N = 16 subshell closure from stability to the neutron drip line

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The Hartree-Fock-Bogoliubov method implemented with the Gogny-D1S force and used in a systematic study of the even-even N = 16 isotones from stability to the neutron drip line strongly suggests that a subshell gap, approximately 4 MeV wide, is opening at N = 16. The first 2⁺ levels as well as $B(E2; 0_{gs}^+ \rightarrow 2^+)$ reduced E2 transition probabilities have also been predicted in configuration mixing and standard random phase approximation calculations using D1S. These results compare favorably with experimental information available for the Z = 10-18 nuclei and with expectations for ²⁴O properties. They also fully support our mean-field predictions, namely that N = 16 is a spherical magic number at the oxygen neutron drip line.

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I. INTRODUCTION

One of the most striking pieces of evidence for quantum mechanics in physics is the existence of magic numbers. From atomic physics to nuclear physics, they are associated with a shell structure and a spherical configuration for which all the available individual levels are filled. In atomic physics, the noble gases are well-known examples of enhanced stability in their chemical properties. In nuclear physics, the same situation indeed occurs but for magic numbers different from those found in atomic physics. The sequence of magic numbers for stable nuclei was explained theoretically for the first time fifty years ago [1,2] and is now well understood. The properties of magic nuclei like ⁴⁰Ca or ²⁰⁸Pb were extensively studied and are the first observables to validate the shell model picture. However, the magnitude of shell gaps may evolve from stability toward drip lines. It may decrease, as already observed at neutron number N = 20 for which spherical magicity vanishes at the neutron drip line [3]. This magnitude may increase as well and give rise to a new magic number at the drip line, one that is not observed for stable nuclei. Recent experimental investigations far away from the valley of stability highlight modifications in the shell model structure. Several observables show that N = 16 neutron-rich isotones present a higher stability compared to their neighbors on the *N*-*Z* chart, namely,

- (i) the N = 16 isotones ²²C, ²³N, and ²⁴O are the last bound nuclei of their respective isotopic chains [4,5],
- (ii) the large energy gap around the Fermi surface for the $N = 16^{28}$ Mg and ²⁶Ne nuclei causes a break in the neutron separation energy S_n systematics, and
- (iii) the first 2⁺ level of ²⁴O has been suggested in a recent γ -spectroscopy experiment [6] to lie at higher excitation energy than the one-neutron separation threshold $S_n = 3.7 \text{ MeV}$ [7]. Such an excitation energy is much higher than the experimental $E(2^+)$'s observed for the other oxygen isotopes.

For these reasons, N = 16 is proposed to be a magic number for the most neutron-rich isotones [8]. This means an enhancement of the spherical gap between the $s_{1/2}$ and the $d_{3/2}$ subshells of the neutron *sd* shell compared to its value for stable nuclei [9–11].

In this paper, we study N = 16 as a possible new spherical magic number for neutron-rich nuclei. Our approach mainly relies on the Hartree-Fock-Bogoliubov (HFB) method. We show that the shell evolution may be interpreted and quantitatively reproduced with the Gogny-D1S effective interaction [12,13]. Self-consistent mean-field methods [14,15] have already been used to study light nuclei. In a previous work [16], the evolution from the stability region to the neutron-rich isotopes of the N = 20 and N = 28 shells has been discussed. Here, we study the N = 16 isotones from ${}^{22}C$, at the neutron drip line, to ${}^{34}Ar$ through a systematic description of even-even nuclei. Axial and triaxial deformations are investigated.

In the following, we briefly recall the HFB formalism and notation. The results from pure mean-field calculations restricted to axial symmetry are presented as a first step. Pairing energies are discussed. Extending the HFB calculations to triaxial shapes, collective 2^+ levels, and reduced transition probabilities B(E2)'s are derived from the generator coordinate method (GCM) implemented within the Gaussian overlap approximation (GOA). This configuration mixing approach complements the mean-field description and also points to N = 16 as a spherical neutron magic number far away from the stability line. Finally, standard random phase approximation (RPA) calculations using the same effective force are performed for ²⁴O.

II. MEAN-FIELD DESCRIPTION

In the constrained HFB theory, the deformed states $|\Phi_q\rangle$ of a nucleus are described as quasiparticule (qp) vacua. The $|\Phi_q\rangle$ states are deduced from the minimization of the energy functional

$$\delta \langle \Phi_q | \hat{H} - \lambda_N \hat{N} - \lambda_Z \hat{Z} - \sum_i \lambda_i \hat{Q}_i | \Phi_q \rangle = 0, \qquad (1)$$

where $|\Phi_q\rangle$ is the HFB wave function of deformation q, and \hat{H} is the many-body nuclear Hamiltonian. In this study, the finite range Gogny-D1S interaction is used, and both mean

and pairing fields are calculated in a consistent way. The λ 's in Eq. (1) are the Lagrange multipliers associated with the constraints on nucleon numbers N and Z and average deformations Q_i :

The \hat{Q}_i 's are taken as the quadrupole operators \hat{Q}_{20} and \hat{Q}_{22} , related to the axial and triaxial Bohr deformations β and γ by

$$\beta = \sqrt{\frac{\pi}{5}} \frac{\sqrt{(Q_{20}^2 + 3Q_{22}^2)}}{AR^2}, \ \gamma = \arctan\frac{\sqrt{3}Q_{22}}{Q_{20}}, \tag{3}$$

where A and $R = \sqrt{\frac{3}{5}}(1.2A^{1/3})$ (fm) are the mass number and radius of the nucleus under consideration, respectively.

We first investigate within constrained axial calculations the N = 16 isotones to test their magicity. At this stage, we consider that neutron (proton) magicity occurs when two criteria are fulfilled: (i) The nucleus is spherical; that is, the potential energy is minimum at zero deformation, and (ii) a large energy gap separates the levels above and below the Fermi surface. The second condition is satisfied when the neutron (proton) pairing energy cancels at zero deformation in this pure mean-field description. In the HFB formalism, the neutron (proton) pairing energy is given by $E_{n(p)} = \frac{1}{2} \text{Tr} \Delta \kappa$, where Δ and κ are the pairing field and the pairing tensor, respectively [17]. Such a correlation energy is not the net gain between the HFB and Hartree-Fock (HF) treatments. However, when the pairing field vanishes, HF and HFB descriptions give rise to the same total potential energy and no correlation energy is found at the mean-field level. Consequently, the study of pairing energies $E_{n(p)}$ provides direct information about pairing correlations.

In the present work, the mean-field equations are solved using basis sets of deformed harmonic oscillators including $N_0 = 9$ major shells. Such a large basis ensures stability of level energies at all deformations up to $\beta \simeq 1$. As an example, Fig. 1 illustrates the convergence of the neutron pairing energy E_n and the potential energy $V = \langle \Phi_q | \hat{H} | \Phi_q \rangle$ for ²⁶Ne with the basis size increasing from $N_0 = 7$ to $N_0 = 13$. Calculations with $N_0 = 7$ clearly do not reach the convergence, whereas those with $N_0 = 9$, 11, and 13 give the same results within 400 keV for V and 200 keV for E_n . Then, a reasonable convergence is achieved for $N_0 = 9$. This result holds true for all the other isotones under study.

All the axial potential energy surfaces (PESs) of the N = 16 isotones, shown in Fig. 2, display a minimum at $\beta = 0$. Thus, the first criterion (i) for magicity is fulfilled for all isotones at the pure mean-field level. Nevertheless, this minimum is shallow for nuclei from ²⁶Ne to ³⁴Ar and their neutron pairing energies, shown in Fig. 3, do not vanish at zero deformation. According to criterion (ii), these nuclei are not magic. The neutron pairing energy vanishes at $\beta = 0$ only for ²⁴O and ²²C. Then, criteria (i) and (ii) are simultaneously verified only for these two nuclei, which are predicted to be magic.

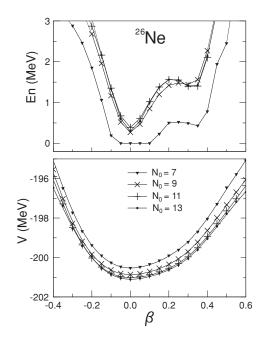


FIG. 1. Neutron pairing energies E_n (upper panel) and potential energy surfaces V (lower panel) of ²⁶Ne for different basis sizes as functions of the axial deformation parameter β . The numbers of major shells considered are $N_0 = 7, 9, 11$, and 13.

Furthermore, proton pairing energy vanishes at $\beta = 0$ for ²⁴O, which is then found to be doubly magic.

A gradual evolution is observed in Fig. 3 for the neutron pairing energy: Neutron pairing energy goes to zero around $\beta \simeq 0.4$ for ³⁴Ar, ³²S, and ³⁰Si. ³²S presents a shoulder at $\beta = 0$. Moving toward the drip line, this shoulder gets more pronounced and becomes a local minimum for ³⁰Si, ²⁸Mg, and ²⁶Ne. The neutron pairing energy vanishes at $\beta = 0$ for ²⁴O and ²²C. The evolution of these two pairing energy minima is directly linked to the existence of prolate and

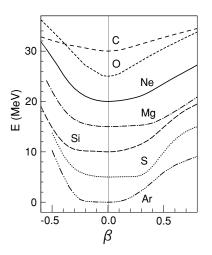


FIG. 2. Potential energy surfaces of the N = 16 isotones from argon to carbon, as functions of the axial deformation parameter β . The minimum potential energy for argon is set to zero. The minima of the curves are arbitrarily separated by 5 MeV. Stable nuclei are ³²S and ³⁰Si.

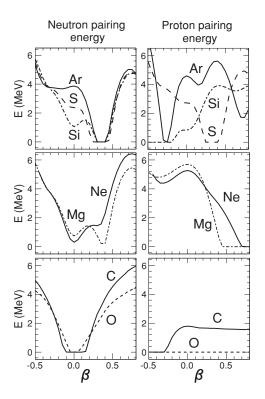


FIG. 3. Neutron (left) and proton (right) pairing energy curves for N = 16 isotones from argon to carbon, as functions of the axial deformation parameter β .

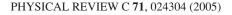
spherical gaps between the $K^{\pi} = 1/2^+$ orbitals (where *K* is the projection of the angular momentum on the intrinsic symmetry axis) from the $s_{1/2}$ and $d_{3/2}$ neutron subshells. This subshell structure is shown in the lower panel of Fig. 4 for ³²S and ²⁶Ne. Whereas ³²S has a 6-MeV-wide prolate gap and a 3.3-MeV-wide spherical gap, ²⁶Ne presents an inverse trend: These gaps are 3.5 and 4.2 MeV wide, respectively.

The upper panel of Fig. 4 shows the opening of the spherical gap δ between the $s_{1/2}$ and $d_{3/2}$ neutron subshells for nuclei from Z = 18 to Z = 6. For ²⁴O, this gap is $\delta \simeq 4$ MeV wide. To check the stability of this prediction, δ has also been calculated using the older D1 [12] parametrization of the Gogny force. The results do not depend significantly on the force parametrization (see Table I). The difference between the two calculations does not exceed 600 keV for δ . Both D1 and D1S parametrizations lead to the prediction of N = 16 as a magic number at the neutron drip line.

To summarize, N = 16 is predicted to be a spherical magic number at the neutron drip line for ²⁴O and ²²C in a pure mean-field approach. It preexists as a strong shell effect in nuclei of the stability region.

III. BEYOND THE MEAN FIELD

To complement the description of these nuclei, an approach beyond the mean field is considered. A dynamical description of the ground state and low-lying 2⁺ collective state is derived from GCM considering rotational and vibrational degrees of freedom in the triaxial plane. Collective states are taken as



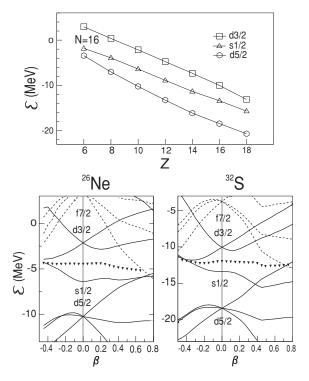


FIG. 4. (Upper panel) Neutron single-particle energies for eveneven N = 16 isotones at $\beta = 0$. (Lower panel) Neutron individual levels around the Fermi surface of ³²S (left) and ²⁶Ne (right) as functions of the axial deformation parameter β . Triangles represent the chemical potential.

linear combinations of constrained HFB basis states $|\Phi_q\rangle$, namely,

$$|\Psi_k\rangle = \int dq \ f_k(q) \, |\Phi_q\rangle \,. \tag{4}$$

Here, q is a five-dimensional generator coordinate, with two dimensions (β, γ) for vibrations and three Euler angles for rotation, and $f_k(q)$ are weight functions, solutions of the Hill-Wheeler equation [18]. The GOA transforms the integrodifferential Hill-Wheeler equation into a computationally tractable Schrödinger-like equation [19]:

$$\hat{\mathcal{H}}_{\text{coll}} g_k(q) = E_k g_k(q), \tag{5}$$

where $g_k(q)$ is the Gauss transform of $f_k(q)$ and eigenstate of the collective Hamiltonian $\hat{\mathcal{H}}_{coll}$ [20,21].

It is a widespread practice to perform calculations of mass and inertia parameters using the Inglis and Beylaev approximation [22,23]. This is a reasonable assumption at low

TABLE I. Spherical energy gap δ between the $s_{1/2}$ and $d_{3/2}$ subshells at $\beta = 0$ predicted using the D1S and D1 interactions in HFB calculations.

	δ (MeV)									
	³⁴ Ar	³² S	³⁰ Si	²⁸ Mg	²⁶ Ne	²⁴ O	²² C			
D1S D1	2.58 2.33	3.32 2.89	4.00 3.24	4.11 3.53	4.25 3.87	4.30 4.17	4.75 4.74			

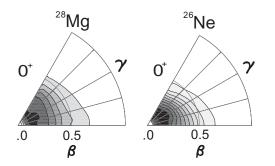


FIG. 5. Collective wave functions of the ²⁶Ne and ²⁸Mg ground states in the (β, γ) triaxial plane.

spin and excitation energy only for medium- and heavy-mass nuclei. As has been shown recently, improvements over the Belyaev formula for moments of inertia were necessary to provide a good description of 40,42,44 S low-energy spectra [24]. In the present work focusing on still lighter mass nuclei, two sets of moments of inertia, $J^A_{x,y,z}$ and $J^B_{x,y,z}$, have been calculated over the (β, γ) plane as follows. In the first set, $J^A_{x,y,z}$ is defined as

$$J_{x,y,z}^{A} = \frac{\langle \Phi_{q,\omega_{x,y,z}} | I_{x,y,z} | \Phi_{q,\omega_{x,y,z}} \rangle}{\omega_{x,y,z}}, \tag{6}$$

where $\hat{I}_{x,y,z}$ are the angular momentum projections in the intrinsic system of coordinates x, y, and z, and $\omega_{x,y,z}$ are the associated rotational frequencies for a nucleus with rotational energy E(I). First, J_z^A is calculated from minimization of the energy functional $\langle \Phi_{q,\omega_z} | \hat{H} - \lambda_N \hat{N} - \lambda_Z \hat{Z} - \omega_z \hat{I}_z | \Phi_{q,\omega_z} \rangle$ in the rotating frame [25], with the additional constraint $\langle \Phi_{q,\omega_z} | \hat{I}_z | \Phi_{q,\omega_z} \rangle = \hbar^2 \sqrt{I(I+1)}$. Here we use I = 2. The moments J_x^A and J_y^A are then calculated after permutation of the principal axes. Finally, the second set $J_{x,y,z}^B$ results from using the Thouless-Valatin approximation [26], namely,

$$J_{x,y,z}^{B} = \lim_{\omega_{x,y,z} \to 0} \frac{\langle \Phi_{q,\omega_{x,y,z}} | \hat{I}_{x,y,z} | \Phi_{q,\omega_{x,y,z}} \rangle}{\omega_{x,y,z}}.$$
 (7)

Equations (6) and (7) form two prescriptions for calculating moments of inertia. Each prescription has been tested separately while solving $\hat{\mathcal{H}}_{coll}$, and their figure of merit is discussed in the following. Using configuration mixing method outlined here, we find that ²⁴O and ²²C are spherical in their ground states whereas all the other nuclei from ³⁴Ar to ²⁶Ne are deformed. An illustration is provided for ²⁶Ne and ²⁸Mg in Fig. 5, where are shown their ground-state wave functions over the (β , γ) coordinates. The topology of these surfaces suggests γ -unstable deformed shapes with mean quadrupole deformations $\langle \beta \rangle = 0.31$ and 0.39, and $\langle \gamma \rangle = 24^{\circ}$ and 22° for ²⁶Ne and ²⁸Mg, respectively. This illustration provides a posteriori justification for our method in which the five collective quadrupole coordinates are explicitly treated.

The first 2^+ level energy and the $B(E2; 0_{gs}^+ \rightarrow 2^+)$ reduced transition probability values calculated with J^A and J^B moments of inertia as collective model inputs are shown in Fig. 6 as open circles and stars, respectively, for the N = 16 isotones from Z = 10 to Z = 18. Standard RPA calculations [27] have also been performed for the doubly magic ²⁴O

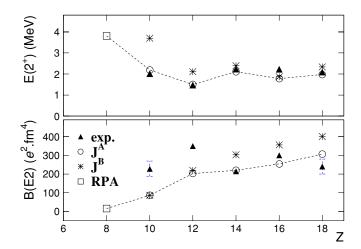


FIG. 6. (Color online) Energies of the first 2^+ excited state for even-even N = 16 isotones from Z = 18 to Z = 8. (Lower panel) Reduced transition probabilities $B(E2; 0_{gs}^+ \rightarrow 2^+)$. Experimental data (triangles) are available down to Z = 10. Both calculations with J^A (open circles) and J^B (stars) are represented from Z = 18 to Z = 10. Squares are standard RPA predictions for ²⁴O. Open circles and squares are connected by dashed lines to guide the eye.

nucleus using the Gogny-D1S force. In Fig. 6, RPA results (squares) are shown for completeness. Table II includes most of these predictions as well as experimental data [28] and mean deformations. As can be seen in the upper part of Fig. 6, the two sets of predictions for $E(2^+)$ (i.e., stars and open circles) significantly differ from each other, the differences being at a maximum for Z = 10 and Z = 12. However, as expected, the improvement in our determination of moments of inertia brings the $E(2^+)$ predictions marked by open circles in close agreement with each measured values (triangles) and with the pattern they display as Z increases from Z = 10 to Z = 18. The lowering observed for the 2^+ level in ²⁸Mg is here interpreted as stemming from a stronger mean deformation (see Table II).

Experimental information available on the structure of ²⁴O is rather scarce. In recent in-beam γ -ray spectroscopy measurements [6], the $2^+ \rightarrow 0^+_{gs} \gamma$ decay was not observed,

TABLE II. First 2⁺ level and $B(E2; 0_{gs}^+ \rightarrow 2^+)$ properties of even-even N = 16 isotones. J^A inertia parameters are adopted in the configuration mixing calculations for Z = 10-18 nuclei (see text). Standard RPA predictions using D1S are given for ²⁴O. Experimental data are from [28].

	$\langle eta angle$ gs 2^+		<i>E</i> (2 ⁺) (MeV)		$B(E2; 0^+_{gs} \rightarrow 2^+)$ $(e^2 \text{ fm}^4)$	
	Theoret.		Theoret.	Expt.	Theoret.	Expt.
³⁴ Ar	0.27	0.31	1.98	2.09	306	240(40)
^{32}S	0.31	0.32	1.78	2.23	254	300(13)
³⁰ Si	0.32	0.37	2.11	2.23	220	215(10)
²⁸ Mg	0.39	0.43	1.50	1.47	202	350(5)
²⁶ Ne	0.31	0.37	2.19	2.02	86	228(41)
²⁴ O	_	_	3.81	_	15	_

suggesting that the first 2^+ excited state lies above the neutron decay threshold $S_n = 3.7$ MeV. An excitation energy $E(2^+) \ge 3.7$ MeV would be roughly twice those for the Z = 10-18 members of the N = 16 isotonic chain and would be consistent with the present RPA prediction of $E(2^+) = 3.81$ MeV. These results together with those described earlier for the Z = 10-18 members of the N = 16 isotonic chain point to the existence of a subshell closure at the oxygen neutron drip line.

IV. CONCLUSION

In this paper, a systematic structure study of the even-even N = 16 isotones has been performed from stability to the neutron drip line. This has been achieved at both mean-field and configuration mixing levels using the Gogny-D1S effective interaction. Standard RPA calculations have also been performed for ²⁴O to complement our survey. These methods allow us to investigate single-particle and pairing properties at spherical shape and along axial and triaxial quadrupole deformations, as well as collective properties.

The mean-field calculations show that a spherical subshell gap is opening at N = 16 (see Fig. 4). Its size ($\delta \simeq 4$ MeV) does not significantly depend on whether the D1 or D1S parametrization of the Gogny force is used. Furthermore the

neutron pairing energy at spherical shape vanishes only for ²⁴O and ²²C. Since the potential energy surfaces of these nuclei also show a minimum at zero deformation, we conclude that N = 16 is a magic number at the neutron drip line, that is, for ²⁴O and ²²C.

Results from configuration mixing calculations bring confirmation that both nuclei are spherical in their ground states. In contrast, similar calculations suggest that all the Z = 10-18isotones show strong mean deformations. A new prescription tailored to calculate moments of inertia in light nuclei leads to improved $E(2^+)$ and $B(E2; 0_{gs}^+ \rightarrow 2^+)$ predictions that are in good overall agreement with experimental data available for the Z = 10-18 nuclei. Our predictions suggest that (i) collectivity of the $0_{gs}^+ \rightarrow 2^+$ transitions gradually decreases as Z gets away from Z = 18 and reaches a minimum at Z = 8, where it is 20 times weaker than at Z = 18, and (ii) $E(2^+)$ energies rise sharply from 1.98 MeV (³⁴Ar) to 3.81 MeV (²⁴O). These collective properties are strongly suggestive of N = 16as a magic number at the neutron drip line.

Experimental determination of the 2^+ excitation energy for ²⁴O is needed to challenge the reliability of our model predictions based on the Gogny-D1S force. It is now a key issue to precisely understand the reasons why this force is providing a subshell gap at the N = 16 neutron drip line. Work along this line is in progress.

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