{E}_\alpha = (5.151 - i0.814 \times 10^{-13})$ MeV for ^{104}Te . The outgoing spherical Coulomb-Hankel function H^+ was calculated using the code in Ref. [63].

F. Calculation of the spectroscopic factor

The radial integration in the expressions for the spectroscopic factor (23) and the formation amplitude in the normalized SGB (36) have been carried out using 200 Gauss-Legendre mesh-points with the maximum radius of 20 fm.

The s.p. core wave functions entering Eq. (32) are those of the s.p. Hamiltonian of Table I. The radial mesh R_k defining the normalized SGB (26) was taken at equidistant points $R_k = k \Delta R$. In order to determine the step ΔR we expanded the s.p. core states $u(r)$ in the normalized SGB: $\tilde{u}(r) \equiv \sum_{k=1}^M a_k [r \tilde{F}_0(r, R_k)]$. Under the condition that $u_{\text{diff}}(r) = |u(r) - \tilde{u}(r)| < 0.005$ fm $^{-1/2}$ we found that $0.44 \lesssim \Delta R \lesssim 0.57$ fm and $R_{\text{max}} \gtrsim 14$ fm. For this range of ΔR and R_{max} the normalized SGB is orthonormal with an accuracy better than 10^{-9} . To illustrate the quality of the resulting expansion, Fig. 1 shows $u_{\text{diff}}(r)$ for the neutron core states in ^{208}Pb .

To calculate the modified formation amplitude $G(R)$, one needs to determine the eigenvalue cutoff n_{min} . To this end, we show in Figs. 2 and 3 typical distribution of the eigenvalues n_ν of the norm kernel (25) for ^{212}Po and ^{104}Te , respectively, for different values of ΔR . One may observe that a significant fraction of them accumulate at zero [53,64]. To eliminate these spurious eigenvectors, we define the cutoff at the value where the eigenvalue distribution changes slope. For ^{212}Po and ^{104}Te this happens at n_ν around 10^{-3} . Consequently, in our calculations, we adopt the cutoff value of $n_{\text{min}} = 0.001$.

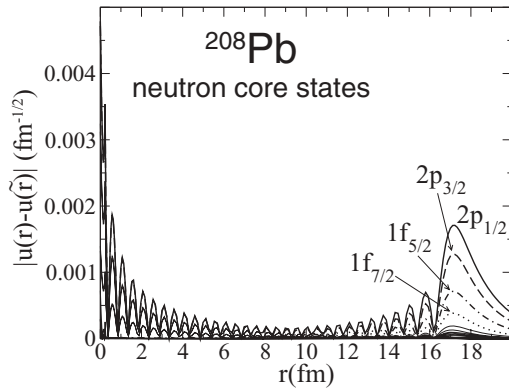


FIG. 1. $u_{\text{diff}}(r) = |u(r) - \bar{u}(r)|$ for the neutron core states in ^{208}Pb for $\Delta R = 0.5$ fm, $M = 30$, and $R_{\text{max}} = 15$ fm.

The eigenfunctions $u_v^L(R)$ of the norm kernel (34) are orthonormal with an accuracy of 10^{-10} for all eigenvalues. The eigenfunctions with $n_v < n_{\text{min}}$ oscillate inside the nuclear volume and vanish outside the surface region. To further check the quality of $u_v^L(R)$ we compute expression (35) by assuming $n_{\text{min}} = 0$ and $n_v = 1$ for all v . In this case, Eq. (35) formally reduces to $g(R)$. Figure 4 shows $g(R)$ for ^{212}Po calculated in this way. The agreement with the original formation amplitude is excellent, except for a small deviation close to $R = 0$ and a small oscillation around and beyond the nuclear surface, which is not visible in the scale of Fig. 4.

Next we study the sensitivity of \mathcal{S} to the choice of R_{max} , ΔR , and n_{min} . For this analysis we relax the condition for $u_{\text{diff}}(r)$ in order to access a wider range of ΔR . First, we study the sensitivity of \mathcal{S} as a function of R_{max} for various values of ΔR . Figure 5 shows the result for $0.53 \leq \Delta R \leq 0.59$ fm for ^{212}Po in the model space M4 and $n_{\text{min}} = 0.001$. Except for a small value of $\Delta R = 0.53$ fm, which does not produce stable results, a plateau in R_{max} is reached around 14 fm.

The dependence of \mathcal{S} on ΔR displayed in Fig. 5 reflects the fact that for too small values of the step the basis functions become numerically linearly dependent, while for too large ΔR 's the basis cannot capture high Fourier components [43, 53, 64]. Figure 6 shows \mathcal{S} for ^{212}Po in the model space M4 and $n_{\text{min}} = 0.001$ as a function of ΔR . In general, appreciable

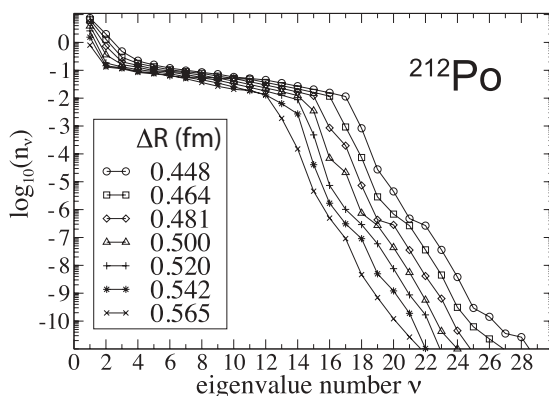


FIG. 2. Eigenvalues of the norm kernel (25) for ^{212}Po for $R_{\text{max}} = 13$ fm for different values of ΔR .

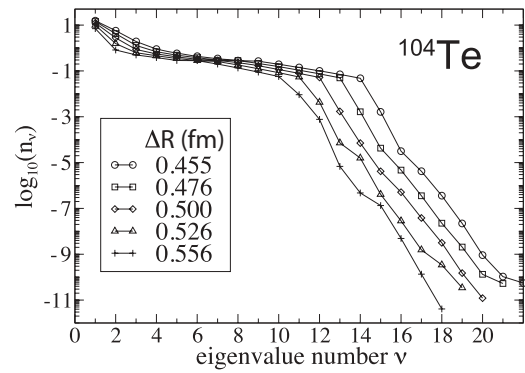


FIG. 3. Similar as in Fig. 2 except for ^{104}Te and $R_{\text{max}} = 10$ fm.

oscillations of \mathcal{S} can be seen except for the “safe” region $0.54 \leq \Delta R \leq 0.59$ fm, where results weakly depend on R_{max} .

Finally, Fig. 7 shows the behavior of \mathcal{S} as a function of the eigenvalue cutoff n_{min} for $\Delta R = 0.57$ fm. The cutoff used in Figs. 5 and 6 corresponds to $n_{\text{min}}^{-1/2} = (0.001)^{-1/2} \approx 31.5$.

G. Integral over intrinsic coordinates

The multidimensional integral (12) depends on the nucleonic coordinates, which are parametrized in terms of intrinsic variables through Eqs. (9) and (10). The integration over φ_i can easily be done analytically. Since the coordinates of particles 1 and 2 depend only on the relative coordinates 1 and 3, and the particle coordinates 3 and 4 depend only on the relative coordinates 2 and 3, one can greatly simplify the remaining six-dimensional integral by making first the integration over the relative coordinates 1 and 2 and then the integration over the coordinate 3:

$$\int d\rho_3 \left[\cdots \left(\int \cdots d\rho_1 \right) \left(\int \cdots d\rho_2 \right) \right]. \quad (45)$$

The integration has been carried out using the Gauss-Legendre quadrature using ten points for the radial integrals and eight points for the the angular coordinates. This guarantees the convergence up to the fourth significant digit.

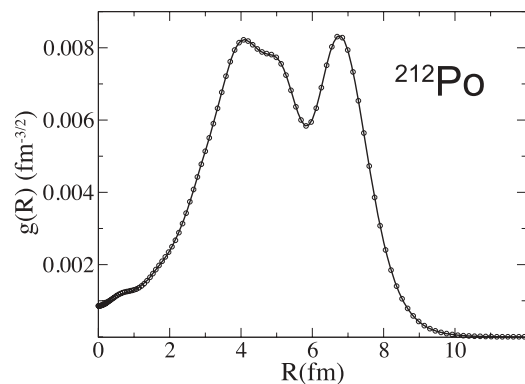


FIG. 4. Formation amplitude $g(R)$ for ^{212}Po in the M4 model space expanded in eigenfunctions of the norm kernel for $\Delta R = 0.500$ fm and $R_{\text{max}} = 13$ fm.

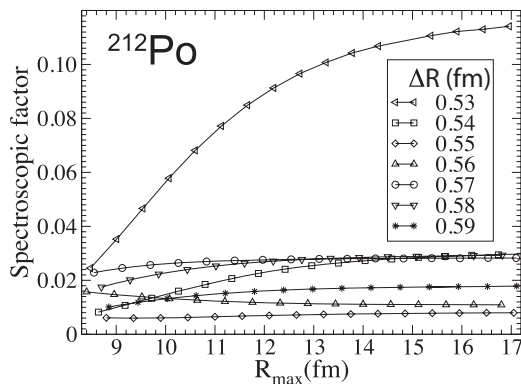


FIG. 5. Convergence of \mathcal{S} for ^{212}Po (model space M4) as a function of R_{max} of the normalized SGB for different values of ΔR (with $n_{\text{min}} = 0.001$).

IV. REDUCED WIDTH FOR ^{212}Po

A. Single- j configuration

Following Rasmussen [32], it is instructive to compute relative reduced widths assuming a pure single- j_n shell-model orbital assignment for the neutron pair, while the proton pair fills the $0h_{9/2}$ shell. For simplicity, the results are expressed relative to the ^{210}Po reference (a neutron pair in $2p_{1/2}$).

In the δ -function approximation of Sec. II B, the ratio r_δ of the reduced widths is given by a simple expression [32]:

$$r_\delta = \frac{\gamma_{j_n}^2}{\gamma_{2p_{1/2}}^2} = \frac{2j_n + 1}{2} \left(\frac{u_{j_n}(R)}{u_{2p_{1/2}}(R)} \right)^4. \quad (46)$$

In a more general case expressed by Eq. (11), the ratio r depends on the proton wave function:

$$r = \frac{\gamma_{j_n, 0h_{9/2}}^2}{\gamma_{2p_{1/2}, 0h_{9/2}}^2} = \frac{2j_n + 1}{2} \left(\frac{I_{j_n, 0h_{9/2}}(R)}{I_{2p_{1/2}, 0h_{9/2}}(R)} \right)^2. \quad (47)$$

Table VII compares the ratio r_δ given by Eq. (46) using the WS wave functions with that of Table I of Rasmussen [32] based on the rounded square well potential of Blomqvist and Wahlborn [65] for several neutron configurations at $R = 9.5$ fm. We find excellent agreement between these two

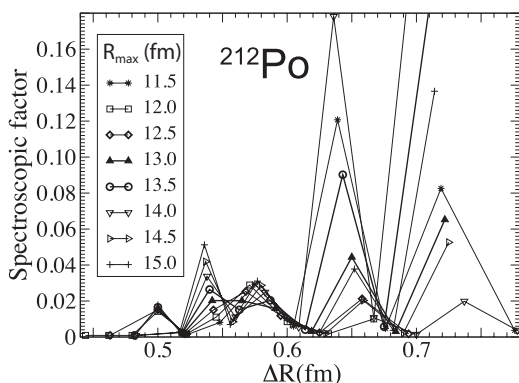


FIG. 6. Similar as in Fig. 5 but as a function of the step size ΔR for different values of R_{max} .

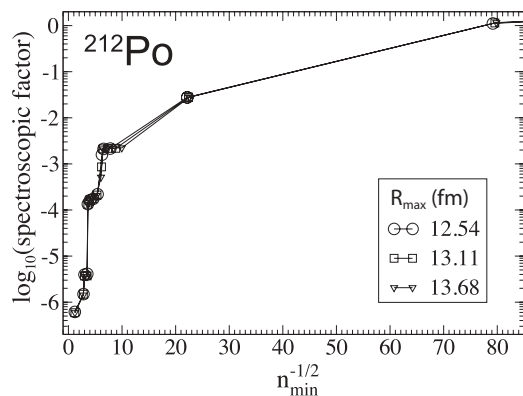


FIG. 7. Similar as in Fig. 5 but as a function of n_{min} for different values of R_{max} and $\Delta R = 0.57$.

calculations, and we checked that this agreement also holds for $R = 9.0$ fm. This is not surprising as both calculations employ finite-depth potentials. The fourth column of Table VII displays the ratio r given by Eq. (47) using the WS wave functions; they are compared with the h.o. values of Ref. [66] (last column). It is seen that h.o. calculations underestimate WS values for high- j orbits by a factor 2–3.

It has been early recognized [28,32] that the δ -function approximation overestimates the contributions of high- j orbitals. One can see it clearly by comparing the values of r_δ of Eq. (46) with those of r (47), i.e., the third and fourth columns of Table VII. To cure this deficiency, a correction factor B_v (15) was introduced [32] in Eq. (14) that depends on the relative angular momentum.

B. Enhancement due to configuration mixing

As was first shown by Harada [14], the reduced width at the surface region is strongly enhanced by the configuration mixing because contributions from various shell-model orbits add coherently. To assess the effect of collective enhancement due to the configuration mixing, we carried out calculations in the M2 space. For $R = 8$ fm, our WS calculations yield the enhancement factor of $\zeta = 8.5$ with respect to the valence-shell configuration M0. This is to be compared with $\zeta = 11$ obtained in the δ -function approximation; $\zeta = 10$ obtained by Rasmussen [32]; and $\zeta = 5.5$ of Harada [14] using h.o. wave functions.

TABLE VII. Single- j alpha reduced width ratios at $R = 9.5$ fm. Shown are: r_δ of Ref. [32], r_δ of Eq. (46), r of Eq. (47), and r of Ref. [66].

Orbital j_n	r_δ [32]	r_δ	r	r [66]
$0i_{13/2}$	0.44	0.46	0.20	0.10
$0i_{11/2}$	0.32	0.34	0.21	0.08
$1g_{9/2}$	7.50	7.50	6.50	3.73
$1f_{5/2}$	0.73	0.74	0.58	0.55
$2p_{3/2}$	1.89	1.89	1.73	1.89

TABLE VIII. The relative reduced width θ^2 (48) obtained in Ref. [15] and this work.

Model space	R (fm)	Ref. [15]	This work
M0	8.4	6.3×10^{-6}	0.60×10^{-6}
M3	8.5	4.4×10^{-5}	0.48×10^{-5}
M4	9.0	2.9×10^{-4}	0.41×10^{-4}

For the model space M3 of Glendenning and Harada [55], obtained by adding the intruder neutron state $0j_{15/2}$ to M2, we obtain $\zeta = 21$. This should be compared with $\zeta = 24$ obtained in the δ -function approximation and $\zeta = 30$ obtained in Ref. [55] (also within the δ -function approximation) using a fairly rich wave function that also includes proton-neutron correlations and $J > 0$ two-particle couplings. It is worth noting that our enhancement is around 80% of that by Glendenning and Harada, and that the seniority-zero component in their wave function is also 80%.

C. Extended shell-model space

Due to the strong collective enhancement of the reduced width due to configuration mixing, it is important to consider extended shell-model space by taking into account higher-lying orbitals [15]. For finite-depth shell-model potentials, such as the WS potential used in this study, this necessitates a proper treatment of the particle continuum. An appropriate representation to deal with the continuum space is the complex Berggren ensemble representing bound and unbound s.p. states [67,68].

Here we consider the large configuration space M4 of Tonzuka and Arima [15], i.e., all s.p. orbits up to the $N = 7$ harmonic oscillator shell except for broad resonances with widths greater than 1 MeV. The shell-model amplitudes were taken from Ref. [15] and renormalized to the reduced model space. For the sake of comparison with Ref. [15], we consider the relative reduced width

$$\theta^2(R) = \frac{\gamma^2(R)}{\gamma_W^2(R)}, \quad (48)$$

where $\gamma_W^2(R) = \frac{3\hbar^2}{2\mu R^2}$ is the Wigner limit [69].

Table VIII compares our WS results for $\theta^2(R)$ with those of Ref. [15] obtained in the h.o. basis for several values of R . Generally, the reduced width obtained in the WS model is about one order of magnitude smaller than that in the h.o. basis. This is because the h.o. basis knows nothing about the particle thresholds, and the radial behavior at large distances is solely determined by the oscillator length. For that reason, calculations based on the h.o. wave functions show large sensitivity to this parameter [70].

The formation amplitude obtained in this work is shown in Fig. 8 for the configuration spaces M0, M1, M2, M3, and M4. Compared with the formation amplitudes of Ref. [15], the maximum of the formation amplitudes obtained in the WS model are significantly larger, and appear at lower values of R , than in the h.o. model. Also the overall shape of the formation amplitude is very different in the two cases. A characteristic

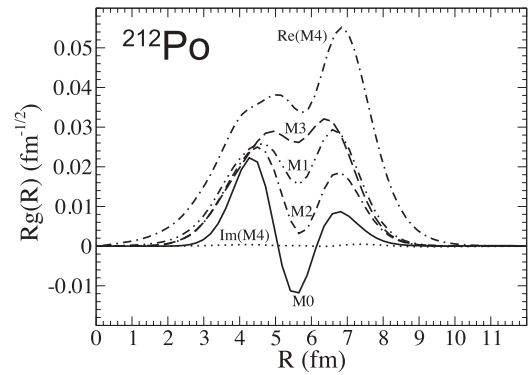


FIG. 8. Formation amplitude $g(R)$ for ^{212}Po obtained in this work in the model spaces M0 to M4 as defined in Table V. The imaginary part of the formation amplitude in M4 (dotted line) is also shown.

two-humped shape of $g(R)$ calculated in M4 resembles the formation amplitude $G(R)$ obtained in Refs. [15,42]. A similar result was also obtained in Refs. [20,21]. It is indeed interesting to see that a two-humped behavior of the formation amplitude for ^{212}Po has been obtained by considering a large configuration space and the Berggren ensemble of the WS potential.

Figure 8 also shows that the formation amplitude in the M4 model space has a small imaginary part. This is because our calculations are carried out in the pole approximation that ignores the nonresonant continuum [68,71–73]. This spurious component of $g(R)$ results in a very small imaginary contribution to the reduced width, which can be safely neglected considering the expected accuracy of our model.

V. ABSOLUTE ALPHA-DECAY WIDTH OF ^{212}Po

The g.s. alpha-decay width of ^{212}Po has been determined in the seniority-zero approximation using three different model spaces listed in Table V: M0, M3, and M4. The corresponding four-particle shell-model wave function contains one configuration in the M0 space, 12 configurations in M3, and 144 seniority zero configurations in M4.

The absolute width from Eq. (1) should not depend on the channel radius R . However, in R -matrix studies involving approximations, such as the one-channel R -matrix treatment, this condition cannot be met [22]. Therefore, in practical calculations, in which the dependence of Γ_L on R around the nuclear surface is small relative to the appreciable R dependence of the formation amplitude, one is trying to meet the plateau condition for $\Gamma_L(R)$ in which the absolute width varies weakly around the nuclear surface [74]. Figure 9 shows the dependence of the R -matrix width (1) on the channel radius. It is seen that the plateau condition is met only in the case of the extended configuration space M4 involving particle continuum. Here, we find a fairly weak variation of $\Gamma(R)$ between 7 and 11 fm.

As seen in Fig. 9, and discussed in Sec. IV B and Refs. [15, 70,75], the width strongly increases with the size of the shell-model space. Indeed, in the surface region, $\Gamma(R)$ obtained in M3 shows an enhancement of ~ 15 with respect to M0, and in

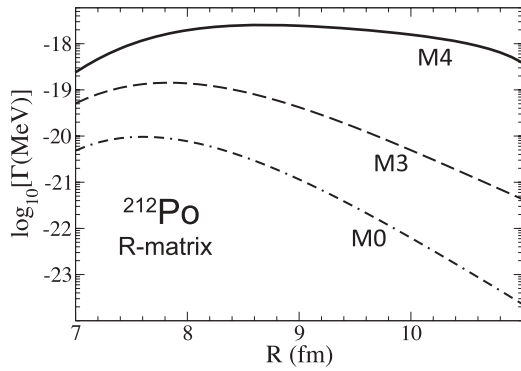


FIG. 9. Dependence of the absolute alpha-decay width (1) of ^{212}Po on the R -matrix channel radius R for three different model spaces M0, M3, and M4.

the extended space M4 the enhancement is ~ 260 . Compared to experimental value, however, the width obtained in M4 is still 600 times smaller than the experimental value $\Gamma_{\text{exp}} = 0.153 \times 10^{-14}$ MeV [54].

A further enhancement in the reduced width is due to the antisymmetrization and normalization of the channel decay [38,42]. This is achieved by replacing the standard formation amplitude $g(R)$ with the modified formation amplitude $G(R)$ of Eq. (24). Figure 10 shows $G(R)$ calculated in the M4 model space with $\Delta R = 0.56$ fm, $R_{\text{max}} = 11.76$ fm ($M = 21$), and $n_{\text{min}} = 0.001$. A small oscillation at the tail of $G(R)$ can be seen. The amplitude of this oscillation, around the asymptotic behavior given by $H_0^+(\eta, kR)$, varies very little with R_{max} for this value of ΔR . As discussed in, e.g., Refs. [3,42,45], the behavior of $g(R)$ and $G(R)$ is generally very different. This can be seen by comparing Figs. 8 and 10.

The absolute alpha-decay width obtained by using the R -matrix expression (1) with the formation amplitude $G(R)$ of Fig. 10 is shown in Fig. 11. There appears a small plateau in the region of nuclear surface that corresponds to $\Gamma \approx 0.0042 \times 10^{-14}$ MeV. This value is ~ 36 times smaller

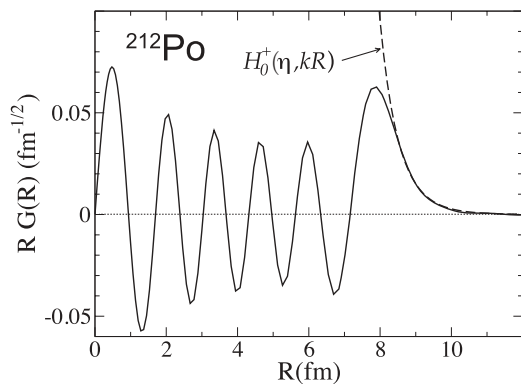


FIG. 10. Modified formation amplitude $G(R)$ of Eq. (24) in the extended model space M4 with $n_{\text{min}} = 0.001$, $\Delta R = 0.56$ fm, and $R_{\text{max}} = 11.76$. Unlike $g(R)$, $G(R)$ properly accounts for the normalization and antisymmetrization of the decay channel. The asymptotic behavior of $G(R)$ is given by the Coulomb-Hankel function at the alpha-decay energy $Q_\alpha = 8.986$ MeV (dashed line).

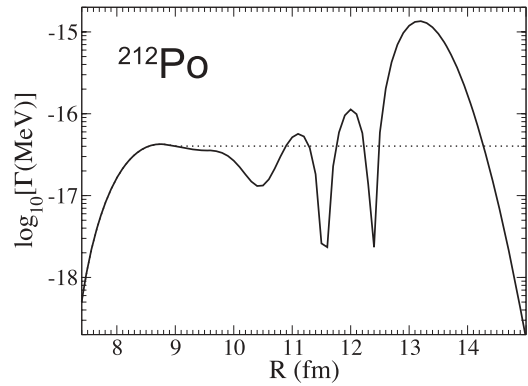


FIG. 11. Absolute width from R -matrix expression (1) calculated in the M4 model space using the modified formation amplitude $G(R)$ of Fig. 10. At $R > 9$ fm, the result obtained by assuming $G(R) \propto H_0^+(\eta, kR)$ is marked by a dotted line.

than Γ_{exp} . At larger distances $R > 9$ fm, the result is affected by spurious oscillations of $G(R)$ around $H_0^+(\eta, kR)$; i.e., it is quite unreliable.

The absolute width can also be obtained from expression (2), which involves the alpha-particle spectroscopic factor \mathcal{S} and the s.p. decay width. Figure 12 shows the result of the current expression (20) for Γ^{sp} as a function of the channel radius. As discussed in Ref. [36], Γ^{sp} calculated this way should be independent of R if R is large enough. This is precisely what is seen in Fig. 12: the s.p. width converges beyond the range of the WS potential to $\Gamma^{sp} = 0.1247 \times 10^{-12}$ MeV, which is indeed very close to the value of $-2\text{Im}(\mathcal{E}_\alpha) = 0.1265 \times 10^{-12}$ MeV given by the imaginary part of the Gamow resonance.

Using the modified formation amplitude $G(R)$ of Fig. 11, we compute the spectroscopic factor $\mathcal{S} = 0.011$, which—combined with the value of Γ^{sp} above—yields $\Gamma = 0.14 \times 10^{-14}$ MeV. Using $\Delta R = 0.55$ fm we obtain $\mathcal{S} = 0.0080$ and $\Gamma = 0.10 \times 10^{-14}$ MeV. Both these values are close to $\Gamma_{\text{exp}} = 0.153 \times 10^{-14}$ MeV.

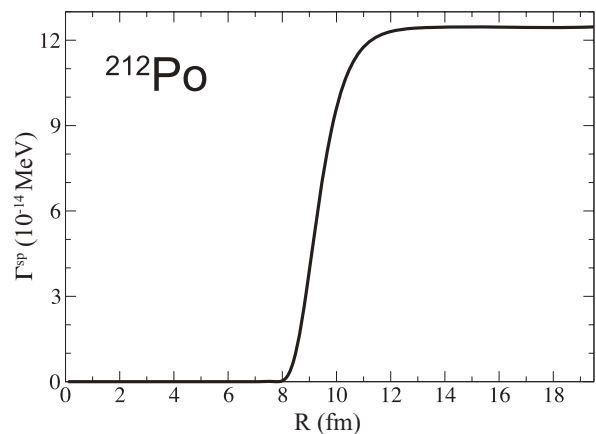


FIG. 12. Single-particle width of ^{212}Po from current expression (20).

TABLE IX. Alpha-decay spectroscopic factor and absolute width for ^{212}Po and ^{104}Te computed in the configuration spaces M1 and M4, with $n_{\min} = 0.001$, $\Delta R = 0.56$, and $R_{\max} = 11.76$ fm ($M = 21$).

Model space	\mathcal{S}		$\Gamma \times 10^{14}$ MeV	
	^{212}Po	^{104}Te	^{212}Po	^{104}Te
M1	0.00032	0.00046	0.0040	0.0075
M4	0.011	0.0051	0.14	0.083

VI. COMPARISON BETWEEN GROUND-STATE ALPHA DECAY OF ^{212}Po AND ^{104}Te

To compare absolute widths of ^{212}Po and ^{104}Te in a consistent way, we consider similar M1 and M4 model spaces for both nuclei. The norm kernel eigenvalues n_ν do not depend on the model space in which $g(R)$ is calculated, so we take the cutoff $n_{\min} = 0.001$.

Let us begin with ^{212}Po by making a convergence analysis of \mathcal{S} in the M1 model space as a function of ΔR and R_{\max} (as in Fig. 5). For $\Delta R = 0.53, 0.54, 0.55$, and 0.56 fm, we found $\mathcal{S} = 0.0041, 0.0011, 0.00030$, and 0.00032 , respectively. The resulting converged value $\mathcal{S} = 0.0003$ is too small, as expected from Fig. 8. This deficiency is related to the poor quality of the interaction used to describe ^{212}Po in M1. To better understand this fact, let us take a look of the spectroscopic factor in terms of the spectral representation of the norm kernel,

$$\mathcal{S} = \sum_{\nu} \frac{g_{\nu}^2}{n_{\nu}}, \quad (49)$$

where the sum is truncated by the condition $n_{\nu} > n_{\min}$. The summation range and eigenvalues n_{ν} are the same for M1 and M4; the only difference comes from g_{ν} . Because of the rapid oscillation of the eigenfunctions inside the nucleus, only the eigenfunctions which are peaked at and beyond the nuclear surface will contribute significantly to the sum. But—because $g(R)$ in M1 is small in the surface region—the overlap with those eigenfunctions is small, and this gives rise to a very reduced value of \mathcal{S} .

By making a similar analysis for ^{104}Te in M1, we found $\mathcal{S} = 0.067, 0.024, 0.0066$, and 0.00046 for $\Delta R = 0.53, 0.54, 0.55$, and 0.56 fm, respectively. In the model space M4 we found $\mathcal{S} = 0.21, 0.088, 0.032$, and 0.0051 for the same values of ΔR . Clearly, the convergence in \mathcal{S} has not been achieved for ^{104}Te . We would like to attribute this to the impact of the proton continuum on g_{ν} , which results in increased oscillations of $G(R)$ in the surface area. Table IX compares the values of \mathcal{S} and the corresponding absolute widths for ^{212}Po and ^{104}Te at $\Delta R = 0.56$ fm. (The single-particle width for Te is $\Gamma^{sp} = 0.162 \times 10^{-12}$ MeV.)

It is interesting to compare our current results for ^{104}Te with the estimates of phenomenological alpha-decay models based on a semiclassical approximation [76–78]. The assumed large value of $Q_{\alpha} = 6.12$ MeV in Ref. [76] results in a very short half-life of 7×10^{-11} s. The alpha-decay energies of 5.05 MeV [77] and 5.42 ± 0.07 MeV [78] result in $T_{1/2} \sim 10^{-7}$ s and $\sim 5 \times 10^{-9}$ s, respectively, and these estimates are not inconsistent with our value (M4 model space)

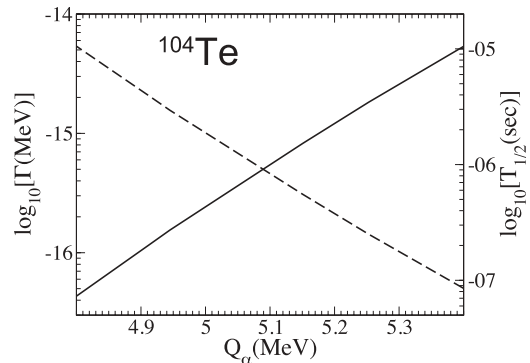


FIG. 13. Ground-state alpha-decay width (left scale) and half-life (right scale) in ^{104}Te as functions of the decay energy.

$T_{1/2} = 5.5 \times 10^{-7}$ s ($Q_{\alpha} = 5.151$ MeV). As the value of Q_{α} in ^{104}Te is very uncertain, we show in Fig. 13 the absolute width and half-life $T_{1/2}$ as a function of Q_{α} for the model space M4.

Our predicted spectroscopic factors in M4 for ^{104}Te and ^{212}Po are about 0.5% and 1%, respectively. As mentioned above, a fairly small value of \mathcal{S} in ^{104}Te could be a consequence of the proximity of the proton continuum. Indeed, all the valence proton shells are resonances. The small value of \mathcal{S} in ^{104}Te could also be attributed to the poor quality of the valence interaction assumed, and the neglect of the $T = 0$ force. The effect of the proton-neutron interaction was examined in, e.g., Refs. [14,79] for ^{212}Po and was found to be minor due to the fact that neutrons and protons in ^{212}Po occupy different shells. This is no longer true in the $N = Z$ nucleus ^{104}Te , in which the major enhancement of \mathcal{S} is expected due to $T = 0$ correlations. Therefore, our predictions for \mathcal{S} and Γ in ^{104}Te given in Table IX should be considered as a very conservative lower limit.

VII. CONCLUSIONS

The g.s. alpha decay of ^{212}Po has been studied within the complex-energy shell-model framework with the Berggren ensemble of the average Woods-Saxon potential. We applied the pole approximation by considering s.p. resonant states only. The overlap integral involving alpha-cluster nucleons was computed exactly, without resorting to the δ -function approximation. We considered the large valence space of Tonozuka and Arima that is necessary to produce the collective enhancement of the formation amplitude.

The absolute alpha-decay width was computed using the reduced width obtained in the framework of the R -matrix theory and also from the alpha spectroscopic factor. The latter approach yielded results consistent with experimental value, but only after considering the antisymmetrization and normalization of the decay-channel wave function. The R -matrix estimate underestimates the experimental width by a factor of ~ 36 . The R -matrix expression depends on the asymptotic value of the formation amplitude, which is very sensitive to the size of the configuration space. On the other hand, the reaction-theory expression (2) involves the

spectroscopic factor—an integral quantity that depends less on the size of the basis used. It is very encouraging to see that a reasonable agreement with the experimental width of ^{212}Po has been obtained without explicitly considering the alpha-cluster component in the wave function of the parent nucleus. In this context, we believe that the improved treatment of the particle continuum has been essential.

We have also provided an estimate of the alpha-decay rate in ^{104}Te . Unfortunately, due to the fact that the valence proton shells in this nucleus lie in the continuum, no fully convergent result has been achieved. We hope to improve the situation in the future by inclusion of the nonresonant continuum space that will remove some of the undesired oscillations in $G(R)$ at large distances. In addition, since the residual interaction employed in our work neglects the proton-neutron components, and the wave function has a seniority-zero character based on $T = 1$ nucleonic pairs, the predicted alpha width in this $N = Z$ nucleus should be viewed as a conservative low limit. Indeed, the inclusion of $T = 0$ correlations is expected to increase the value of Γ significantly.

The calculations presented in this study should be considered as an important step towards an improved microscopic understanding of the alpha-decay process. Still, as this work

demonstrates, further improvements are needed. The neglect of the nonresonant continuum, i.e., complex-energy scattering states in the Berggren ensemble, slightly violates the completeness relation at a one-body level. This results in small imaginary contributions to spectroscopic factors and reduced widths, and—most importantly—can affect the behavior of formation amplitudes at very large distances. The second crucial development will be the use of large-scale shell-model calculations, including realistic $T = 0$ and $T = 1$ interactions, to compute wave function amplitudes. This will enable us to provide a more meaningful estimate of ^{104}Te alpha-decay rate. The work in both directions is underway.

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