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Two-Neutron Halo Structure and Anti-halo Effect in ³¹F

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Abstract We perform a detailed analysis for the structure of 31 F, which is a candidate of the halo nucleus. We calculate the radius and total reaction cross section using a three-body wave function of 29 F+n+n and discuss how the competition between the neutron-pairing and the single-particle energy induces structural changes of 31 F. The present analysis further clarifies a new aspect of the anti-halo effect that suppresses the halo structure in the ground state.

1 Introduction

Characteristic nuclear structure with loosely bound nucleons is called "halo" and has been extensively studied from both the theoretical and experimental sides [1,2]. In recent years, a dripline nucleus of the fluorine isotopes ³¹F has been observed [3]. Because of the small neutron separation energy and the subsystem ³⁰F (²⁹F+n) being unbound, ³¹F is expected to have a halo structure as found in Borromean nuclei, e.g., ⁶He, ¹¹Li, ¹⁴Be, and ¹⁹B. On the other hand, although development of the halo structure is closely related to the small neutron separation energy, a mechanism to suppress the nuclear radius in weakly bound systems has been discussed as the "pairing anti-halo" effect [4,5].

To give a theoretical prediction to the structure of 31 F, we performed three-body calculations using a 29 F+n+n model and showed the matter radius and reaction cross-sections of 31 F in Ref. [6]. This analysis revealed that the energy gap between the p-orbit and f-orbit in the subsystem 30 F is an essential ingredient to determine the radius of 31 F. Furthermore, a new aspect to the anti-halo effect was proposed. In this paper, we discuss the mechanism of the newly proposed anti-halo effect of 31 F more detail and clarify a general condition for the formation of the halo structure and the occurrence of the anti-halo effect.

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2 Theoretical Model and Interactions

We calculate the structure of ${}^{31}F$ in a ${}^{29}F+n+n$ three-body model. To describe ${}^{31}F$ as a loosely-bound Borromean system, the calculation method is required to be able to accurately take into account the continuum states. For this purpose, we employ the cluster-orbital shell model (COSM) [7] with the Gaussian expansion method (GEM) [8,9].

The Hamiltonian and basis functions in COSM with GEM for the $^{29}F+n+n$ system are briefly explained below. The Hamiltonian by removing the center of mass motion in COSM is formulated as follows:

$$\hat{H} = \sum_{i=1}^{2} (\hat{T}_i + \hat{V}_i + \lambda \hat{\Lambda}_i) + (\hat{t}_{12} + \hat{v}_{12}). \tag{1}$$

Here $\lambda \hat{\Lambda}$ is introduced to eliminate the Pauli forbidden states in the procedure of the orthogonality condition model [10], and \hat{t}_{12} is the recoil term induced by the subtraction of the center of mass motion.

We solve the eigenvectors of the Hamiltonian (1) using the Gaussian basis $\Phi_{pq\ell j}^{JM}$ in a variational way as $\Psi_{JM} = \sum_{pq\ell j} C_{pq\ell j} \Phi_{pq\ell j}^{JM}$ for the 0^+ state. Since we assume the core nucleus 29 F as a spin-less particle for simplicity, the spin-parity of 31 F is determined by the valence nucleons part $\Phi_{pq\ell j}^{JM} \equiv \mathcal{A}\left\{ [\phi_{p\ell j}(1) \otimes \phi_{q\ell j}(2)]_{JM} \right\}$, where \mathcal{A} is an antisymmetrizer. For the basis functions, we take the maximum orbital angular momentum ℓ up to 5, and the width parameters of the Gaussian are prepared using a geometric progression manner, where 20 bases are applied for each coordinate. Therefore, for example, a typical basis size is 2310 for the 0^+ state.

For the $^{29}\text{F}+n$ interaction \hat{V}_i , we employ the Woods-Saxon (WS) potential referring to the parameter sets to reproduce the total reaction cross sections of ^{31}Ne [11], which is the next nucleus of ^{30}F by a proton. Since there is no information other than ^{30}F being unbound, we adjust the strength parameter V_0 using two different values for each set of $\{r_0, a\}$ in WS under the condition that ^{30}F is unbound and ^{31}Ne is bound. This leads to 12 different parameter sets for the calculation [6]. For the neutron-neutron potential \hat{v}_{12} , we use the central part of the Minnesota potential [12] with the exchange parameter u=1.0.

3 Results and Discussion

First, we perform calculations using the 12 different parameter sets. As a result, the two-neutron separation energies S_{2n} are obtained in the range of 0.44 - 1.37 MeV, and the matter radius $R_{\rm rms}$ varies in the range of 3.48 - 3.70 fm. For the maximum radius case $R_{\rm rms} = 3.70$ fm, the halo structure is well developed. We also calculate the total reaction cross section at incident energies of 240 and 900 MeV/nucleon using the Glauber model (NTG) [13–18] and obtain large cross-sections as 1530 and 1640 mb [6]. This result further supports the large matter radius of 31 F, i.e., the halo structure, is realized under a certain condition in 30 F.

Next, to clarify the condition, we analyze the characteristics of the results for ${}^{31}F$ using the 12 parameter sets. To this end, we introduce the Pearson's correlation coefficient (PCC) between two variables x and y,

$$r_{xy} = \frac{\sum_{i=1}^{M} (x_i - \overline{x})(y_i - \overline{y})}{\{\sum_{i=1}^{M} (x_i - \overline{x})^2\}^{1/2} \{\sum_{i=1}^{M} (y_i - \overline{y})^2\}^{1/2}}.$$
 (2)

Contrary to an expectation from the standard picture that a small binding energy leads to a large nuclear radius due to the extension of the wave function in the asymptotic region, the obtained correlation between S_{2n} and R_{rms} becomes not so strong, i.e. $r_{xy} = -0.801$. Therefore, a question arises, "which physical quantity is the explanatory variable to determine the nuclear radius?"

To answer this question, we plot the calculated results of the 12 parameter sets in Fig. 1 in the following manner. We take the x-axis for the energies of the p-orbit and the f-orbit in 30 F and the y-axis for the matter radius of 31 F. The symbols are assigned to different diffuseness parameters as a = 0.65, 0.70, and 0.75 fm, and the distance between the solid and open symbols connected by a dashed-line corresponds to the energy gap between the $p_{3/2}$ orbit and the $f_{7/2}$ orbit in 30 F. As seen from Fig. 1, it can be considered that the radius of 31 F is strongly correlated to the energy gap in 30 F.

In order to confirm the above investigation, we calculate PCC between the energy gap $\Delta \varepsilon$ in 30 F and the radius $R_{\rm rms}$ of 31 F. Here, the energy gap is defined as $\Delta \varepsilon \equiv \varepsilon(f) - \varepsilon(p)$, where $\varepsilon(f)$ and $\varepsilon(p)$ are the real part of the resonant energies of the lowest orbits of $7/2^-$ and $3/2^-$ in 30 F, respectively. The obtained PCC shows

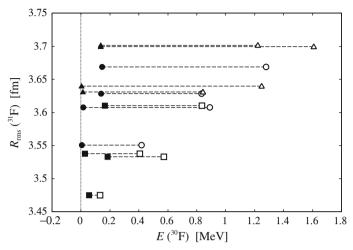


Fig. 1 Calculated matter radius of 31 F in different parameter sets of $\{V_0, a, r_0\}$ with the WS potential of 30 F. The x-axis shows the energy of 30 F, and the triangles, circles and squares are the results for the parameter sets of $r_0 = 0.65, 0.70$, and 0.75 fm, respectively. Dashed-line indicates the pair of 30 F energies for $p_{3/2}$ (solid) and $f_{7/2}$ (open) orbits. For detail, see text

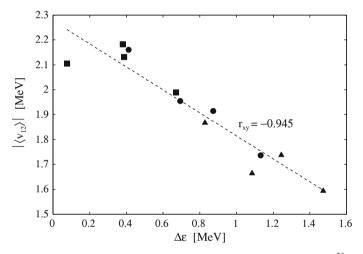


Fig. 2 The correlation between the absolute value of the neutron-neutron interaction $\langle \hat{v}_{12} \rangle$ in the ³¹F system and the energy gap $\Delta \varepsilon$ in 30 F. The definition of the triangles, circles and squares are the same with those applied in Fig. 1

very strong correlation as $r_{xy} = 0.935$, which is stronger than that obtained for S_{2n} and R_{rms} as $r_{xy} = -0.801$. Therefore, we consider the energy gap $\Delta \varepsilon$ is a key variable to indicate the structural change of ³¹F.

The energy gap also has a close relationship to the expectation value of the neutron numbers for the valence orbits in ³¹F. We calculate these numbers for the $p_{3/2}$ orbit (N(p)) and the $f_{7/2}$ orbit (N(f)), where N(f) and N(p) are added up to 2 including other orbits. In the largest radius case $R_{\rm rms} = 3.70$ fm, N(p) and N(f) are obtained as 1.63 and 0.19, respectively, i.e. ³¹F is p-wave dominant. On the other hand, in the smallest radius case $R_{\text{rms}} = 3.48$ fm, 31 F becomes f-wave dominant as N(p) = 0.38 and N(f) = 1.53. As shown in Fig. 3 of Ref. [6], N(p) and N(f) change inseparably and intersect each other at about $\Delta \varepsilon = 0.4$ MeV. Hence, it is confirmed that the energy gap $\Delta \varepsilon$ is related to the neutron numbers and also to the radius. However, even in the smallest radius case, a small separation energy $S_{2n} = 1.37$ MeV leads to the small radius $R_{\rm rms} = 3.48$ fm. Therefore, we consider a new mechanism to shrink the nuclear radius should occur in ³¹F as an "anti-halo" effect.

Next, we investigate how the energy gap affects to the pairing strength in the Borromean system, since the anti-halo effect has been discussed in terms of the pairing correlation [4,5]. Figure 2 shows the absolute value of $\langle \hat{v}_{12} \rangle$ in the Hamiltonian (1) with respect to the change of the energy gap $\Delta \epsilon$. As shown in Fig. 2, $\Delta\epsilon$ is strongly correlated to $|\langle \hat{v}_{12} \rangle|$, where PCC is obtained as $r_{xy} = -0.945$. It is noted that the correlation **20** Page 4 of 5

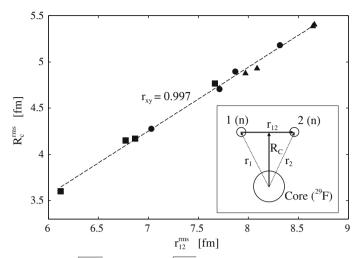


Fig. 3 The correlation between $r_{12}^{\text{rms}} \equiv \sqrt{\langle r_{12}^2 \rangle}$ and $R_c^{\text{rms}} \equiv \sqrt{\langle R_c^2 \rangle}$. The right-bottom panel shows the coordinate system of the $^{29}\text{F}+n+n$ three-body system in COSM (r_1 and r_2), and the definition of r_{12} and R_c . The definition of the triangles, circles and squares are the same with those applied in Fig. 1

between the separation energy S_{2n} and $|\langle \hat{v}_{12} \rangle|$ is very weak by considering PCC obtained as $r_{xy} = 0.660$. Combined with the result for the $\Delta \varepsilon$ dependence of N(p) and N(f), we can confirm that $|\langle \hat{v}_{12} \rangle|$, which can be an measure of the neutron-neutron correlation, increases for a small energy gap, and ^{31}F becomes f-wave dominant. As a result, the radius of ^{31}F shrinks even though the separation energy is still small.

To realize the strong correlations between $\Delta\epsilon$ and other quantities such as $R_{\rm rms}$, N(p), N(f), and $|\langle \hat{v}_{12} \rangle|$, and the shrinkage of the radius induced by $\Delta\epsilon$, there is another important condition that is the inversion of the valence orbits. $^{30}{\rm F}$ and $^{31}{\rm F}$ are considered to be placed on the "island of inversion", where the order of the single-particle orbits is inverted from the normal shell-model one. If the f-orbit ($f_{7/2}$) lies above the p-orbit ($f_{3/2}$), a competition occurs between the energy loss from the gap $\Delta\epsilon$ and the gain from the nucleon-nucleon interaction $|\langle \hat{v}_{12} \rangle|$, and consequently, $\Delta\epsilon$ becomes an explanatory variable to determine $|\langle \hat{v}_{12} \rangle|$, N(f), N(p), and $R_{\rm rms}$.

From the above discussion, the key ingredient that determines the nuclear radius of 31 F is considered to be the energy gap between the $p_{3/2}$ orbit and the $f_{7/2}$ orbit in 30 F. The formation of the halo structure in 31 F depends on whether the energy gap is large enough to overcome the energy loss of the nucleon-nucleon interaction part. This situation in general occurs for systems that the valence particle orbitals are inverted, i.e. a smaller angular momentum orbit lies below a higher angular momentum orbit. As the energy gap decreases, neutrons begin to occupy the orbit with the higher angular momentum, and the nucleon-nucleon interaction part gives larger contribution to overcome the energy loss from the gap. As a result, the radius becomes small even for a small separation energy. We consider this mechanism is a novel anti-halo effect.

Finally, in addition to the anti-halo effect, we investigate the possibility of the di-neutron like localization in 31 F. The correlation between the neutron-neutron distance $r_{12}^{\rm rms} \equiv \sqrt{\langle r_{12}^2 \rangle}$ and distance from the core to the center of mass of valence neutrons $R_c^{\rm rms} \equiv \sqrt{\langle R_c^2 \rangle}$ is shown in Fig. 3. If the ratio of $r_{12}^{\rm rms}$ to $R_c^{\rm rms}$ deviates from the line, for example, in the case of a small $r_{12}^{\rm rms}$ and a large $R_c^{\rm rms}$, the neutrons are supposed to have a di-neutron like localization. However, the correlation becomes very strong as $r_{xy} = 0.997$, which is almost the liner correlation, and it is difficult to find a possibility of the di-neutron like localization for the parameter sets of this calculation.

4 Summary

We study the structure of 31 F, which is the dripline nucleus of the fluorine isotopes, using COSM with GEM in a 29 F+n+n three-body model. Since there is an ambiguity to determine the 29 F+n potential, the strength parameter of WS can be changed by keeping the consistency for 31 Ne and 30 F. The correlation analysis is performed from the results using the 12 different parameter sets by keeping the consistency. As a result, we

found that a key ingredient to determine the nuclear radius of 31 F is the energy gap between the $p_{3/2}$ orbit and $f_{7/2}$ orbit in 30 F. This is considered as a general condition for other nuclei near the dripline, where the valence orbits are inverted, and it is expected that the cross over of the halo structure and the occurrence of the anti-halo effect can be determined by the energy gap between these inverted valence orbits.

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