

Radial basis function approach in nuclear mass predictions

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The radial basis function (RBF) approach is applied in predicting nuclear masses for eight widely used nuclear mass models, ranging from macroscopic-microscopic to microscopic types. A significantly improved accuracy in computing nuclear masses is obtained, and the corresponding rms deviations with respect to the known masses are reduced by up to 78%. Moreover, strong correlations are found between a target nucleus and the reference nuclei within about three units in distance; these play critical roles in improving nuclear mass predictions. Based on the latest Weizsäcker-Skyrme mass model, the RBF approach can achieve an accuracy comparable with the extrapolation method used in atomic mass evaluation. In addition, the necessity of new high-precision experimental data to improve the mass predictions with the RBF approach is emphasized as well.

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I. INTRODUCTION

Nuclear mass plays an important role not only in studying our knowledge of nuclear structure [1] but also in understanding the origin of elements in the universe [2,3]. With the construction and upgrade of radioactive ion beam facilities, measurements of nuclear masses have made great progress in recent years. During the past decade, hundreds of nuclear masses were measured for the first time or with higher precisions [4].

The astrophysical calculations involve thousands of nuclei far from the β -stability line. However, most of these nuclei are still beyond experimental reach. One could use local mass relations such as the Garvey-Kelson (GK) relations [5,6] and the residual proton-neutron interactions [7–10] to predict unknown masses. However, intrinsic error grows rapidly when local mass relations are used to predict nuclear masses in an iterative way [11,12]. Therefore, theoretical predictions for nuclear masses are inevitable to perform astrophysical calculations. The early theoretical studies of nuclear masses are mainly macroscopic models, such as the famous Weizsäcker mass formula [13]. It is known that this kind of mass model neglects microscopic effects, and hence it shows systematic deviations for nuclei near the shell closure or those with large deformations. In order to better describe nuclear ground-state properties, the macroscopic-microscopic and microscopic theoretical models have been developed for mass predictions.

By including the microscopic correction energy to the macroscopic mass formula, the macroscopic-microscopic mass model can well take into account the important microscopic corrections. During the past decades, a number of

macroscopic-microscopic mass models have been developed; these include the finite-range droplet model (FRDM) [14], the extended Thomas-Fermi plus Strutinsky integral (ETFSI) model [15], and the Koura-Tachibana-Uno-Yamada (KTUY) model [16]. These macroscopic-microscopic mass models have similar accuracy for mass prediction and their root-mean-square (rms) deviation with respect to data in the atomic mass evaluation of 2012 (AME12) [4] is about 0.7 MeV. Guided by the Skyrme energy density functional, a semiempirical nuclear mass formula, the Weizsäcker-Skyrme (WS) model, was proposed based on the macroscopic-microscopic method [17–19]. For the latest version of the WS model (WS3) [19], the rms deviation with respect to 2353 known nuclear masses in AME12 is significantly reduced to 0.335 MeV.

On the other hand, great progress has been achieved for microscopic mass models with the rapid development of computer technology in the new century. Based on Hartree-Fock-Bogoliubov (HFB) theory with a Skyrme or Gogny force, a series of microscopic mass models have been proposed with accuracy comparable with that of traditional macroscopic-microscopic mass models [20–22]. Apart from the nonrelativistic microscopic model, the relativistic mean-field (RMF) model has also received wide attention due to its many successes in describing lots of nuclear phenomena [23–35] as well as successful applications in astrophysics [36–44]. A systematic study of the ground-state properties for all nuclei from the proton drip line to the neutron drip line with $Z, N \geq 8$ and $Z \leq 100$ was performed for such a model several years ago, and the rms deviation with respect to known masses is about 2 MeV [45]. However, it should be noted that the effective interaction of this RMF mass model was only optimized with the properties of a few selected nuclei. By carefully adjusting the effective interaction of the RMF model with the properties of more selected nuclei, the deviation can be remarkably reduced. For the 575 even-even nuclei with $8 \leq Z \leq 108$, the rms deviation with respect to known masses in the atomic mass evaluation of 2003 (AME03) is reduced to 1.24 MeV for the effective interaction PC-PK1 [46].

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Moreover, the PC-PK1 predictions well reproduce the new and accurate mass measurements from Sn to Pa [47] with an rms deviation of 0.859 MeV [48] and also successfully describe the Coulomb displacement energies between mirror nuclei [49]. In addition, inspired by the shell model, the Duflou-Zuker (DZ) mass model [50,51] has been quite successful in describing nuclear masses with an accuracy of about 0.5 MeV.

Although these theoretical models can well reproduce the experimental data, there are still large deviations among the mass predictions of different models, even in the region close to known masses. A number of investigations on the accuracy and predictive power of these nuclear mass models have been made so far in the literature (e.g., Refs. [1,45,52–54]). To further improve the accuracy of nuclear mass models, image reconstruction techniques based on the Fourier transform have been applied to the nuclear mass models, significantly reducing the rms deviation to the known masses with the CLEAN algorithm [55]. Later on, the radial basis function (RBF) approach was developed to improve the mass predictions of several theoretical models [56]. Compared with the CLEAN reconstruction, the RBF approach more effectively reduces the rms deviations with respect to the masses first appearing in AME03 [56].

To improve the mass prediction of a nucleus, thousands of nuclei with known masses are involved in the RBF approach [56]. However, do all the nuclei involved play effective roles in the improvement of mass prediction for a given nucleus? What are the key nuclei that have to be included in the RBF approach? In other words, how far away from the measured region of nuclear mass could we predict with satisfactory accuracy in the RBF approach? These questions were not addressed in previous investigations [56]. Therefore, it is interesting to investigate the mass correlations between a certain nucleus and those nuclei involved in the RBF approach and hence to evaluate the predictive power of the RBF approach.

In this work, we will carefully evaluate the predictive power of the RBF approach based on eight widely used nuclear mass models, ranging from macroscopic-microscopic to microscopic types. Special attention will be paid to the mass correlations among various nuclei. The paper is organized as follows. In Sec. II, a brief introduction to the RBF approach including numerical details is given. In Sec. III, the mass correlations are first carefully investigated, and then the predictive power of the RBF approach based on different mass models will be evaluated. Finally, a summary is presented in Sec. IV.

II. RADIAL BASIS FUNCTION APPROACH AND NUMERICAL DETAILS

The RBF approach has been widely applied in surface reconstruction and its solution is written as

$$S(x) = \sum_{i=1}^m \phi(\|x - x_i\|) \omega_i, \quad (1)$$

where x_i denotes the point from the measurement, ω_i is the weight of the center x_i , ϕ is the radial basis function, $\|x - x_i\|$

is the Euclidean norm, and m is the number of data points to be fitted. Given m samples (x_i, d_i) , one wishes to reconstruct the smooth function $S(x)$ with $S(x_i) = d_i$, i.e.,

$$\begin{pmatrix} d_1 \\ d_2 \\ \dots \\ d_m \end{pmatrix} = \begin{pmatrix} \phi_{11} & \phi_{12} & \dots & \phi_{1m} \\ \phi_{21} & \phi_{22} & \dots & \phi_{2m} \\ \dots & \dots & \dots & \dots \\ \phi_{m1} & \phi_{m2} & \dots & \phi_{mm} \end{pmatrix} \begin{pmatrix} \omega_1 \\ \omega_2 \\ \dots \\ \omega_m \end{pmatrix}, \quad (2)$$

where $\phi_{ij} = \phi(\|x_i - x_j\|)$ ($i, j = 1, \dots, m$). Then the RBF weights are determined to be

$$\begin{pmatrix} \omega_1 \\ \omega_2 \\ \dots \\ \omega_m \end{pmatrix} = \begin{pmatrix} \phi_{11} & \phi_{12} & \dots & \phi_{1m} \\ \phi_{21} & \phi_{22} & \dots & \phi_{2m} \\ \dots & \dots & \dots & \dots \\ \phi_{m1} & \phi_{m2} & \dots & \phi_{mm} \end{pmatrix}^{-1} \begin{pmatrix} d_1 \\ d_2 \\ \dots \\ d_m \end{pmatrix}. \quad (3)$$

Once the weights are obtained with the m samples (x_i, d_i) , the reconstructed function $S(x)$ can be calculated with Eq. (1) for any point x .

As in Ref. [56], the Euclidean norm is defined to be the distance between nuclei (Z_i, N_i) and (Z_j, N_j) on the nuclear chart:

$$r = \sqrt{(Z_i - Z_j)^2 + (N_i - N_j)^2}. \quad (4)$$

The basis function $\phi(r) = r$ is adopted in this work, since the mass deviation can be reconstructed relatively better with $\phi(r) = r$ than with other basis functions [56]. Then the mass difference $D(Z, N) = M_{\text{exp}}(Z, N) - M_{\text{th}}(Z, N)$ between the experimental data M_{exp} and those predicted with nuclear mass models, M_{th} , could be reconstructed with Eq. (2). Once the weights are obtained, the reconstructed function $S(Z, N)$ can be calculated with Eq. (1) for any nucleus (Z, N) . Then the revised mass for nucleus (Z, N) is given by

$$M_{\text{th}}^{\text{RBF}}(Z, N) = M_{\text{th}}(Z, N) + S(Z, N). \quad (5)$$

For training the RBF with Eq. (2), only those nuclei between the minimum distance R_{min} and maximum distance R_{max} are involved, i.e., $R_{\text{min}} \leq r \leq R_{\text{max}}$. If the reconstructed function $S(Z, N)$ for a nucleus is obtained by training the RBF including itself, i.e., $R_{\text{min}} = 0$, it is clear that $S(Z, N)$ is just $D(Z, N)$ and hence $M_{\text{th}}^{\text{RBF}}(Z, N) = M_{\text{exp}}(Z, N)$. Therefore, to test the predictive power of the RBF approach, the function $S(Z, N)$ for a known nucleus should be reconstructed with $R_{\text{min}} \geq 1$.

To evaluate the predictive power of the RBF approach, the rms deviation, i.e.,

$$\sigma_{\text{rms}} = \sqrt{\frac{1}{n} \sum_{i=1}^n (M_i^{\text{th}} - M_i^{\text{exp}})^2}, \quad (6)$$

is employed, where M_i^{th} and M_i^{exp} are the theoretical and experimental nuclear masses, respectively, and n is the number of nuclei contained in a given set. In this investigation, we only consider nuclei with $N \geq 8$ and $Z \geq 8$ and the experimental data are taken from AME12 [4], unless otherwise specified. For the theoretical mass models, we take RMF [45], HFB-21 [22], DZ10 [50], DZ31 [51], ETFSI-2 [15], KTUY [16], FRDM [14], and WS3 [56] mass models as examples, with

the rms deviation spanning from 2.2 to 0.3 MeV with respect to experimental data in AME12. For convenience, the standard deviations $\sigma_{\text{rms}}(\text{Model})$ and $\sigma_{\text{rms}}(\text{Model} + \text{RBF})$ of the nuclear mass models are denoted as $\sigma_{\text{rms0}}(\text{Model})$ and $\sigma_{\text{rmsR}}(\text{Model})$, hereafter.

III. RESULTS AND DISCUSSION

For testing the predictive power of the RBF approach, we first reconstruct the function $S(N, Z)$ for a known nucleus based on the remaining known masses in AME12 and the predictions in nuclear mass models. In other words, we take $R_{\text{min}} = 1$ and $R_{\text{max}} = 1000$ (i.e., no limits on R_{max}) for training the RBF. The corresponding results are given in Table I. It is found that the reduction of the rms deviation exceeds 25% for all mass models considered here. In particular, the largest improvement, with 78% reduction of the rms deviation, is obtained for the RMF mass model. The corresponding rms deviation is reduced from 2.2 to 0.5 MeV, which is comparable to the corresponding rms deviation in the microscopic HFB-21 mass model. Therefore, the reduction of rms deviations clearly shows that the predictive accuracy of the nuclear mass models can be significantly improved by combining the model with the RBF approach.

In the calculations of Table I, the reconstructed function $S(N, Z)$ for a nucleus is obtained by training the RBF with the remaining known masses in AME12. For better understanding the predictive power of the RBF, it is necessary to investigate the mass correlations between this nucleus and those nuclei used in training the RBF. In Fig. 1, the rms deviations σ_{rmsR} and the relative rms deviations $\sigma_{\text{rmsR}}/\sigma_{\text{rms0}}$ with respect to the known masses in AME12 are shown as a function of R , which is the distance between the selected nucleus and the nuclei used in training the RBF. For clarity, only the results of RMF, DZ31, FRDM, and WS3 mass models are shown in the figure. In fact, other mass models show similar trends and their corresponding results are almost between the results of RMF and WS3 mass models. From Fig. 1, it is clear that σ_{rmsR} generally increases as R increases, while the order of σ_{rmsR} generally remains the same as that of σ_{rms0} . For the WS3, DZ31, and FRDM mass models, the RBF approach

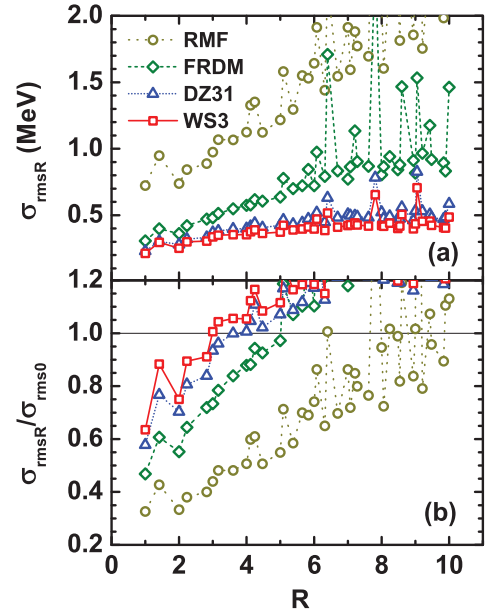


FIG. 1. (Color online) The rms deviations σ_{rmsR} and the relative rms deviations $\sigma_{\text{rmsR}}/\sigma_{\text{rms0}}$ with respect to the known masses in AME12 for different mass models. For training the RBF for a certain nucleus, only those nuclei with $r = R$ are involved.

ceases to improve the mass predictions, i.e., $\sigma_{\text{rmsR}}/\sigma_{\text{rms0}} > 1$, when the distances R are larger than 3, 4, and 5, respectively. However, the RBF approach can improve the mass prediction of the RMF model even with some nuclei around $R = 10$. This long correlation may imply that some important physics correlations are missing in this RMF mass model and hence a larger rms deviation with respect to the known masses.

Furthermore, we investigate the cumulative rms deviations by training the RBF with nuclei at distance between R_{min} and R_{max} . By fixing $R_{\text{min}} = 1$, the rms deviations σ_{rmsR} and the relative rms deviations $\sigma_{\text{rmsR}}/\sigma_{\text{rms0}}$ as a function of R_{max} are shown in Fig. 2 for different mass models. The points at $R_{\text{max}} = 0$ mean no masses are included in training the RBF, so σ_{rmsR} is just σ_{rms0} . From Fig. 2, one can see that the nuclei at distance $r = 1$ play an important role in improving the predictive accuracy of different mass models. By further including the nuclei in the range of $1 < r \leq 3$ for training the RBF, the rms deviations can be slightly reduced. However, the improvement in the predictive accuracy is almost negligible with the inclusion of nuclei at $r > 3$ for all mass models, although the nuclei at $r > 3$ can still help the RBF approach improve the mass predictions for some mass models, such as the FRDM and RMF models. This indicates that the RBF approach can well extract most mass correlations only from those nuclei with $r \leq 3$.

On the other hand, for fixed $R_{\text{max}} = 1000$ (i.e., no limits on R_{max}), the rms deviations σ_{rmsR} and the relative rms deviations $\sigma_{\text{rmsR}}/\sigma_{\text{rms0}}$ as a function of R_{min} are shown in Fig. 3 for different mass models. The points at $R_{\text{min}} = 0$ mean that the reconstructed function $S(Z, N)$ for one nucleus is obtained by training the RBF including itself, so $M_{\text{th}}^{\text{RBF}} = M_{\text{exp}}$ and hence $\sigma_{\text{rmsR}} = 0$. The points at $R_{\text{min}} = 1$ just correspond to those σ_{rmsR} values in Table I. If nuclei with $r = 1$ are excluded

TABLE I. The rms deviations in MeV between known masses in AME12 and predictions of various nuclear mass models without (the second column) and with (the third column) the RBF approach. The fourth column gives the reduction of the rms deviations after combining the model with the RBF approach.

Model	σ_{rms0}	σ_{rmsR}	Reduction (%)
RMF	2.217	0.488	78%
HFB-21	0.572	0.410	28%
DZ10	0.591	0.225	62%
DZ31	0.397	0.204	49%
ETFSI-2	0.719	0.360	50%
KTUY	0.701	0.210	70%
FRDM	0.654	0.268	59%
WS3	0.335	0.207	38%

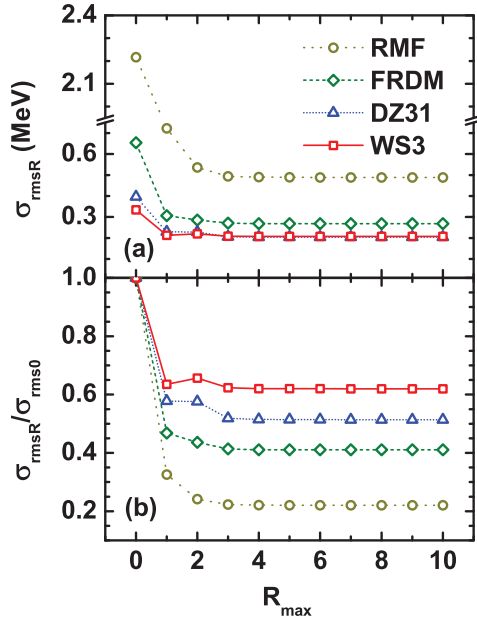


FIG. 2. (Color online) The rms deviations σ_{rmsR} and the relative rms deviations $\sigma_{\text{rmsR}}/\sigma_{\text{rms0}}$ with respect to the known masses in AME12 for different mass models as a function of R_{max} . For training the RBF, R_{min} is fixed to be $R_{\text{min}} = 1$.

from training the RBF, the rms deviation σ_{rmsR} increases for the DZ31, FRDM, and WS3 mass models, while it decreases for the RMF mass model. However, the influence on σ_{rmsR} is unremarkable, so the RBF approach can also remarkably improve the mass predictive accuracy only with nuclei of $r \geq 2$. Furthermore, if we exclude nuclei with $r \leq 2$, σ_{rmsR}

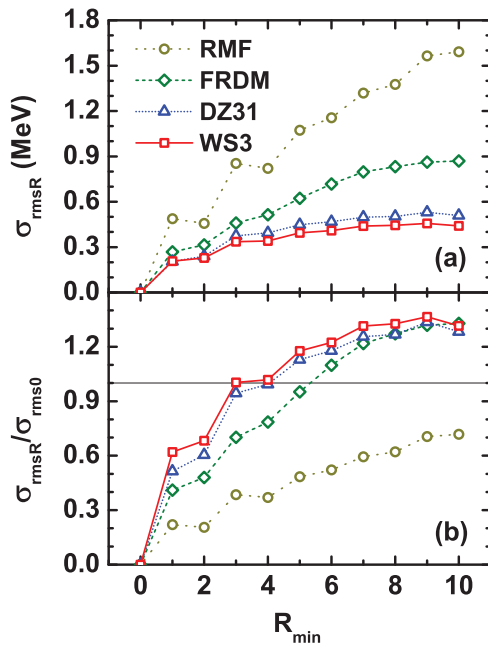


FIG. 3. (Color online) The rms deviations σ_{rmsR} and the relative rms deviations $\sigma_{\text{rmsR}}/\sigma_{\text{rms0}}$ with respect to the known masses in AME12 for different mass models as a function of R_{min} . For training the RBF, there are no limits on R_{max} .

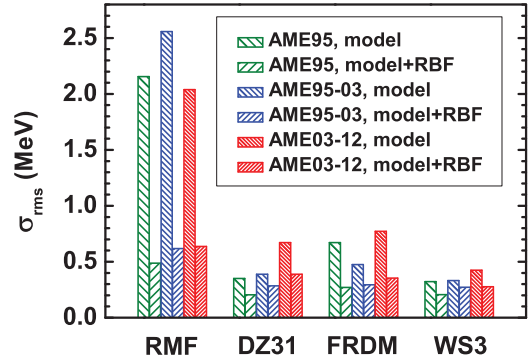


FIG. 4. (Color online) The rms deviations σ_{rms0} and σ_{rmsR} with respect to the known masses for different mass models in the AME95-03-12 test. For training the RBF, only those nuclei in AME95 are employed, while their masses are taken from AME12.

is systematically increased for all mass models. For mass models with smaller σ_{rms0} , i.e., DZ31 and WS3 models, the rms deviations σ_{rmsR} at $R_{\text{min}} = 3$ are similar to σ_{rms0} , which means that the RBF approach ceases to improve the model predictions effectively. However, the RBF approach is still effective for the FRDM and RMF mass models, even if nuclei with $r \leq 3$ are excluded. From Fig. 1, it is known that these two mass models have relatively longer mass correlations even with nuclei at $r > 3$, so the RBF approach still remarkably reduces their model deviations. However, it should be noted that the order of σ_{rmsR} among different models almost remains unchanged at various R_{min} in Fig. 3, i.e., $\sigma_{\text{rmsR}}(\text{WS3}) < \sigma_{\text{rmsR}}(\text{DZ31}) < \sigma_{\text{rmsR}}(\text{FRDM}) < \sigma_{\text{rmsR}}(\text{RMF})$.

AME95-03 has been extensively employed to check the predictive power of nuclear mass models in the literatures [26,56] and we extend this test to AME95-03-12 in this work. In the AME95-03-12 test, the nuclei in AME12 are separated into three subsets, i.e., the 1758 nuclei in the atomic mass evaluation of 1995 (AME95) [57], the 381 nuclei first appearing in AME03 [58], and the 214 “new” nuclei appearing in AME12. The rms deviations σ_{rms0} and σ_{rmsR} for different mass models in the AME95-03-12 test are shown in Fig. 4. For training the RBF, only those nuclei in AME95 are employed, while their masses are taken from AME12. From this figure, it is clear that the RBF approach significantly reduces the rms deviations of different mass models, especially for those models with large σ_{rms0} . The rms deviations σ_{rmsR} with respect to the masses of 1758 nuclei in AME95 are all within 0.5 MeV and the best predictive accuracy can reduce this to 0.206 MeV based on the WS3 mass models. The rms deviations σ_{rmsR} with respect to the masses of 381 nuclei first appearing in AME03 are slightly increased, while they are still remarkably smaller than the corresponding σ_{rms0} . The 214 “new” masses appearing in AME12 are not used in determining the effective interactions in any of mass models considered here, so these new data in AME12 are worthwhile for testing the predictive power of the RBF approach. It is found that the rms deviations σ_{rmsR} with respect to the 214 “new” masses appearing in AME12 are also remarkably reduced, and the best predictive accuracy is 0.277 MeV based on the WS3 mass models. From Fig. 5, it is shown that the 381 nuclei first appearing in AME03

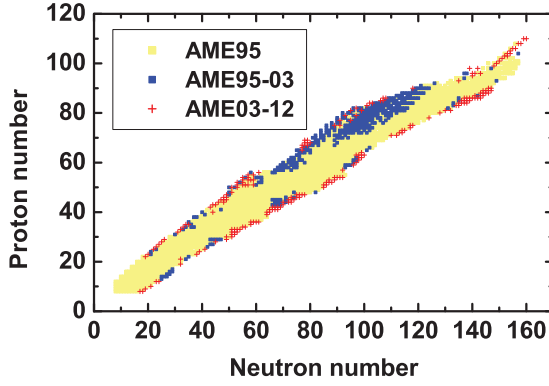


FIG. 5. (Color online) Positions of nuclei in the AME95-03-12 test.

and the 214 “new” nuclei appearing in AME12 are mainly around the nuclei in AME95 with $r \lesssim 3$, so the RBF approach can remarkably reduce the rms deviations for these nuclei.

By using the systematic trends in the mass surface and its derivative, the mass evaluation method in the atomic mass evaluation (AME) provides the best short-range mass extrapolation [1]. Therefore, it is interesting to compare the accuracy between the RBF approach and the method used in the AME. By taking the nuclei whose masses are evaluated values in AME03 (marked by “#” in the mass table of AME03) as an example, the rms deviations with respect to the new experimental data in AME12 are 0.398 MeV. For comparison with the method in the AME, the experimental data employed in training the RBF should be taken from AME03 as well, and the corresponding results based on different mass models are given in the third column of Table II. In addition, the rms deviations σ_{rms0} for these nuclei are given in the second column of Table II. Clearly, the RBF approach also remarkably improves the predictive power of various mass models. It should be pointed out that the rms deviation σ_{rmsR} based on the WS3 mass model is even smaller than that from the method in the AME. In Fig. 6, the experimental uncertainties in AME12 and AME03 are presented as a function of the isospin asymmetry $I = (N - Z)/A$. It is found that the experimental uncertainties are significantly improved in AME12 compared

TABLE II. The rms deviations in MeV with respect to the known masses in AME12 for the nuclei whose masses are evaluated values in AME03 (marked by “#” in the mass table of AME03). The second column represents the rms deviations σ_{rms0} for different mass models. The third and fourth columns both represent the rms deviations σ_{rmsR} by training the RBF with the nuclei in AME03, while the data are taken from AME03 and AME12, respectively.

Model	σ_{rms0}	$\sigma_{\text{rmsR}}(\text{AME03})$	$\sigma_{\text{rmsR}}(\text{AME12})$
RMF	1.956	0.604	0.620
HFB-21	0.631	0.559	0.483
DZ10	0.876	0.419	0.327
DZ31	0.673	0.430	0.325
ETFSI-2	0.690	0.484	0.415
KTUY	1.102	0.415	0.325
FRDM	0.771	0.469	0.371
WS3	0.425	0.375	0.268

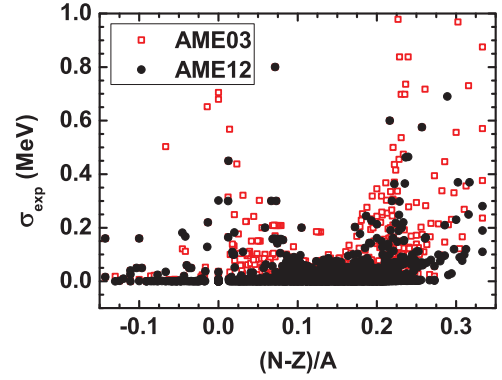


FIG. 6. (Color online) The experimental uncertainties as a function of the isospin asymmetry $I = (N - Z)/A$. The squares and circles represent the experimental uncertainties in AME03 and AME12, respectively.

with those in AME03, especially for those nuclei around the border region of experimental data with $I \lesssim 0.1$ and $I \gtrsim 0.2$. Therefore, we further update the data in training the RBF with those in AME12, and the corresponding results are shown in the fourth column of Table II. It is clear that σ_{rmsR} can be significantly reduced for most mass models. Based on the WS3 mass model, the predictive accuracy has been reduced to 0.268 MeV. Therefore, the new high-precision experimental data are also very important for improving the nuclear mass models with the RBF approach.

To investigate the radial basis function corrections in detail, the reconstructed functions $S(Z, N)$ based on the measured masses in AME12 are shown in Fig. 7 by taking the RMF and WS3 mass models as examples. It is clear that the reconstructed functions are sensitive to the nuclear mass models. For the RMF mass model, $S(Z, N)$ values are about 3 MeV for nuclei near $(Z, N) = (50, 50)$, $(Z, N) = (58, 82)$,

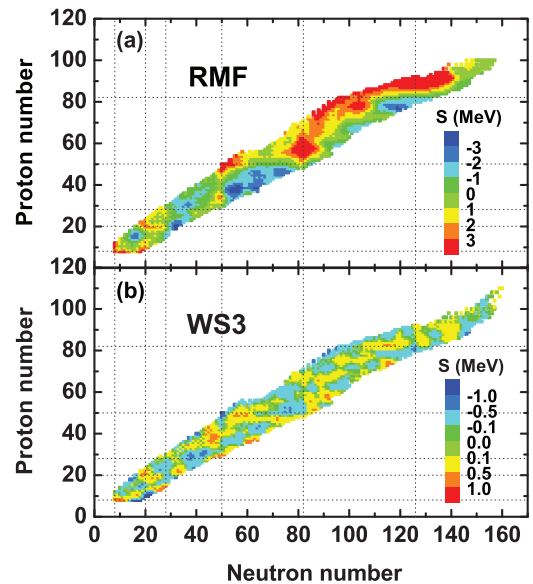


FIG. 7. (Color online) The reconstructed functions $S(Z, N)$ based on the measured masses in the AME12 for the RMF (a) and WS3 (b) mass models. The dotted lines denote the magic numbers.

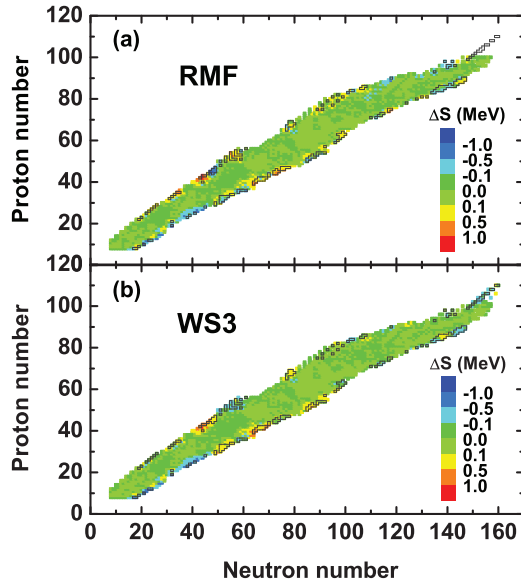


FIG. 8. (Color online) The differences of the reconstructed functions $S(N, Z)$ between those based on the measured masses in the AME12 and those based on the measured masses in AME03 for (a) RMF and (b) WS mass models, respectively. The “new” nuclei in AME12 are indicated by the black contours.

$(Z, N) = (78, 92)$, and $(Z, N) = (92, 126)$. This just corresponds to nuclei whose masses are underestimated in the RMF model [45]. Moreover, the overestimations of nuclear masses in the regions near $(Z, N) = (38, 60)$ and $(Z, N) = (78, 120)$ in the RMF model are also well improved by using the RBF approach with $S(Z, N) \sim -3$ MeV in these two regions. It better describes the nuclear masses for the WS3 mass model than for the RMF mass model, while there still exist small but systematically correlated errors [19,59]. With the RBF approach, these systematic correlations can be well extracted as well. To further investigate the influence of the “new” masses in AME12 on improving the nuclear mass models with the RBF approach, the differences of the reconstructed functions $S(N, Z)$ between those based on the measured masses in AME12 and those based on the measured masses in AME03 are shown in Fig. 8. It is found that the differences of reconstructed functions $S(N, Z)$ are generally within 100 keV for most nuclei, while they are relatively larger for those nuclei around the border region of experimental data, especially for those “new” nuclei in AME12. This can be well understood since significant improvements in the mass measurements have been made for nuclei near the border region in recent years [4]. Therefore, the RBF approach is sensitive to the experimental

masses and it is necessary to adopt high-precision experimental data to improve the nuclear mass models.

IV. SUMMARY AND PERSPECTIVE

In this work, the mass correlations in the radial basis function approach are carefully investigated based on eight widely used nuclear mass models, ranging from macroscopic-microscopic to microscopic types. The mass correlations usually exist between a nucleus and its surrounding nuclei with distance $r \lesssim 3$. However, the correlation distance is dependent on the nuclear mass model, which can go up to a distance of $r \sim 10$ for mass models with larger rms deviations, such as the RMF model. To extract these mass correlations, it is shown that nuclei at distance $r \leq 3$ should be included in the training of the RBF approach. In this way, the RBF approach can make significant improvements in the mass predictions for different mass models. The AME95-03-12 test further shows that the RBF approach provides a very effective tool for improving mass predictions significantly in regions not far from known nuclear masses. Based on the latest Weizsäcker-Skyrme mass model, the RBF approach can achieve an accuracy comparable with the extrapolation method used in the atomic mass evaluation, which can be further improved by the incorporation of new measurements. As claimed in the introduction, the effective interaction PC-PK1 remarkably improves the mass prediction of the RMF model. Therefore, it will be interesting to investigate the predictive power of the PC-PK1 mass model with the help of the RBF approach when the calculated masses with PC-PK1 for all nuclei in AME12 are available in the future. In addition, given the success in improving the nuclear mass predictions, the RBF approach has great potential for improving theoretical calculations of other physical quantities, such as nuclear β -decay half-lives, fission barriers, and excitation spectra.

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