

Effects of nuclear masses from different mass models on α -decay properties of superheavy nuclei*

Yao-Hui Ding (丁耀辉) Niu Wan (万牛)[†]

School of Physics and Optoelectronics, South China University of Technology, Guangzhou 510641, China

Abstract: Using four representative nuclear mass models, namely WS4, FRDM, DZ10, and KTUY, we perform a systematic investigation on how nuclear masses affect α -decay properties of superheavy nuclei, including decay energies, α -cluster preformation factors, and corresponding half-lives. The α -cluster preformation factors are obtained from two types of cluster-formation model (CFM) and also extracted from experimental decay half-life. All mass models reproduce the known α -decay energies with small root-mean-square errors, while WS4 and FRDM show the highest accuracy. Strong correlations among preformation factors from different mass models are identified in both CFM and extracted results, although the exponential dependence of half-lives on decay energy weakens correlations between the two approaches. For possible α -decay chains of superheavy nuclei, the decay energy systematically decreases and the predicted half-life increases with the decreasing proton number. This tendency is also observed from the calculations for superheavy isotopes with same proton number. These results indicate that isotopes of superheavy elements with more neutrons are expected to exhibit enhanced stability, providing theoretical reference for future synthesis of elements with $Z=119$ and 120 .

Keywords: alpha decay, nuclear mass model, superheavy nuclei, preformation factor

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I. INTRODUCTION

The study of radioactive decay remains one of the most fundamental areas in nuclear physics, which provides direct insights into nuclear structure, stability, and the limits of the nuclear landscape. As one significant decay mode in heavy and superheavy nuclei, α decays offer reliable information on ground-state energies, half-lives, and shell effects, thereby serving as a sensitive probe of the nuclear many-body system at extreme proton numbers. The study of α decay can be traced back to the discovery of natural radioactivity by Becquerel in 1896 [1], which opened a new era in the study of unstable nuclei. Shortly thereafter, Rutherford identified α particles as helium nuclei and established α decays as a universal decay mode for heavy elements [2, 3]. The empirical Geiger–Nuttall law further revealed a strikingly simple linear relationship between the logarithm of the half-life and the inverse square root of the decay energy [4], highlighting the exponential sensitivity of the decay rates to α -decay energies [5–8]. A microscopic explanation was later provided by Gamow’s as well as Gurney and Condon’s quantum tunneling theory, which described α decay as the penetration of an α cluster through

the potential barrier [9–11]. These developments established α decay as the earliest and one of the most profound demonstrations of quantum mechanics on a macroscopic scale. Over the decades, α decay has been studied using a wide range of theoretical approaches. Early models such as the liquid-drop description captured gross features but failed to reproduce detailed systematics [12–16]. Shell-model calculations revealed the important role of closed shells and pairing correlations in stabilizing nuclei [17]. Cluster models emphasized the preformation of the α particle inside the parent nucleus [18]. More sophisticated approaches, including generalized liquid-drop models [19, 20], fission-like models [21, 22], microscopic treatments based on shell model and self-consistently density functional theory such as the relativistic Hartree-Bogoliubov theory [23, 24] have progressively improved the description of nuclear decays.

In current theoretical frameworks, α decay is usually treated as a two-step process: (i) the formation of an α cluster inside the parent nucleus, and (ii) the quantum tunneling through the potential barrier. The preformation factor P_α characterizes the likelihood of forming an α cluster in the parent nucleus. Once formed, the cluster tunnels through a barrier consisting of nuclear, Coulomb,

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[†] E-mail: wanniu@scut.edu.cn

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and centrifugal contributions. The tunneling factor is evaluated via the semiclassical Wentzel-Kramers-Brillouin (WKB) approximation [25–27], while the decay width Γ as well as the half-life $T_{1/2}$ can be calculated using the two-potential approach (TPA) introduced by Gurvitz in 1987 [28, 29]. On the other hand, since the direct calculation on P_α involves complicated quantum many-body problems, it is usually assumed to be a constant value or extracted by reproducing the experimental data. There have been several studies to calculate the preformation factor from systematics or semi-microscopic models [30–35]. By combining the shell model and cluster model, the α -cluster preformation factor in the parent nucleus ^{212}Po with a doubly-magic core ^{208}Pb has been microscopically calculated lying in the range of 0.23–0.30 [36, 37], as constrained by the reproduction of the experimental α -decay width. Under reasonable approximations for the many-body system, the quarteting wave-function approach is proposed to microscopically calculate the preformation factor in heavy and superheavy nuclei [38, 39]. From the interactions among the valence nucleons in the parent nucleus, the cluster-formation model (CFM) is proposed to calculate the preformation factor of even-even parent nuclei with experimental separation energies [40]. Later, the CFM is further developed to study the odd- A and odd-odd nuclei by considering the effects from the odd-valence nucleons and the blocking [41–44]. This model has been used to systematically calculate the α -cluster preformation factor for heavy and superheavy nuclei, where the consistent value 0.221 for ^{212}Po is also obtained.

Besides the preformation factor P_α , the accuracy of the calculated α -decay half-lives $T_{1/2}$ also depends critically on the α -decay energies Q_α which can be determined by direct experimental data or available nuclear masses. Since the decay half-life depends exponentially on the decay energy, small uncertainty in nuclear masses can propagate into the predicted half-lives by orders of magnitude differences. Several global nuclear mass models, such as the Weizsäcker-Skyrme model (WS4) [45–49], the finite-range droplet model (FRDM) [50–52], the Duflo–Zuker model (DZ10) [53–55], and the Koura Tachibana–Uno–Yamada model (KTUY) [56], are widely used to reproduce the experimental masses given in the Atomic Mass evaluation (AME2020) [57–59] throughout the nuclide chart and predict theoretical masses for possible nuclides. These models incorporate liquid-drop contributions, shell corrections, pairing, and diffuseness effects, and can reproduce known α -decay energies consistently. Recently, the microscopic density functional theory approaches, particularly the relativistic mean-field with PC-PK1 parametrization, have achieved great success in describing nuclear masses across the nuclear chart [60–67]. All these models can reproduce the experimental data very well, however, their extrapolations into re-

gions lacking experimental data can diverge significantly, especially in the superheavy region.

The superheavy element region is of particular interest because it involves the existence limit of protons in atomic nuclei. Since the landmark synthesis of the element with $Z=118$ in 2007 using ^{48}Ca -induced reactions [68, 69], much attention has been paid to the next candidates, namely the elements with $Z=119$ and $Z=120$ [70–74]. Experimental progress has been rapid by selecting ^{50}Ti , ^{51}V or ^{54}Cr as projectiles [75–78] and the direct identification of new elements relies almost exclusively on the observations of relevant α -decay chains. Considering the particular role of α decay in confirming the discovery of new elements [79], it is necessary to provide more accurate theoretical predictions of the α -decay half-lives for experimental searches. However, there exist large uncertainties for α decays of superheavy elements from the unknown decay energies Q_α as well as the cluster preformation factor P_α . Hence, to perform more consistent theoretical calculations for α decays with known superheavy nuclei and more reliable predictions for unknown ones, it is necessary to study the effects of the predicted α -decay energies and the deduced α -cluster preformation factors from different nuclear mass models.

In this work, we use the density-dependent cluster model (DDCM) to study α decays throughout the calculations [80–83], which has been widely used for α decays by inserting the decay energy Q_α and the preformation factor P_α . The latter is usually obtained from two types of CFM by combining the relevant theoretical nuclear masses [42, 44]. Under different nuclear mass models to estimate the decay energies Q_α , the preformation factors P_α , and the decay half-lives $T_{1/2}$, we perform a systematic evaluation of α -decay properties for superheavy nuclei. Using nuclear masses from WS4, FRDM, DZ10, and KTUY as inputs, we calculate α -decay systematics for known and extrapolated superheavy elements, while the data in AME2020 are used as benchmark. We assess their consistency, quantify correlations among preformation factors, and explore the stability trends toward superheavy isotopes with more neutrons. This study provides both the predictive power of these mass models and practical predictions for ongoing experimental searches for new elements. This paper is organized as follows. Section 2 details the theoretical formalism of DDCM and CFM. In Section 3, the corresponding α -decay half-lives along with correlations among theoretical predictions for various nuclear mass models are presented and discussed. A summary is given in Section 4.

II. THEORETICAL FORMALISM

A. Density-dependent cluster model

Within the density-dependent cluster model (DDCM),

the total potential $V(R)$ between the α cluster and the daughter nucleus is the sum of the nuclear potential $V_N(R)$, the Coulomb potential $V_C(R)$, and the centrifugal potential $V_\ell(R)$:

$$V(R) = V_N(R) + V_C(R) + V_\ell(R), \quad (1)$$

where R represents the distance between the centers of mass of the α cluster and the daughter nucleus. The nuclear potential $V_N(R)$ and the Coulomb potential $V_C(R)$ are calculated using the double-folding model [82–85]:

$$V_N(R) = \lambda \int \int \rho_1(\mathbf{r}_1) \rho_2(\mathbf{r}_2) g_{00}(s, \epsilon) d\mathbf{r}_1 d\mathbf{r}_2, \quad (2)$$

$$V_C(R) = \int \int \rho_{1p}(\mathbf{r}_1) \rho_{2p}(\mathbf{r}_2) \frac{e^2}{|\mathbf{s}|} d\mathbf{r}_1 d\mathbf{r}_2, \quad (3)$$

where ρ_1 and ρ_2 represent the mass density distributions of the α cluster and the daughter nucleus, while ρ_{1p} and ρ_{2p} denote their proton density distributions, respectively. The density distribution of the α cluster is usually taken as the Gaussian form [82, 84, 85]

$$\rho_1(\mathbf{r}_1) = 0.4229 \exp(-0.7024 r_1^2), \quad (4)$$

and the density distribution of the daughter nucleus is taken as the standard two-parameter Fermi form

$$\rho_2(r_2) = \frac{\rho_{20}}{1 + \exp\left(\frac{r_2 - c}{a}\right)}, \quad (5)$$

where the half-radius $c = 1.07 A_2^{1/3}$ fm and the diffuseness parameter $a = 0.54$ fm are taken from textbooks. The value of ρ_{20} is determined by integrating the density distribution equivalent to the mass number A_2 of the daughter nucleus. The quantity $g_{00}(s, \epsilon)$ in Eq. (2) denotes the M3Y-Reid effective nucleon-nucleon interaction

$$g_{00}(s, \epsilon) = 7999 \frac{e^{-4s}}{4s} - 2134 \frac{e^{-2.5s}}{2.5s} - 276(1 - 0.005\epsilon)\delta(s), \quad (6)$$

where $\epsilon = Q_\alpha/A_1$ with $A_1 = 4$ being the mass number of α cluster. The centrifugal potential $V_\ell(R)$ is defined as

$$V_\ell(R) = \frac{\hbar^2 \ell(\ell+1)}{2\mu R^2}, \quad (7)$$

where ℓ denotes the orbital angular momentum carried by the α cluster and μ represents the reduced mass of the two-body system. For favored α decays with no changes of angular momentum and parity between the parent and daughter nuclei, $\ell = 0$ is adopted, which applies to the

majority of superheavy nuclei considered in this work. For unfavored decays, the minimum allowed ℓ consistent with angular momentum and parity conservation is used. The nuclear potential strength λ in Eq. (2) is determined to satisfy the Bohr-Sommerfeld quantization condition [86–88]

$$\int_{R_1}^{R_2} \sqrt{\frac{2\mu}{\hbar^2} |Q_\alpha - V(R)|} dr = (G - \ell + 1) \frac{\pi}{2}, \quad (8)$$

where R_1 and R_2 are the first two classical turning points determined by $V(R) = Q_\alpha$. The global quantum number G is usually taken as [83, 89]

$$G = \begin{cases} 22 & N > 126, \\ 20 & 82 < N \leq 126, \\ 18 & N \leq 82. \end{cases} \quad (9)$$

The α -cluster penetration factor through the potential barrier is calculated using the well-known WKB approximation [25–27]

$$P_T = \exp \left[-2 \int_{R_2}^{R_3} \sqrt{\frac{2\mu}{\hbar^2} |Q - V(R)|} dR \right], \quad (10)$$

where R_3 is the third classical turning point. Then the α -decay width Γ can be obtained using the TPA

$$\Gamma = P_\alpha F \frac{\hbar^2}{4\mu} P_T, \quad (11)$$

where P_α is the α -cluster preformation factor in the parent nucleus, which will be discussed in detail in the next section. The normalization factor F is defined by the TPA and can be calculated by [28, 29]

$$F \int_{R_1}^{R_2} \frac{1}{2 \sqrt{\frac{2\mu}{\hbar^2} |Q - V(R)|}} dR = 1. \quad (12)$$

Finally, the α -decay half-life can be calculated using the famous relationship $T_{\text{cal}} = \hbar \ln 2 / \Gamma$. By fixing the α -cluster preformation factor as unity in the calculation of the decay half-life T_{ext} , it can be also extracted from the experimental decay half-life T_{expt} by

$$P_\alpha^{\text{ext}} = \frac{T_{\text{ext}}}{T_{\text{expt}}}. \quad (13)$$

Hence, we can systematically compare the deduced preformation factors from theoretical mass models and the extracted ones from experimental decay half-lives.

B. Cluster formation model

Within the DDCM framework, α decay is treated as the preformed α cluster inside the parent nucleus penetrating the potential barrier. Hence, the preformation factor P_α is necessary for the α -decay process. Total microscopic calculations are generally impractical for heavy and superheavy nuclei because of the complexity of many-body correlations. Here we use the recently developed cluster formation model (CFM) to provide systematic evaluations.

The CFM is originally proposed to calculate α -cluster preformation factor for even-even parent nuclei with the ratio of the effective formation energy $E_{f\alpha}$ and the relevant total energy E of the α cluster

$$P_\alpha^{\text{CFM}} = \frac{E_{f\alpha}}{E}, \quad (14)$$

where the above two energies can be calculated by using the relevant proton-, neutron-, and α -cluster separation energies. Shortly after, this approach is further developed to study the α decays for odd- A and odd-odd nuclei by including the odd-nucleon effects with two separate treatments. By considering the blocking effects of unpaired nucleons on the separation energies, Deng *et al.* proposed to use the relevant two-nucleon separation energies instead of the single-nucleon energies for the effective formation energy $E_{f\alpha}$ [42]. Here we call this CFM framework as Type I. The corresponding formulas for $E_{f\alpha}$ and E can be denoted as

$$E_{f\alpha} = \begin{cases} 2S_p(N, Z) + 2S_n(N, Z) - S_\alpha(N, Z), & e-e \\ 2S_p(N, Z) + S_{2n}(N, Z) - S_\alpha(N, Z), & e-o \\ S_{2p}(N, Z) + 2S_n(N, Z) - S_\alpha(N, Z), & o-e \\ S_{2p}(N, Z) + S_{2n}(N, Z) - S_\alpha(N, Z), & o-o \end{cases} \quad (15)$$

and

$$E = S_\alpha(N, Z), \quad (16)$$

where S_p , S_n , S_{2p} , S_{2n} , and S_α represent the separation energies of a single proton, single neutron, two protons, two neutrons, and an α cluster, respectively. They can be calculated from the relevant binding energies of the corresponding nuclei by

$$\begin{cases} S_p(N, Z) = B(N, Z) - B(N, Z-1), \\ S_n(N, Z) = B(N, Z) - B(N-1, Z), \\ S_{2p}(N, Z) = B(N, Z) - B(N, Z-2), \\ S_{2n}(N, Z) = B(N, Z) - B(N-2, Z), \\ S_\alpha(N, Z) = B(N, Z) - B(N-2, Z-2) \end{cases} \quad (17)$$

On the other hand, by considering the blocking effects of the unpaired nucleons on the residual nuclei, Ahmed *et al.* [44] introduced the odd-nucleon effects into both the separation energies and the corresponding total energies E . We call this CFM framework as Type II and the explicit expressions can be denoted as

$$E_{f\alpha} = \begin{cases} 2S_p(N, Z) + 2S_n(N, Z) - S_\alpha(N, Z), & e-e \\ 2S_p(N, Z-1) + 2S_n(N, Z-1) \\ \quad - S_\alpha(N, Z-1), & e-o \\ 2S_p(N-1, Z) + 2S_n(N-1, Z) \\ \quad - S_\alpha(N-1, Z), & o-e \\ 2S_p(N-1, Z-1) + 2S_n(N-1, Z) \\ \quad - S_\alpha(N-1, Z), & o-o \end{cases} \quad (18)$$

and

$$E = \begin{cases} B(N, Z) - B(N-2, Z-2), & e-e \\ B(N, Z) - B(N-2, Z-3), & e-o \\ B(N, Z) - B(N-3, Z-2), & o-e \\ B(N, Z) - B(N-3, Z-3), & o-o \end{cases} \quad (19)$$

From above formulism, we can see that the α -cluster preformation factor from both the two types of CFM are closely related to the relevant binding energies. With the binding energies from nuclear mass models, both the α -decay energies Q_α and the preformation factors P_α can be deduced. Under the double-folding potential within the DDCM, the corresponding α -decay half-lives can be calculated. Hence, different mass predictions with various nuclear mass models can result in diverse α -decay energies and α -cluster preformation factors, as well as the corresponding α -decay half-lives, especially in superheavy element region with relatively large uncertainty of theoretically predicted nuclear masses.

III. RESULTS AND DISCUSSION

Within the DDCM and CFM framework, systematic evaluations of α -decay properties can be achieved across known and extrapolated superheavy nuclei by combining nuclear mass models. Using theoretical nuclear masses from WS4, FRDM, DZ10, and KTUY, we computed α -decay energies Q_α , α -cluster preformation factors P_α , and α -decay half-lives $T_{1/2}$. The decay energies deduced from the experimental nuclear masses given in AME2020 are employed as benchmark calculation.

Since the decay energy is crucial for α decays, we first check the consistency between the theoretical decay energies and the experimental data deduced from AME2020 for all the known α -decay superheavy parent

nuclei from ^{228}Th to ^{295}Og . In Table I, we show the corresponding root-mean-square error (RMSE) for the α -decay energies, which is calculated by

$$\text{RMSE} = \sqrt{\frac{1}{N} \sum_{i=1}^N (Q_{\alpha,i}^{\text{theo}} - Q_{\alpha,i}^{\text{expt}})^2}, \quad (20)$$

where N denotes the number of the involved experimental α -decay energies. From Table I we can see that in general, all the nuclear mass models can well reproduce the deduced experimental decay energies with very small RMSEs. In detail, the WS4 and FRDM results achieve the smallest errors, typically below 0.4 MeV, whereas the DZ10 model displays the largest error and the KTUY model lies between them. The large differences for the α -decay energies from the DZ10 model could result in relatively large discrepancies for the corresponding theoretical α -decay half-lives.

To study the correlations for the α -decay energies among different nuclear masses, the corresponding Pearson correlation coefficient P is calculated, which is defined as

$$P = \frac{\sum (x_i - \bar{x})(y_i - \bar{y})}{\sqrt{\sum (x_i - \bar{x})^2} \sqrt{\sum (y_i - \bar{y})^2}}. \quad (21)$$

The Pearson coefficient P is widely used in statistics to measure the degree of linear dependence between two sets of variables $\{x_i\}$ and $\{y_i\}$, where $\bar{x}$ and $\bar{y}$ are their averaged values, respectively. The absolute value of the Pearson coefficient P is in the range of 0–1, where $|P| \sim 1.0$ indicates a very strong correlation and $|P| \sim 0.0$ demonstrates a very weak correlation between $\{x_i\}$ and $\{y_i\}$. In Fig. 1, we give the results of the Pearson coefficients for the decay energies among different nuclear masses. It is obvious that all the values of the Pearson coefficient between the theoretical mass models and the AME2020 data are very close to unity, indicating very good consistency for reproducing the experimental data by all the nuclear mass models. In addition, the very large Pearson coefficients among the different mass models also demonstrate the consistency to reproduce the available data. Relatively, the DZ10 model corresponds to the

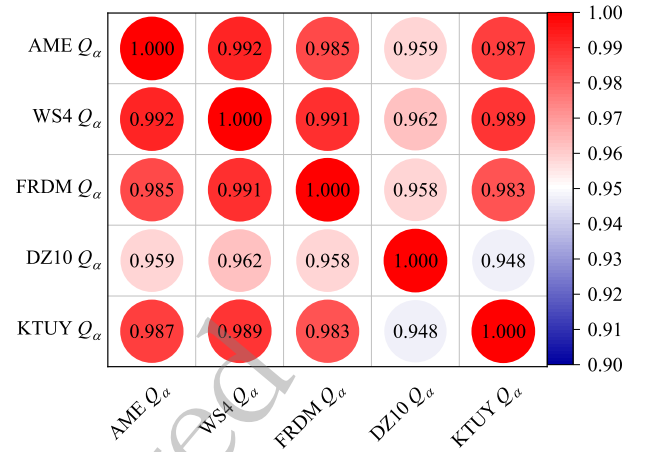


Fig. 1. (color online) Pearson correlations among different decay energies Q_α deduced from WS4, FRDM, DZ10, and KTUY mass models and the experimental masses given in AME2020.

smallest Pearson coefficients, which is accordant with the largest RMSE shown in Table I.

With these obtained α -decay energies, we can further extract the α -cluster preformation factors P_α^{ext} from the experimental decay half-lives with Eq. (13). Using the nuclear masses from different theoretical models, we can also calculate the preformation factors P_α^{CFM} with Eq. (14) under the CFM. In Fig. 2, we show the Pearson coefficients among the preformation factors from both the experimental extractions and the CFM calculations. The preformation factors within CFM in Fig. 2(a) are calculated by the Type I Eqs. (15)–(16) and those in Fig. 2(b) are calculated by the Type II Eqs. (18)–(19).

In the upper left block of Fig. 2(a) are the Pearson coefficients among the CFM-calculated preformation factors with the experimental masses from AME2020 (AME P_α^{CFM}) and with the theoretical masses from corresponding models (model P_α^{CFM}). It can be seen that the Pearson coefficients are generally very close to unity. In detail, the results with WS4, FRDM, and KTUY are all larger than 0.9, while those with DZ10 are relatively small, which is consistent with the larger decay energy deviation from the experimental data. In the lower right block of Fig. 2(a) are the Pearson coefficients among the extracted preformation factors with the deduced experimental decay energies from AME2020 (AME P_α^{ext}) and with the theoretical decay energies from corresponding mass models (model P_α^{ext}). It can be seen that these Pearson coefficients for the extracted preformation factors are generally smaller than those from the CFM calculations with nuclear mass models. This is mainly because of the exponential dependence of the decay half-life on the decay energy. Hence, the small difference in the nuclear mass and the decay energy can lead to large discrepancy in the decay half-life, resulting in large difference in the

Table 1. Root Mean Square Errors (RMSEs) of the decay energy Q_α deduced by different mass models compared with those from the experimental masses given in AME2020 for all the known α -decay superheavy parent nuclei from ^{228}Th to ^{295}Og .

	WS4	FRDM	DZ10	KTUY
RMSE	0.2455	0.3679	0.5319	0.4371

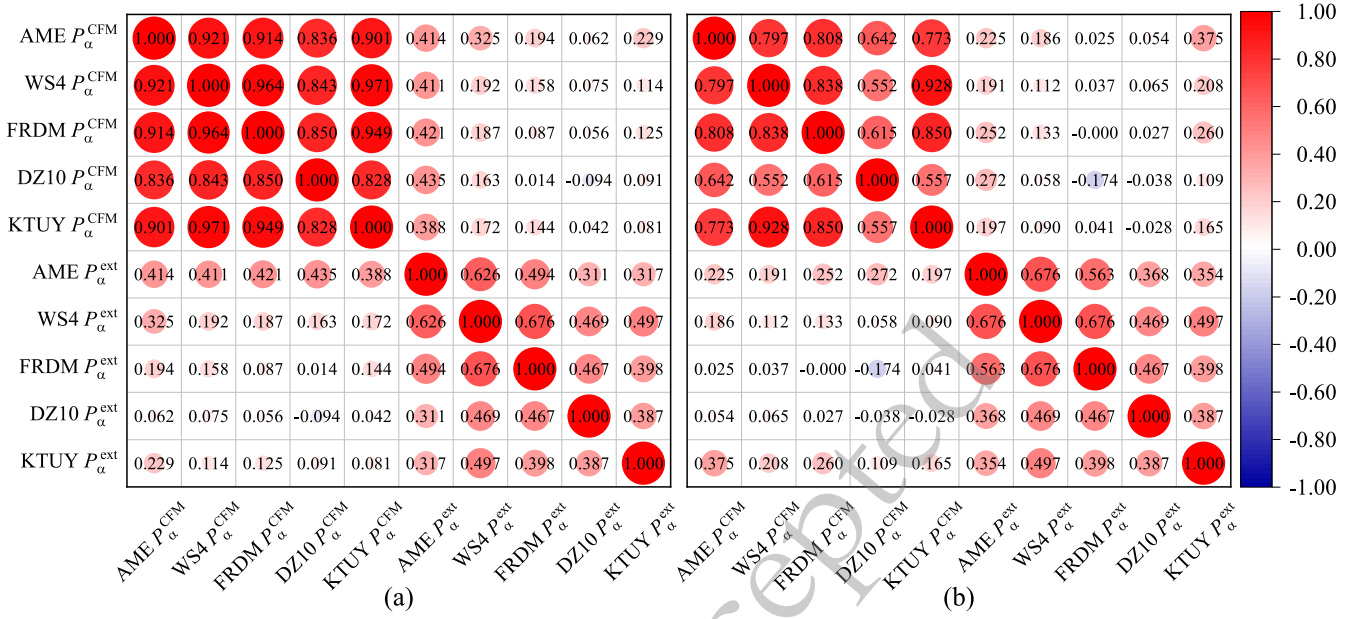


Fig. 2. (color online) Pearson correlations among the preformation factors obtained by the extractions from experimental half-lives with Eq. (13) and by the CFM calculations with Eq. (14). The CFM results in panel (a) are calculated by Type I formulas with Eqs. (15)-(16) and those in panel (b) are calculated by Type II formulas with Eqs. (18)-(19).

extracted preformation factors and small corresponding Pearson coefficients. However, the values of the Pearson coefficients are generally around 0.5, indicating consistency among these extracted preformation factors.

The results in the lower left block (conjugate in the upper right block) of Fig. 2(a) demonstrate the correlations among the preformation factors deduced from nuclear masses with CFM (P_α^{CFM}) and those extracted with the experimental decay half-lives (P_α^{ext}). The values of the Pearson coefficients are generally small. Such kind of weak correlations among them originates not only the exponential dependence on the decay energy Q_α but also the discrepancies from the theoretical nuclear masses. Considering the exponential sensitivity of the decay half-life T_{cal} to the decay energy Q_α , small discrepancies in Q_α can lead to orders-of-magnitude changes in T_{cal} . Hence, the correlation between P_α^{ext} and P_α^{CFM} will be significantly affected by the accuracy of the decay energy and the nuclear masses. Besides, the differences of the theoretical nuclear masses will also contribute to these weak correlations among the mass models via the decay energy Q_α and the CFM-calculated preformation factors. However, when both experimental masses and experimental decay half-lives are used, the extracted preformation factors ($AME P_\alpha^{ext}$) show obvious correlations with those from nuclear masses under CFM (model P_α^{CFM}), as shown by the Pearson coefficients around 0.4 in the first line of this block. This is because the mass models can reproduce well the experimental masses as well as the deduced experimental decay energies. Hence, the correlations are enhanced by reducing the discrepancy from the

decay energies. It is worth noting that the extracted WS4 preformation factors ($WS4 P_\alpha^{ext}$) have a relatively large Pearson coefficient 0.325 with the factors $AME P_\alpha^{CFM}$, which indicates the very good accuracy of the WS4 mass model to reproduce the experimental masses and decay energies.

In Fig. 2(b), the Pearson coefficients for CFM calculations in the upper left block from the Type II formulas are generally a little smaller than those from the Type I formulas, while the values for extractions in the lower right block from the Type II are generally a little larger than those from the Type I. That is to say, though the Type II treatment on the odd-nucleon effect weakens the correlations among the preformation factors within different mass models, it promotes the correlations from extractions within experimental decay half-lives. However, the overall correlations shown in Fig. 2(b) agree very well with those in Fig. 2(a), indicating the very consistent inclusion of the odd-nucleon effect on the preformation factors between these two types of CFM formulas. Hence, we take the type I CFM formulas as an example to calculate the α -cluster preformation factors P_α throughout the following calculations.

With the deduced preformation factors under the CFM and corresponding decay energies, we can further calculate the α -decay half-lives within the DDCM. In Figs. 3-6, we present the theoretical α -decay energies and calculated α -decay half-lives for different mass models as well as the available experimental data for possible α -decay chains with even-even, even-odd, odd-even, and odd-odd superheavy parent nuclei, respectively. The dotted

lines denote the α -decay energies Q_α and the solid lines represent the corresponding decay half-lives $T_{1/2}$. The black lines represent the available experimental α -decay half-lives given in AME2020 with known α -decay branch ratios, while the colored lines correspond to the theoretical calculations where the decay energies and preformation factors are both deduced from the nuclear masses. Besides, all panels are plotted from large mass and proton numbers to small ones, demonstrating the possible α -decay chains. Generally, the theoretical α -decay energies in all figures decrease with the decreasing mass and proton numbers, while the tendency for the theoretical decay

half-lives is inverse. This is because the larger decay energy will enhance the probability of α -cluster penetrating the potential barrier, leading to smaller decay half-life. Besides, from Figs. 3-6 we can also see that among all models, the WS4 and FRDM masses provide the most consistent half-life predictions with the AME2020 data, while the DZ10 exhibits systematic fluctuations in the superheavy region. The higher accuracy of the former two mass models originates from their refined macroscopic-microscopic frameworks, which include updated shell and deformation corrections optimized for heavy nuclei. In contrast, while successful in medium-mass regions, the

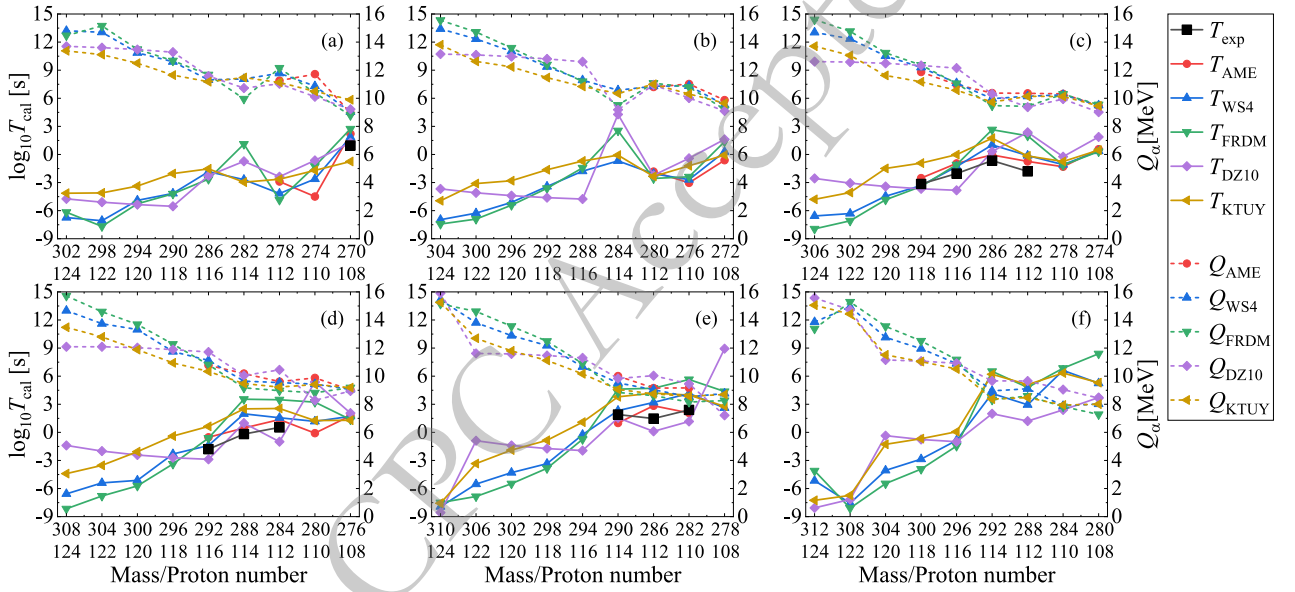


Fig. 3. (color online) Theoretical α -decay energies and calculated α -decay half-lives from different mass models as well as the available experimental data given in AME2020 for possible α -decay chains with even-even superheavy parent nuclei.

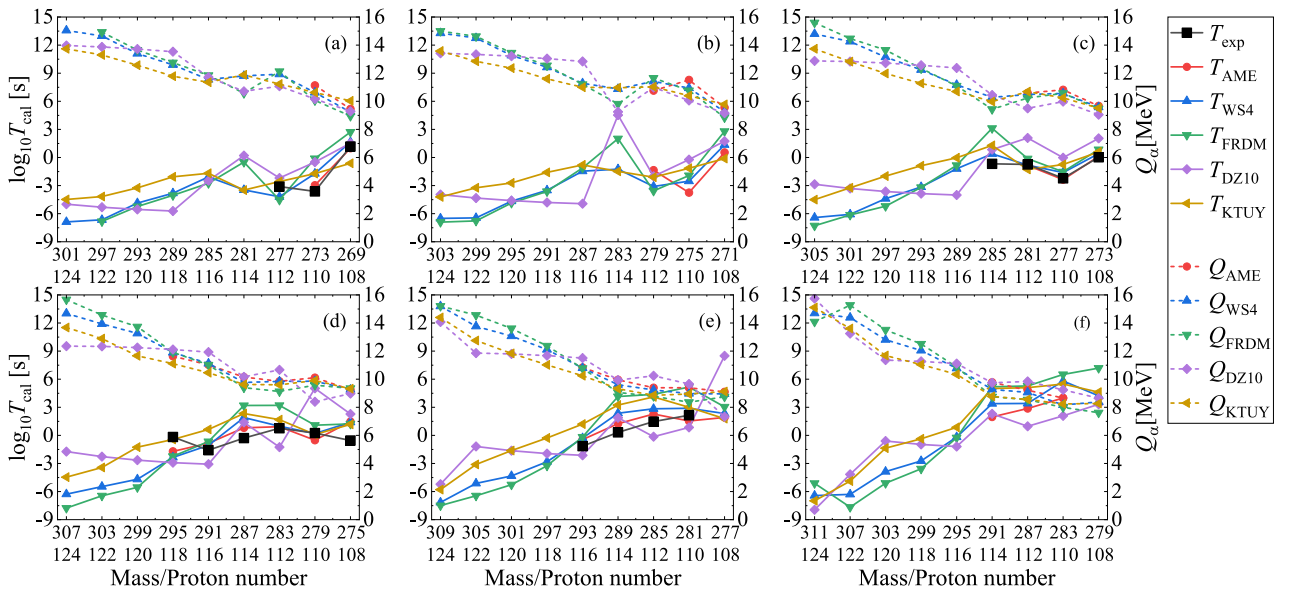


Fig. 4. (color online) Same as Fig. 3 but for even-odd superheavy parent nuclei.

DZ10 model tends to estimate the nuclear masses in the superheavy domain with large uncertainty, resulting in larger deviations in decay energies and half-life predictions. The detailed numerical values for the experimental α -decay half-lives and those calculated with the DDCM using AME2020, WS4, FRDM, DZ10, and KTUY nuclear masses are listed in Table 2 for the known nuclei in Figs. 3-6. It is shown that the theoretical calculations generally agree with the experimental data, especially the AME2020 results with a smallest RMSE of 0.8. Besides, the results with WS4 masses are more consistent with the experimental data than those with FRDM, DZ10 and KTUY masses.

Next let's concentrate on the results in Fig. 3. From each panel of Fig. 3 for even-even nuclei, it can be seen that for the superheavy parent nuclei with available deduced experimental decay energies Q_{AME} (red dashed lines), the theoretical decay energies (colored dashed lines) from different mass models agree with each other very well. However, there are discrepancies among the theoretical predictions for the unknown superheavy parent nuclei, especially for those with large mass and proton numbers. For the α -decay half-lives, all the theoretical calculations (solid lines) can well reproduce the available experimental data (black lines), with the discrepancy generally in one order of magnitude. Especially,

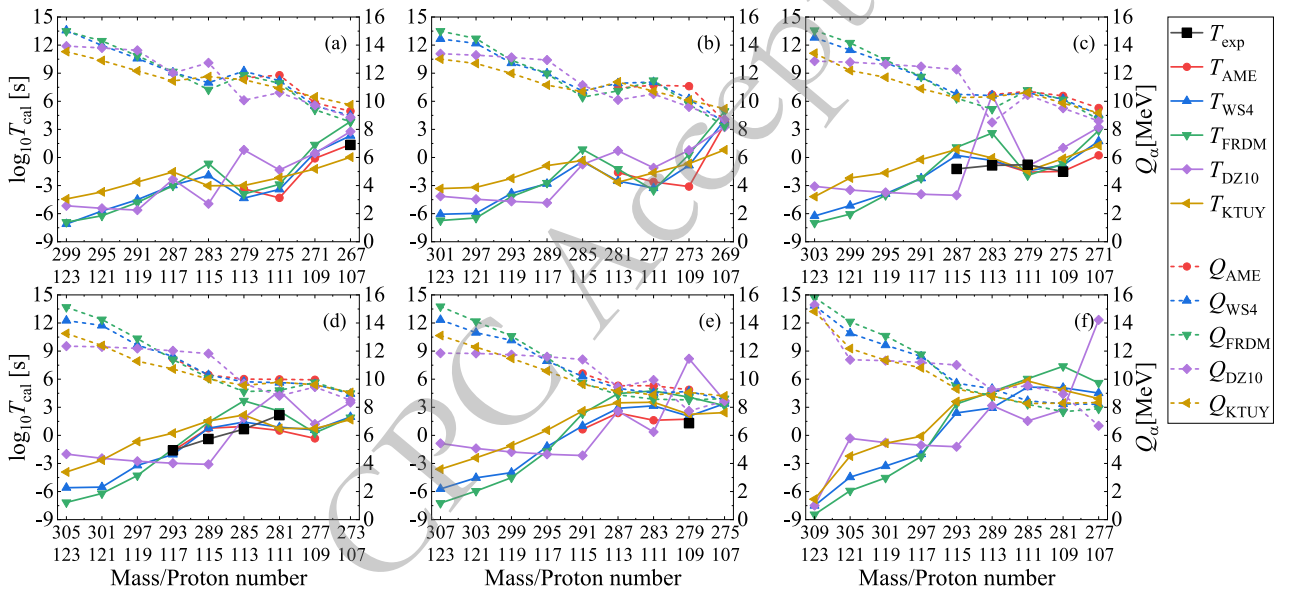


Fig. 5. (color online) Same as Fig. 3 but for odd-even superheavy parent nuclei.

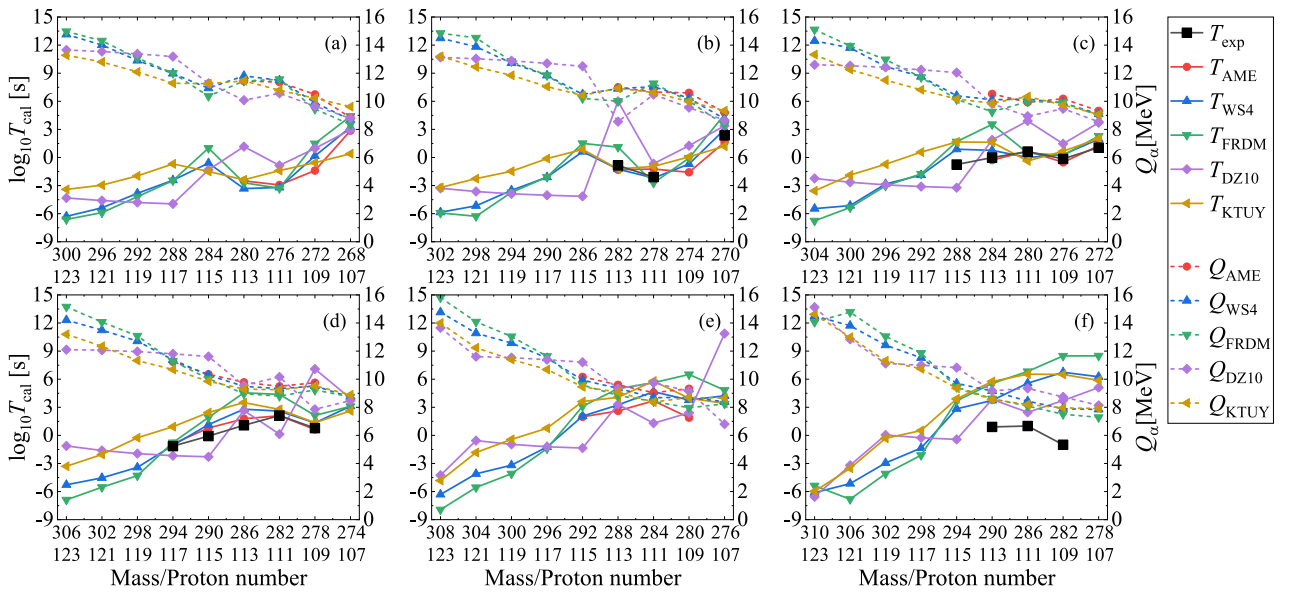


Fig. 6. (color online) Same as Fig. 3 but for odd-odd superheavy parent nuclei.

Table 2. Comparison between the experimental α -decay half-lives and the calculated values under the DDCM using AME2020, WS4, FRDM, DZ10, and KTUY nuclear masses.

A	Z	$\log_{10} T_{\text{exp}}[\text{s}]$	$\log_{10} T_{\text{AME}}[\text{s}]$	$\log_{10} T_{\text{WS4}}[\text{s}]$	$\log_{10} T_{\text{FRDM}}[\text{s}]$	$\log_{10} T_{\text{DZ10}}[\text{s}]$	$\log_{10} T_{\text{KTUY}}[\text{s}]$
270	108	0.9542	2.2033	1.8534	2.7329	1.2590	-0.7488
272	108		-0.6159	0.2882	1.3680	1.6121	-0.0658
274	108		0.5625	0.3597	0.3084	1.8747	0.4616
276	108		1.7719	1.7098	1.3653	2.0505	1.2850
274	110		-4.4981	-2.6237	-1.2527	-0.6394	-1.7032
276	110		-3.0367	-2.6888	-2.3824	-0.4053	-1.2142
278	110		-1.3154	-1.0886	-1.1539	-0.2408	-0.7480
280	110		-0.0900	1.1702	3.2231	5.3795	1.2578
282	110	2.4014	2.0740	4.1514	5.6401	1.1600	3.9835
278	112		-2.8880	-4.1681	-4.8579	-2.3797	-2.6439
280	112		-1.8288	-2.0696	-2.5568	-2.2240	-2.3406
282	112	-1.8037	-0.7425	-0.1146	1.9899	2.3708	-0.1562
284	112	0.5315	1.3605	1.5598	3.4882	-1.0032	2.5436
286	112	1.4771	2.8427	3.1899	4.7132	0.1055	4.1916
286	114	-0.6569	-0.0587	1.0237	2.6452	0.3143	1.7322
288	114	-0.1851	0.4647	1.9768	3.5379	0.9875	2.5002
290	114	1.9031	1.0021	2.3257	4.6088	1.5506	3.7946
290	116	-2.0458	-0.9539	-1.3042	-1.0991	-3.8273	-0.0216
292	116	-1.7959	-0.5264	-1.4071	-0.6114	-2.8858	0.6309
294	118	-3.1549	-2.5281	-3.3218	-3.4187	-3.6680	-0.9201
269	108	1.1761	0.9896	1.6533	2.7232	1.5322	-0.6284
271	108		0.5424	1.3585	2.7936	1.7417	-0.1211
273	108	0.0253	0.1192	0.4103	0.8539	2.0474	0.5906
275	108	-0.5528	1.4691	1.4030	1.2235	2.2748	1.1768
277	108		1.9092	2.3288	3.0438	8.4894	1.7878
273	110	-3.6198	-2.9717	-1.7228	-0.1080	-0.4918	-1.7254
275	110		-3.7542	-2.5221	-1.9273	-0.2207	-1.1495
277	110	-2.2218	-2.3708	-1.6182	-1.5318	-0.0254	-0.7579
279	110	0.2430	-0.5155	0.1897	1.0890	5.0247	0.0969
281	110	2.1461	1.5324	2.8942	5.1638	0.8471	2.9974
283	110		3.9934	5.8096	6.5100	2.0892	5.4835
277	112	-3.1024		-4.1969	-4.5360	-2.1951	-2.5659
279	112		-1.3501	-3.0915	-3.5669	-2.0067	-2.1088
281	112	-0.7447	-0.8491	-0.8134	-0.1266	2.0843	-1.2802
283	112	0.7636	0.9265	0.9830	3.2037	-1.2542	1.6729
285	112	1.4771	2.2832	2.8369	4.3812	-0.1431	4.1085
287	112		2.8689	3.4086	5.2776	0.9645	5.0929
285	114	-0.6778		0.3818	3.1222	0.8405	1.2954
287	114	-0.2924	0.8063	1.9083	3.1927	1.3894	2.3371
289	114	0.3222	1.2447	2.3606	4.1463	1.8832	3.2366
291	114		1.9705	3.3823	5.1876	2.2958	4.9979
291	116	-1.5850	-0.8377	-1.0636	-0.6566	-3.0753	0.6363
293	116	-1.1549	-0.4481	-0.2928	-0.1759	-2.1325	1.2085
295	118	-0.1675	-1.7111	-2.3503	-2.2362	-2.9102	-0.4436

Continued on next page

Table 2-continued from previous page

A	Z	$\log_{10} T_{\text{exp}}[\text{s}]$	$\log_{10} T_{\text{AME}}[\text{s}]$	$\log_{10} T_{\text{WS4}}[\text{s}]$	$\log_{10} T_{\text{FRDM}}[\text{s}]$	$\log_{10} T_{\text{DZ10}}[\text{s}]$	$\log_{10} T_{\text{KTUY}}[\text{s}]$
267	107	1.3424	1.3920	2.3185	3.8407	2.7752	0.0335
269	107		3.5596	3.9890	4.9322	3.2414	0.8114
271	107		0.2222	1.7274	2.8833	3.2121	1.2822
271	109		-0.0946	0.5230	1.3595	0.4446	-1.2262
273	109		-3.1088	-0.8499	0.3244	0.7714	-0.6580
275	109	-1.5086	-1.4891	-0.8299	-0.6856	1.0251	-0.1097
277	109		-0.3207	0.5852	0.2643	1.2018	0.7281
279	111	-0.7696	-1.5661	-1.4275	-1.9151	-0.9333	-1.4368
281	111	2.1648	0.5127	0.8683	2.5834	4.7339	0.7403
283	111		1.6086	3.1470	4.7474	0.3682	3.5424
279	113		-3.3594	-4.3395	-3.8980	0.8057	-2.9949
281	113		-1.5979	-2.5209	-1.2381	0.7137	-2.6460
283	113	-0.8539	-0.3812	-0.3147	2.5972	6.6919	-0.0498
285	113	0.6628	0.9413	1.4239	3.6872	1.8626	2.1744
287	113		2.3858	2.8949	4.4630	2.5884	3.4723
287	115	-1.2218		0.2158	1.0875	-4.0421	0.8652
289	115	-0.3872	0.7644	0.7669	1.4003	-3.0994	1.5415
291	115		0.6461	0.9994	2.3222	-2.1555	2.6052
293	117	-1.6021	-1.6441	-2.0396	-1.4658	-2.9758	0.2322
268	107		2.8578	3.2398	4.4050	2.9598	0.3997
270	107	2.3579	1.6387	2.9934	4.5878	3.3850	1.2087
272	107	1.0531	1.2425	1.9570	2.2745	3.7146	2.0226
272	109		-1.3962	0.1623	1.4835	0.9691	-0.5563
274	109		-1.5835	-0.6992	-0.0690	1.2581	0.0651
276	109	-0.1549	-0.5155	0.3046	0.2320	1.4631	0.6707
278	109	0.7782	0.6384	1.3282	2.1319	7.0820	1.3559
280	109		1.8950	3.7765	6.5020	2.4097	3.9677
282	109	-1.0000		6.7376	8.4851	3.7226	6.5040
276	111		-2.9333	-3.2062	-3.2868	-0.8513	-1.3953
278	111	-2.0969	-1.2334	-2.1979	-2.7220	-0.6644	-0.9623
280	111	0.6335	0.6725	-0.0036	0.4767	3.9036	-0.3641
282	111	2.1139	2.0946	2.5922	4.3775	0.1325	2.7752
284	111		3.6910	4.6629	5.6215	1.3015	5.7912
286	111	1.0000		5.5703	6.8245	2.4572	6.5413
280	113		-2.4996	-3.3176	-2.6973	1.1593	-2.3916
282	113	-0.8539	-1.2978	-1.2468	1.1099	6.0008	-1.2248
284	113	-0.0132	-0.2838	0.7654	3.5324	1.8742	1.6284
286	113	1.0792	1.7789	2.7992	4.5533	2.6541	3.5013
288	113		2.6155	3.2479	4.9915	3.3144	4.0637
290	113	0.9031		3.7270	5.5409	3.8462	5.7562
288	115	-0.7520		0.8855	1.6860	-3.2236	1.6512
290	115	-0.0757	0.7825	1.1555	1.9115	-2.2890	2.4377
292	115		1.9733	2.0576	3.1115	-1.3562	3.6144
294	117	-1.1549	-0.8999	-1.0697	-0.7599	-2.1668	0.9242

from the comparison between the solid red and black lines, we can find that within the decay energies and preformation factors both deduced from the experimental nuclear masses (AME2020), the calculated decay half-lives by the DDCM are in good agreement with the experimental data, indicating the good prediction accuracy of the DDCM. For the superheavy parent nuclei with smaller mass and proton numbers, due to the less uncertainty of the decay energies among different nuclear mass models, their predicted decay half-lives agree with each other very well. However, as shown from the starting points for the much heavier parent nuclei in each panel, the theoretical predicted decay energies are divergent from different mass models, resulting in large differences among these half-life predictions. Besides, for the isotopes from panel (a) to panel (f) in Fig. 3, such as those with the proton number $Z = 108$, we can see that the theoretical decay energies Q_α will decrease if the superheavy parent nuclei have more neutrons. As a result, the corresponding predicted decay half-lives $T_{1/2}$ are also increased for heavier isotopes. For example, the α -decay half-life of $^{270}_{108}\text{Hs}$ is around the order of 10^2 s, while it is as large as 10^8 s for $^{280}_{108}\text{Hs}$. Though there are differences among the predicted decay half-lives, however, all the mass models demonstrate that the superheavy nuclei tend to become more stable with more neutrons compared with the present measured isotopes, which is consistent with the predicted “island of stability” near $N \approx 184$. These results provide quantitative support for ongoing experimental efforts to synthesize superheavy isotopes with more neutrons of elements with $Z=119-120$ using ^{50}Ti - or ^{54}Cr -induced reactions.

The tendency of the results in Figs. 4-6 for even-odd, odd-even, and odd-odd superheavy parent nuclei are very similar to those in Fig. 3 for the even-even nuclei. For the superheavy nuclei with decay energies available, all the theoretical calculations can reproduce the experimental data very well. The predicted decay energies for the superheavy nuclei with small mass and proton numbers are consistent with each other among different mass models, while there are large discrepancies for those with large mass and proton numbers. As a result, the corresponding decay half-lives for the superheavy parent nuclei with available experimental data and with small mass and proton numbers agree with each other well, generally within one order of magnitude discrepancy, while the differences among different mass models are large for the superheavy nuclei with large mass and proton numbers. Hence, the present theoretical mass models can predict relatively more consistent α -decay energies for unknown isotopes than for unknown superheavy elements. In the future, the α -decay properties of unknown isotopes can be employed to check the extrapolation reliability of different nuclear mass models as well as for further predictions of unknown superheavy elements.

IV. SUMMARY

Within the density-dependent cluster model (DDCM), we systematically study the effects of nuclear masses on α -decay properties of superheavy nuclei for different mass models, namely WS4, FRDM, DZ10, and KTUY. The experimental masses given in AME2020 are used for benchmark calculation. The consistency of α -decay energies between different nuclear mass models and the experimental data are checked at first for all the known superheavy parent nuclei from ^{228}Th to ^{295}Og . It is found that the WS4 and FRDM results achieve smaller errors, while the DZ10 model displays a relatively large discrepancy.

Under the framework of cluster formation model (CFM) with two types of formulas to include the odd-nucleon effect, the α -cluster preformation factors for the superheavy parent nuclei are calculated with the experimental masses and the theoretical masses from different mass models. With the deduced experimental decay energies, the corresponding preformation factors are also extracted from the experimental decay half-lives. The correlations among the CFM calculations and the half-life extractions are systematically studied by calculating the corresponding Pearson correlation coefficients. The results of the Pearson coefficients indicate that the CFM preformation factors are consistent with each other very well among different mass models. The extracted preformation factors also agree with each other but with less correlations among them. This is because of the exponential dependence of the decay half-life on the decay energy. Hence, the small differences in the decay energy from different nuclear masses will result in large discrepancies of the decay half-lives. Besides, though the preformation factors between the CFM calculations and the half-life extractions generally show weak correlations for each mass model, the WS4 data set demonstrates remarkably close correlations between them. In addition, the two types of CFM formulas lead to very similar correlations for all data sets, indicating the consistent inclusion of the odd-nucleon effect on the α -cluster preformation factor within the CFM.

Under the DDCM, the decay half-lives of possible α -decay chains with even-even, even-odd, odd-even, and odd-odd superheavy nuclei are calculated. Similar tendency for all the possible decay chains are obtained that with the decreasing mass and proton numbers, the theoretical decay energies decrease and the corresponding decay half-lives increase. In detail, for the superheavy parent nuclei with experimental data available, both the deduced experimental decay energies and half-lives can be reproduced well. Besides, the decay energies and half-lives for the unknown superheavy isotopes with small mass and proton numbers can be predicted with smaller uncertainty, generally with one order of magnitude differences for the half-lives. However, there are larger dis-

crepancies for the decay energies and half-lives for the unknown superheavy elements with large mass and proton numbers. Besides, it is found that the theoretical decay energies Q_α for superheavy isotopes will decrease with the increasing neutron number, leading to the in-

creasing predicted decay half-lives. Hence, the superheavy nuclei tend to become more stable with more neutrons compared with the present available isotopes and the island of superheavy nuclei could be searched by synthesizing the superheavy isotopes with more neutrons.

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