



# Ground-state properties of uranium isotope chain: a relativistic mean field approach

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## Abstract

The study of uranium isotopes plays a crucial role in advancing our knowledge of nuclear physics, particularly in the realm of isospin and exotic nuclei. This study focused on the ground-state properties of uranium isotopes ranging from  $A = 203$  to  $A = 305$ . The key physical quantities examined included binding energy, quadrupole deformation, isotopic displacement, single-particle energy levels, and nucleon density distributions. Recent experimental advancements in uranium isotope studies have emphasized the indispensable role of theoretical models in interpreting experimental data. Moreover, the industrial applications of uranium—especially in nuclear energy production and weapons development—underscore the importance and necessity of accurate theoretical insights. The framework of the finite-range droplet model (FRDM) was utilized for comparative analysis because its predictions closely align with the experimental results. Through an analysis of the single-particle energy levels and continuous-state occupancy, this study identified  $^{207}\text{U}$  as the proton drip line nucleus. This research not only deepens our understanding of uranium isotopes but also provides a solid theoretical foundation to guide future experimental investigations.

**Keywords** Relativistic mean field · BCS theory · Uranium isotopic chain

## 1 Introduction

Exploration of the mass and charge limits of atomic nuclei is a fundamental challenge in nuclear physics. With advancements in heavy-ion accelerators and advanced detection systems, the synthesis of new superheavy elements, such as element 119, following the discovery of element 118, has gained significant attention [1–6]. These efforts are crucial not only for extending the periodic table but also for deepening our understanding of nuclear shell structures. The stability of heavy elements is intricately linked to the arrangement of protons and neutrons, and many nuclear

structure theories, from the shell model proposed by Mayer and Jensen [7] to modern microscopic approaches, have sought to predict these configurations, particularly the magic numbers.

A key element in the study of nuclear structures is the concept of "magic numbers." Magic numbers are specific numbers of nucleons (protons or neutrons) that result in highly stable atomic nuclei due to the closed shells in the nuclear structure, analogous to the electron shells in atoms. Historically, the classic magic numbers have been identified as 2, 8, 20, 28, 50, 82, and 126. However, recent theoretical advances and experimental observations have revealed the existence of new magic numbers, particularly in superheavy regions and exotic nuclei. For example, various nuclear models, such as the Skyrme–Hartree–Fock and relativistic continuum Hartree–Bogoliubov models [8], predict different sets of magic numbers beyond traditional ones. These include proton numbers  $Z = 114, 120$ , and  $126$  and neutron numbers  $N = 172, 184$ , and  $198$  [9, 10], which are expected to enhance the stability of superheavy nuclei. Moreover, recent research by Rydin [11] suggested additional magic numbers based on a geometrical packing approach, which

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implies new magic numbers such as  $Z = 90, 100$ , and  $118$ , as well as neutron numbers such as  $N = 58, 68$ , and  $76$ . These findings indicate that the landscape of magic numbers is far more nuanced than previously thought and that the concept of magicity continues to evolve as new experimental data become available.

Recently, the study of uranium isotopes has played an essential role in both theoretical and applied nuclear physics. Uranium is pivotal for nuclear energy and weapons applications and serves as a potential starting point for synthesizing superheavy nuclei [12] through nuclear decay pathways. Understanding the uranium isotope chain, particularly in terms of ground-state properties such as binding energy, deformation, and density distributions, provides insights into the broader aspects of nuclear stability and shell structure.

Recent experimental advancements have significantly contributed to the study of uranium isotopes, particularly the synthesis and characterization of new neutron-deficient isotopes such as  $^{215}\text{U}$ . In 2015, a new isotope,  $^{215}\text{U}$ , was produced using a complete fusion reaction involving  $^{180}\text{W}$  and  $^{40}\text{Ar}$ , followed by the separation of the evaporation residues using the gas-filled recoil separator SHANS. The identification of  $^{215}\text{U}$  was based on energy–position–time correlations, with an observed alpha-particle energy of  $8.428$  MeV and a half-life of approximately  $0.73$  ms [13]. Similarly, experiments have determined the properties of uranium, showing a consistent trend in the  $\alpha$ -decay behavior of neutron-deficient uranium isotopes [14]. These experimental efforts, aided by facilities such as the Heavy Ion Research Facility in Lanzhou (HIRFL), provide essential data that validate theoretical predictions and extend our knowledge of the stability and decay characteristics of heavy nuclei [15].

Theoretical advancements in nuclear physics have led to the development of several models to describe atomic nuclei, including first-principle methods [16–18], the shell model [19–22], and density functional theory (DFT) [23, 24]. Among these, relativistic mean field (RMF) theory, a variant of DFT, has proved to be a powerful tool for describing nuclear structures. The RMF theory incorporates relativistic effects, providing a more comprehensive description of nucleon interactions and allowing for the prediction of ground-state properties and magic numbers with notable accuracy. Previous studies have shown that the RMF theory, combined with the Bardeen–Cooper–Schrieffer (BCS) approach to treat pairing correlations [25, 26], is effective for describing the ground-state properties of isotopes near the proton drip line.

Despite these advances, several challenges remain in understanding the complete behavior of uranium isotopes, particularly those near the proton drip line, where conventional models face difficulties owing to the intricate interplay between pairing forces and continuum effects. In this study, we employed the RMF theory framework utilizing the TM1

parameter set to systematically investigate the ground-state properties of uranium isotopes ranging from  $A = 204$  to  $A = 305$ . Our work aims to provide new insights into the binding energies, deformation characteristics, and proton drip line behavior by comparing the RMF results with the finite-range droplet model (FRDM) [27, 28] predictions.

This paper is organized as follows: In Sect. 2, we provide a brief overview of RMF theory and its application in our study. Section 3 presents our analysis of the ground-state properties of the uranium isotopes. Section 4 discusses the identification of proton drip line nuclei within the uranium chain. Finally, in Sect. 5, we summarize the results and their implications for future experimental and theoretical studies in this field.

## 2 Methodology and theoretical framework

### 2.1 Relativistic mean fields

We used the Lagrangian density

$$\begin{aligned} \mathcal{L} = & \bar{\psi}_i \{i\gamma^\mu \partial_\mu - M\} \psi_i \\ & + \frac{1}{2} \partial^\mu \sigma \partial_\mu \sigma - U(\sigma) - g_\sigma \bar{\psi}_i \psi_i \sigma \\ & - \frac{1}{4} \Omega^{\mu\nu} \Omega_{\mu\nu} + \frac{1}{2} m_\omega^2 \omega^\mu \omega_\mu - g_\omega \bar{\psi}_i \gamma^\mu \psi_i \omega_\mu \\ & - \frac{1}{4} \vec{R}^{\mu\nu} \vec{R}_{\mu\nu} + \frac{1}{2} m_\rho^2 \vec{\rho}^\mu \vec{\rho}_\mu - g_\rho \bar{\psi}_i \gamma^\mu \vec{\tau} \psi_i \vec{\rho}_\mu \\ & - \frac{1}{4} F^{\mu\nu} F_{\mu\nu} - e \bar{\psi}_i \gamma^\mu \frac{(1 + \tau_3)}{2} \psi_i A_\mu. \end{aligned} \quad (1)$$

The first row of the nucleon terms,  $\psi_i$ , is the wave function of a nucleon, where  $i$  represents a nucleon inside the nucleus. The next three terms are the  $\sigma$  meson,  $\omega$  meson, and  $\rho$  meson terms.  $M$ ,  $m_\sigma$ ,  $m_\rho$ , and  $m_\omega$  are the masses of the nucleus and three mesons, respectively.  $g_\sigma$ ,  $g_\omega$ , and  $g_\rho$  are the coupling constants of the three mesons, respectively. The values of the nonlinear potential  $U(\sigma)$  for the  $\sigma$  term [31] and those of the meson and electromagnetic field tensors are as follows:

$$U(\sigma) = \frac{1}{2} m_\sigma \sigma^2 + \frac{1}{3} g_2 \sigma^3 + \frac{1}{4} g_3 \sigma^4, \quad (2)$$

$$\begin{aligned} \Omega^{\mu\nu} &= \partial^\mu \omega^\nu - \partial^\nu \omega^\mu, \\ R^{\mu\nu} &= \partial^\mu \rho^\nu - \partial^\nu \rho^\mu, \\ F^{\mu\nu} &= \partial^\mu A^\nu - \partial^\nu A^\mu. \end{aligned} \quad (3)$$

The Dirac and Klein–Gordon equations were derived from the Lagrangian density using the variational method. For the numerical solution of these equations, an axially symmetric harmonic oscillator was used as the basis for expanding the wave function in cylindrical coordinates, allowing for the

effective treatment of deformed nuclei. Initially, trial calculations were performed using  $N_f = N_b = 12$  for the selection of the major shell components representing the number of oscillator shells for fermions and bosons. To improve the accuracy of the results, the model parameters were later refined to  $N_f = N_b = 20$ , the iteration limit was set to 1600, and the error value was  $10^{-7}$ , which ensures a more precise representation of the nucleon wave functions, especially in the context of deformed nuclear systems (Table 1).

To further enhance the analysis, three parameter sets—NL1, TM1, and NLSH—were utilized in this study. Each set provides different values for the coupling constants and meson masses, affecting the representation of nuclear interactions and, consequently, nuclear structure predictions. Specifically, the TM1 parameter set, widely acknowledged for its effectiveness in ground-state property calculations, was chosen as the primary model for most calculations presented in this work. For investigations near the proton drip line, comparisons were made among the NL1, TM1, and NLSH sets to assess their respective accuracy and consistency in predicting the properties of exotic nuclei. The parameters for each set are as follows [29–31]:

## 2.2 Ground-state properties

In the following, we outline the methods used to calculate several important nuclear properties, including binding energy, quadrupole deformation, continuum occupation numbers (or single-particle energy levels), and nucleon density distributions. The relationships and formulas used were derived from a study by Gambhir et al. (1990) [32]. These calculations provide comprehensive insight into the ground-state characteristics of uranium isotopes.

The average binding energy, an essential quantity for determining nuclear stability, was obtained by integrating the energy densities of the nucleons and meson fields. The total binding energy was expressed as follows:

$$E(\psi_i^\dagger, \psi_i, \sigma, \omega^0, \rho^0, A^0, v_i) = E_{\text{part}} + E_\sigma + E_\omega + E_\rho + E_e + E_{\text{pair}} + E_{\text{CM}} - AM, \quad (4)$$

where  $E_{\text{part}}$  represents the kinetic energy of the nucleons.  $E_\sigma$ ,  $E_\omega$  and  $E_\rho$  represent the contributions from the interactions mediated by the scalar meson field ( $\sigma$  meson), vector meson field ( $\omega$  meson), and isovector-vector meson field ( $\rho$  meson), respectively, which are responsible for the effective nuclear force between nucleons.  $E_e$  refers to the Coulomb energy, accounting for the electrostatic repulsion between protons.  $E_{\text{pair}}$  describes the pairing energy between the nucleons, which is particularly important for maintaining nuclear stability by minimizing the total energy.  $E_{\text{AM}}$  refers to the center-of-mass correction energy, which ensures the accuracy of the calculation by correcting for the spurious motion of the center of mass.

The quadrupole deformation parameter was used to quantify the shape of the nuclei. Deformation parameters were calculated based on the quadrupole moment of the nucleus.

$$Q = Q_n + Q_p = \sqrt{\frac{16\pi}{5}} \frac{3}{4\pi} A R_0^2 \beta, \quad (5)$$

where  $R_0 = 1.2A^{1/3}(\text{fm})$ . The quadrupole moments were calculated using the following expressions:

$$Q_{n,p} = \langle 2r^2 P_2(\cos \theta) \rangle_{n,p} = \langle 2z^2 - x^2 - y^2 \rangle_{n,p}. \quad (6)$$

For single-particle energy levels and continuum occupation numbers, we used the BCS approach to account for pairing correlations. The occupation number for each single-particle state is given by

$$n_i = v_i^2 = \frac{1}{2} \left( 1 - \frac{\epsilon_i - \lambda}{\sqrt{(\epsilon_i - \lambda)^2 + \Delta^2}} \right), \quad (7)$$

$\epsilon_i$  represents the single-particle energy of the specific state  $i$ .  $\lambda$  is the chemical potential ensuring particle number conservation, and  $\Delta$  is the pairing gap, which quantifies the strength of the pairing interaction between nucleons, such as paired neutrons or protons. The coefficients  $\mu_i$  and  $v_i$  denote the probability amplitudes for a single-particle state  $i$  to be occupied or unoccupied, respectively, and they satisfy  $\mu_i^2 + v_i^2 = 1$ .

RMF theory is an effective framework for describing nuclear structures by considering the interactions between nucleons mediated by various mesons. However, this alone does not fully account for the pairing interactions between nucleons, which play a significant role in determining the stability and deformation properties of nuclei, particularly in open-shell configurations. To address this, the BCS theory [32, 33] was employed, as it allows for the effective treatment of these pairing correlations and provides a more

**Table 1** Parameter sets used in RMF model: TM1, NL1, and NLSH

Parameter	TM1	NL1	NLSH
$m_\sigma$ (MeV)	511.198	492.25	526.059
$m_\omega$ (MeV)	783.0	795.36	783.0
$m_\rho$ (MeV)	770.0	763.0	763.0
$g_\sigma$	10.0289	10.138	10.4434
$g_\omega$	12.6139	13.285	12.945
$g_\rho$	4.6322	4.975	4.382
$b$	$-7.2325 \times 10^{-4}$	$-6.9099 \times 10^{-4}$	$-6.9099 \times 10^{-4}$
$c$	0.6183	$-5.4965 \times 10^{-5}$	0.615

complete and accurate depiction of the nuclear ground-state properties. This is especially important for heavy elements, such as uranium, where pairing interactions influence many key properties, including binding energy and deformation.

Many properties of the nucleus exhibit a parity dependence, leading to distinct pair correlations between protons and neutrons. The RMF approach used here incorporates these pair interactions by employing the BCS theory as a perturbative correction, specifically incorporating the effective pair force constants for protons and neutrons, given by

$$G_n = \frac{21}{A} \left( 1 - \frac{N - P}{2A} \right), \quad (8)$$

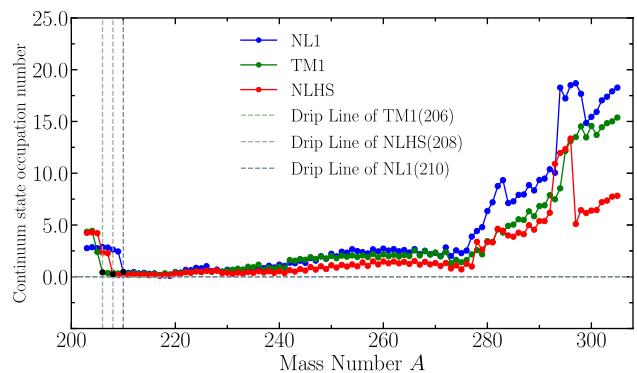
$$G_p = \frac{27}{A} \left( 1 + \frac{N - P}{2A} \right),$$

where  $A$  is the mass number, and  $N$  and  $P$  represent the neutron and proton numbers, respectively. In addition, in systems with an odd number of nucleons, one nucleon remains unpaired, which can significantly affect pairing correlations. Thus, the blocking method is an important complementary technique used when dealing with this case of odd- $A$  nuclei, which ensures that this unpaired nucleon is "frozen" in its specific orbital, characterized by an energy level  $\epsilon_k$  thus preventing it from participating in the overall pairing interaction [34].

Thus, the integration of BCS theory, blocking method, and RMF theory forms a comprehensive framework for studying uranium isotopes. These methods allow for accurate calculation of pairing interactions, single-particle energies, and deformation parameters, which are essential for understanding both the stability and structural nuances of heavy nuclei, including those near the proton drip line.

### 3 Determination of drip line nucleus in U isotopic chain

To determine the proton drip line [35–37], we analyzed the continuum-state occupation number of uranium isotopes calculated with different relativistic mean field parameters: NL1, TM1, and NLSH. As depicted in Fig. 1, the continuum-state occupation number shows distinct behavior for different parameter sets. For isotopes with mass numbers ranging from  $A = 203$  to  $A = 305$ , the occupation number generally remains low for mid-range isotopes, indicating bound systems. However, as shown in the figure, each parameter set predicts a different mass number for the proton drip line nucleus. Specifically, the NL1 parameter set identifies  $A = 210$  as the proton drip line nucleus, whereas the TM1 and NLSH parameter sets predict proton drip lines at  $A = 206$  and  $A = 208$ , respectively. These differences

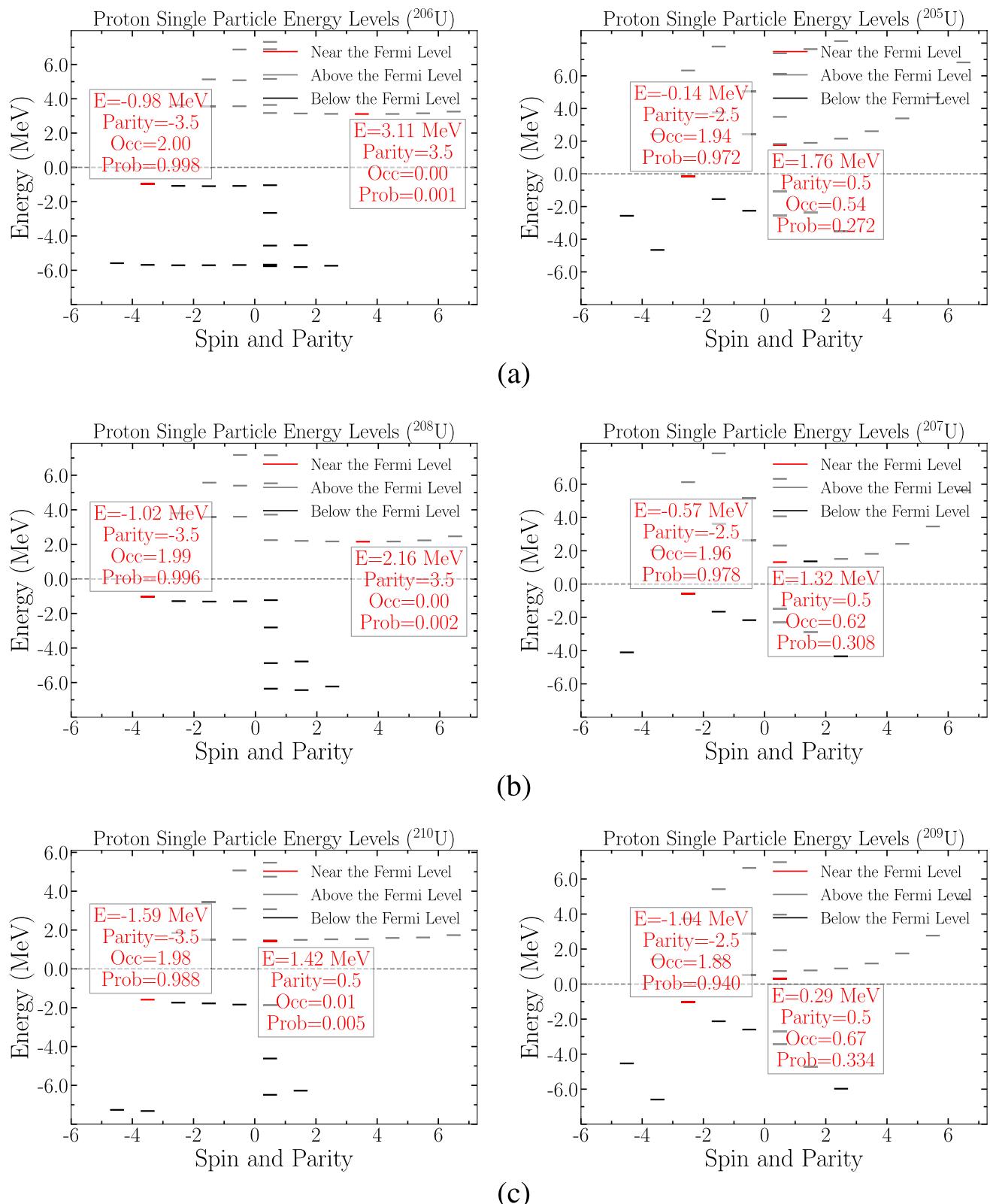


**Fig. 1** (Color online) Continuum-state occupation number and drip line of protons

highlight the sensitivity of drip line predictions to the choice of interaction parameters.

In Fig. 2, the single-particle energy levels near the Fermi surface of the drip line nuclei and their neighboring isotopes are displayed for NL1, TM1, and NLSH parameters. These energy level distributions illustrate that for nuclei at the proton drip line, the occupation of continuum states becomes prominent, signifying their unbound nature. Thus, the proton drip line kernels for the three parameter sets are  $^{209}\text{U}(\text{NL1})$ ,  $^{205}\text{U}(\text{TM1})$ ,  $^{207}\text{U}(\text{NLSH})$ . Importantly, different results for the three parameter sets reflect the sensitivity of the RMF model to the choice of parameters. In the absence of definitive experimental data, it is advisable to consider the results from multiple parameter sets to form a more representative prediction and acknowledge this uncertainty in the analysis.

To describe the region of nuclei with relatively more protons along the isotopic chain by applying the better parameter set NLSH, we consider the proton drip line nuclei as a result of the calculation of the NLSH parameter set  $^{207}\text{U}$ . For the NLSH parameter set, the analysis of  $^{207}\text{U}$  and  $^{208}\text{U}$  reveals significant details regarding the occupation of energy states near the Fermi surface. In  $^{207}\text{U}$ , the first positive-energy state above the Fermi surface, identified as  $\frac{1}{2}^+$ , has an energy of 1.32 MeV and an occupation probability of 30.8%. This suggests that a substantial fraction of protons occupies an unbound state, indicating that  $^{207}\text{U}$  is at the proton drip line for the NLSH parameter set. Conversely, for  $^{208}\text{U}$ , the last proton resides in a negative-energy state, and the first positive-energy state above the Fermi surface shows zero occupancy and occupation probability. This clearly highlights the difference in the binding behavior between  $^{207}\text{U}$  and  $^{208}\text{U}$ , confirming that  $^{207}\text{U}$  lies on the proton drip line, whereas  $^{208}\text{U}$  remains bound and stable. This analysis provides further evidence of the NLSH parameter set's balanced approach in predicting the proton drip line by accurately capturing the transition between bound and unbound states.

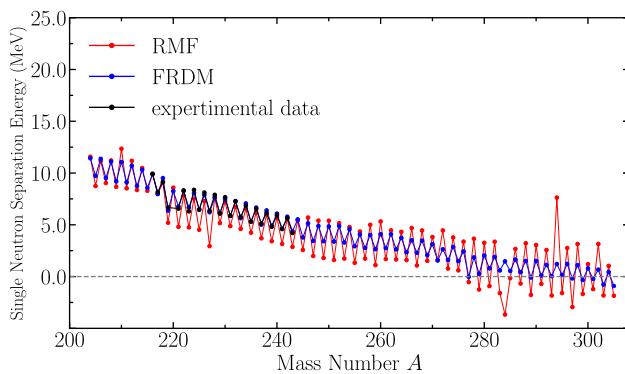


**Fig. 2** Energy levels near Fermi surface of dripline nuclei for NL1, TM1, and NLSH parameters. **a** Single-particle energy levels of  $^{205}\text{U}$  and  $^{206}\text{U}$  for TM1; **b** Single-particle energy levels of  $^{207}\text{U}$  and  $^{208}\text{U}$  for NLSH; **c** Single-particle energy levels of  $^{209}\text{U}$  and  $^{210}\text{U}$  for NL1

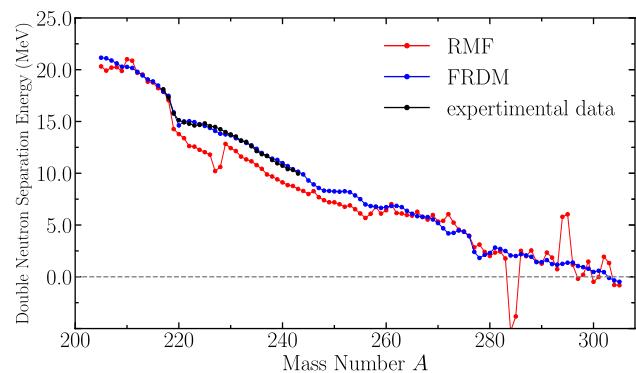
As the occupation number of the continuum state fluctuated significantly on the neutron-rich side of the uranium isotopic chain, a distinct anomaly was observed compared with the proton side. This anomaly is particularly evident at mass numbers  $A = 277$  and  $A = 278$ , where the continuum occupation number exhibits an abrupt change from 0.99 to 3.38. This step-like transition suggests that the mean field strength and pairing correlation strength are comparable in magnitude for these nuclei. In such cases, treating pairing correlations merely as perturbations is no longer valid because their contribution is significant enough to substantially influence nucleon dynamics. Consequently, pairing correlations must be rigorously included in self-consistent equations of motion rather than as a small correction.

The improper perturbative treatment of these pairing effects leaves some neutron-rich nuclei in a state where the last neutrons occupy continuum levels, which is unphysical under these conditions and results in an incorrect representation of the binding properties of the nucleus occupying continuum states. This implies that the pairing correlations should be inherently included in the mean field framework, as these correlations are essential for stabilizing the nuclear system and determining the drip line. Therefore, the sharp increase in the continuum occupation number at  $A = 277$  and  $A = 278$  cannot be interpreted accurately without consistently incorporating these pairing correlations.

To further clarify the determination of the neutron drip line, we analyzed the one-neutron separation energy  $S_n$  and the two-neutron separation energy  $S_{2n}$ , as shown in Figs. 3 and 4. The experimental data were obtained from [38]. The results indicate that for the last nuclei in the uranium chain, neither the one-neutron nor the two-neutron separation energies exhibit a clear trend toward zero, which signifies a drip line. Instead, the separation energies decreased gradually without reaching a definitive cut-off, implying that these nuclei were still marginally bound. This lack of a clear zero crossing in the separation energies suggests that using only



**Fig. 3** (Color online) One-neutron separation energy  $S_n$  of U isotope chains



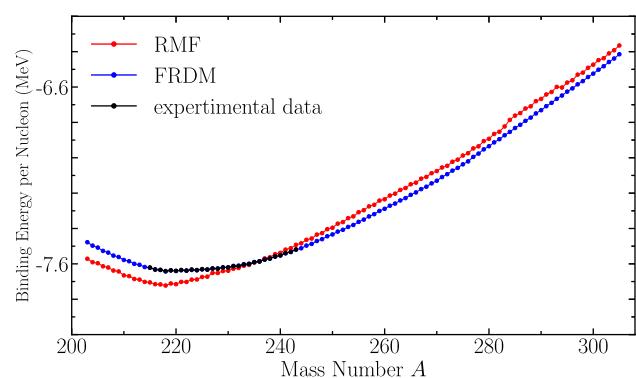
**Fig. 4** Two-neutron separation energy  $S_{2n}$  of U isotope chains

the  $S_n$  and  $S_{2n}$  values is insufficient for precisely identifying the neutron drip line.

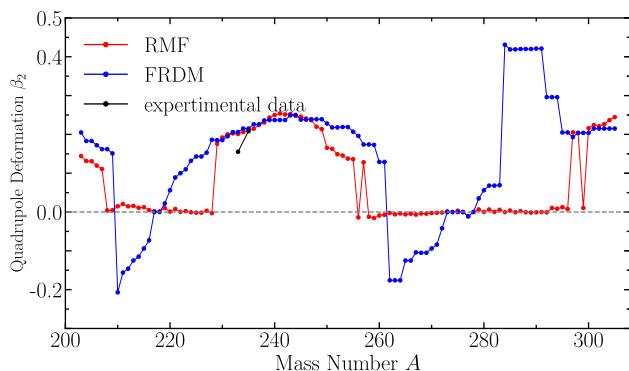
#### 4 Properties of ground state of nuclei in U isotope chain

The average binding energies are presented in Fig. 5. It is evident that our calculations are in good agreement with the finite-range droplet model (FRDM) data, with the lowest binding energy observed at the neutron magic number  $N = 126$ . This agreement indicates that our chosen force constants, treatment of pairing correlations, and implementation of the blocking method for odd- $A$  nuclei provide a reliable theoretical framework. Furthermore, our results show that at  $N = 126$  (corresponding to  $^{218}\text{U}$ ) [39], the average binding energy reaches its maximum value. The binding energy of  $^{218}\text{U}$  calculated with our approach is notably higher than that predicted by the FRDM, suggesting the enhanced representativeness of our model.

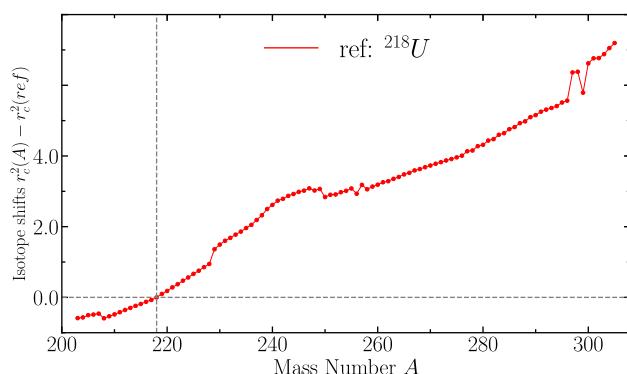
The deformation of the uranium isotope chains is shown in Fig. 6. The red line represents the finite-range droplet model (FRDM) predictions, whereas the black line



**Fig. 5** (Color online) Binding energy per nucleon



**Fig. 6** (Color online) Deformation of nuclei in chain of uranium isotopes for 20 shells



**Fig. 7** Isotope shifts  $r_c^2(A) - r_c^2(\text{ref})$  for U isotope chain

corresponds to the relativistic mean field (RMF) calculations. A comparison reveals that, unlike the FRDM results, the RMF model shows a smoother deformation [40] trend without the abrupt changes observed at mass numbers  $A = 284 - 296$ . This smoother trend suggests that the RMF approach provides a more consistent representation of the quadrupole deformation, particularly in regions where spherical symmetry is expected. Additionally, the deformation pattern indicates that nuclei with mass numbers less than  $A = 208$  exhibit a prolate, elongated ellipsoidal shape, whereas those between  $A = 208 - 228$  are almost spherical and exhibit very small deformations. Beyond  $A = 228$ , the deformation alternates between prolate and spherical, except at  $A = 256$ , where the deformation suddenly decreases to a spherical shape. Throughout the isotopic chain, there are no pronounced oblate or flat ellipsoidal shapes, and the few negative deformation values can be interpreted as spheroidal rather than strongly oblate shapes.

Using  $^{218}\text{U}$ , a semi-magic nucleus, as a reference, we analyzed the isotope shifts presented in Fig. 7. The data exhibited a generally smooth increase, with noticeable flattening for nuclei with mass numbers  $A = 247$  to  $A = 256$ .

In nuclear physics, a "kink" in isotope shift data refers to an abrupt change in the trend of nuclear charge radii as a function of neutron number, typically occurring at neutron magic numbers. This phenomenon has been observed in rare-earth elements, where such kinks appear near neutron magic numbers, indicating changes in nuclear structure and stability [39, 41–46].

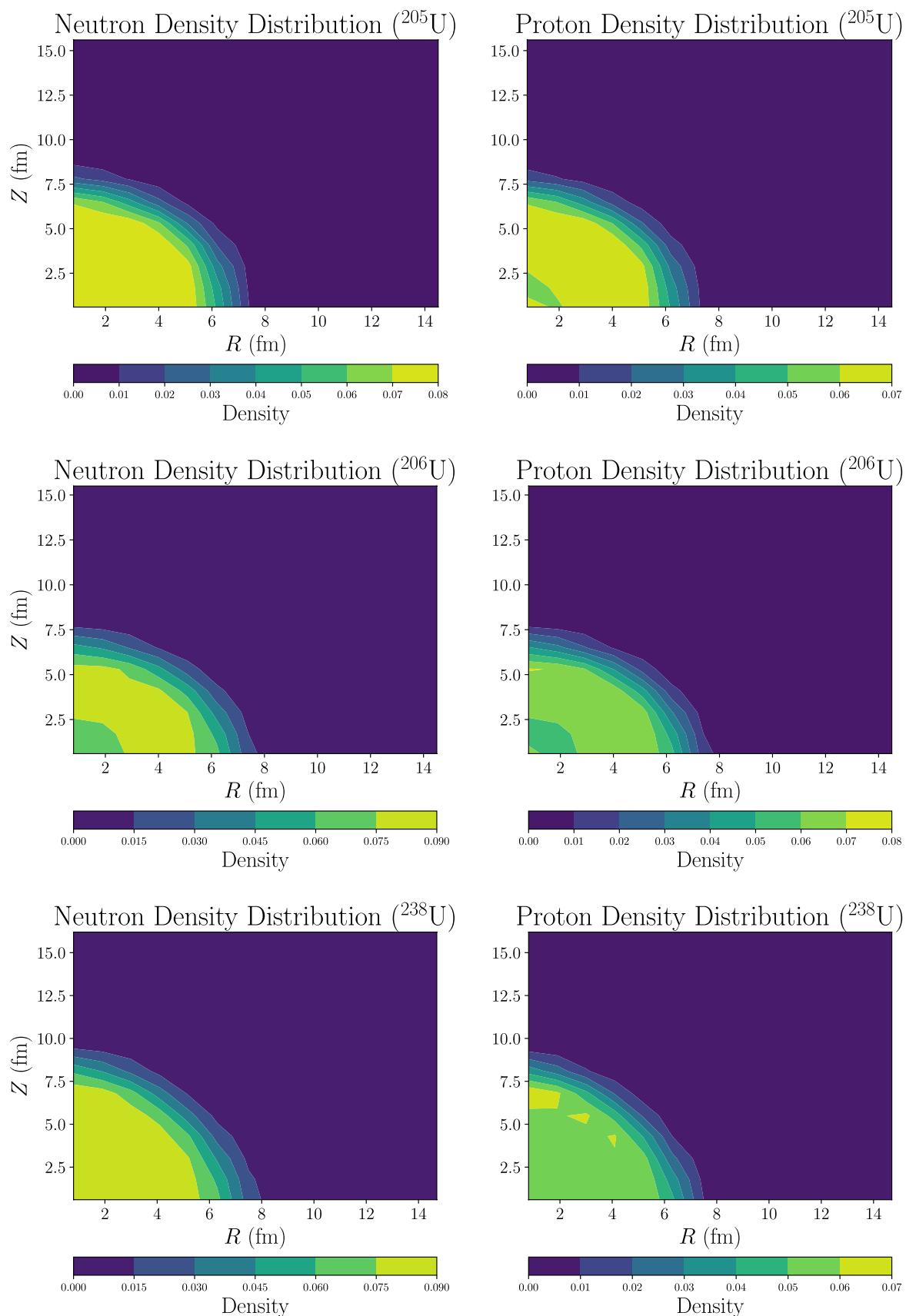
However, in our study of uranium isotopes, no evident kink was observed at  $A = 218$ , which corresponds to the neutron magic number  $N = 126$ . This absence suggests that the expected shell closure effect at  $N = 126$  does not manifest prominently in the U isotope shift data. Consequently, based on the isotope shift analysis alone, we cannot confirm that  $A = 218$  is a magic number nucleus. This finding is consistent with those of similar studies [47, 48], indicating that the manifestation of magic numbers can vary across different elements and isotopic chains (Fig. 8).

The analysis of the neutron and proton density distributions along the major axes of uranium isotopes, including  $^{205}\text{U}$ ,  $^{206}\text{U}$ , and  $^{238}\text{U}$ , provides important insights into their structural characteristics and deformation properties.

First, the density distributions of protons and neutrons are generally centered around the nucleus, with a distinct peak near the core. In all isotopes analyzed, the neutron density extends further than the proton density, indicating the presence of a "neutron skin." The neutron skin, where neutrons dominate the outer regions of the nucleus, is a common feature of neutron-rich heavy nuclei and contributes to the enhanced stability of these uranium isotopes. However, the proton density is more concentrated toward the center, which reflects the influence of Coulomb repulsion, pushing protons inward to counterbalance their mutual repulsive forces.

Among the isotopes,  $^{205}\text{U}$  displayed the smallest deformation, suggesting a nearly spherical shape, whereas  $^{238}\text{U}$  exhibited the largest deformation, characterized by significant elongation along the major axis.  $^{238}\text{U}$  shows intermediate deformations that can be described as prolate and resembling elongated ellipsoids. The neutron skins in these nuclei become more pronounced as the mass number increases, which agrees with the increasing neutron-to-proton ratio [49].

These findings collectively emphasize the significance of neutron excess and nuclear deformation in determining the density profiles of uranium isotopes. The presence of neutron skins in all the analyzed isotopes indicates that neutrons dominate the periphery of these nuclei, which has important implications for understanding nuclear stability, interaction cross sections, and the behavior of these nuclei near the dripline. RMF theory effectively captures these differences in density distributions and provides a comprehensive picture of the underlying nuclear structure of heavy isotopes, contributing to a better understanding of their stability and deformation characteristics.



**Fig. 8** (Color online) Density distribution of neutrons and protons

## 5 Summary

We calculated the ground-state properties of nuclei in the uranium isotope chain using the relativistic mean field (RMF) theory, incorporating pairing correlations through the Bardeen–Cooper–Schrieffer (BCS) approach. Our results, which show good agreement with the finite-range droplet model (FRDM) data, particularly at the neutron magic number  $N = 126$  validate the robustness of our approach in modeling the binding energies and deformation properties of uranium isotopes. By analyzing the Fermi surface and single-particle energy levels, we confirmed that  $^{207}\text{U}$  is a proton dripline nucleus due to its continuum-state proton occupancy, whereas  $^{208}\text{U}$  remains bound, highlighting the precise identification of the dripline.

In the future, the continued development of RMF theory, incorporating more advanced treatments of pairing and beyond-mean-field effects, holds great promise for understanding the properties of nuclei near the dripline, particularly for superheavy elements, which are crucial for practical applications, especially in nuclear fission processes relevant to energy production and reactor safety. Additionally, advancements in experimental facilities will be key to validating theoretical predictions and exploring new regions of the nuclear chart. The synergy between theoretical advancements and experimental verification will deepen our understanding of nuclear structure, stability, and broader implications for nuclear technology.

**Author contributions** All authors contributed to the study conception and design. Material preparation, data collection, and analysis were performed by Di Xu and Hong-Fei Zhang. The first draft of the manuscript was written by Di Xu, and all authors commented on previous versions of the manuscript. All authors read and approved the final manuscript.

**Data availability** The data that support the findings of this study are openly available in Science Data Bank at <https://cstr.cn/31253.11.sciedb.j00186.00717> and <https://doi.org/10.57760/sciedb.j00186.00717>.

## Declarations

**Conflict of interest** The authors declare that they have no conflict of interest.

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