α clustering in nuclei and its impact on the nuclear symmetry energy

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Nuclear symmetry energy is a fundamental quantity currently under intense investigation in both nuclear physics and astrophysics. The *softness* or *stiffness* of symmetry energy is still under debate and the extraction of symmetry energy from neutron skin thickness R_{skin} remains a challenge. Parity-violating measurements PREX and CREX provide important opportunities for constraining R_{skin} in ²⁰⁸Pb and ⁴⁸Ca. We investigate the occurrence of an α cluster at the surface of nuclei and its impact on the extraction of symmetry energy from R_{skin} . Our result indicates that the α -clustering probability in ²⁰⁸Pb is small and the extracted density slope of symmetry energy *L* is almost unaffected. In contrast, the α -clustering probability in ⁴⁸Ca is sizable and the corresponding correction to *L* should be taken into account. This correction progressively increases with the α -clustering probability, leading to a modification of the *L*- R_{skin} correlation, a fact that may have important implications in constraining nuclear symmetry energy.

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Introduction. The formation of compact clusters (e.g., α clusters) is an interesting feature of nuclear quantum manybody systems and plays an essential role in many important problems of astrophysics. The phenomena of α clustering are abundant in excited states of light nuclei close to the decay threshold [1]. One of the famous instances is the 3α -structure Hoyle state in ¹²C, which unlocks the puzzle in the production of heavy elements inside stars [2,3]. In contrast to light nuclei, the α -clustering problem in heavy nuclei is still not fully solved and the microscopic treatment of cluster dynamics beyond mean-field theory is a great challenge [4–10].

Recently, the PREX collaboration reported the measurement of parity-violating asymmetry A_{PV} and deduced a rather large neutron skin thickness R_{skin} in ²⁰⁸Pb (PREX-2) [11]. Thick skin in ²⁰⁸Pb suggests a very stiff symmetry energy in contrast to the previous constraints obtained from many other observations [12–17]. Very recently, the CREX collaboration has successfully conducted the parity-violating experiment in ⁴⁸Ca and deduced a thin R_{skin} [18], suggesting a soft symmetry energy. Much effort has been expended on attempting to reconcile these seemingly contradictory results. Special attention has been devoted to the problem of α clustering at the surface of nuclei that is expected to affect the density slope of symmetry energy L [19].

L is critical for understanding not only the structure of rare isotopes and the reaction mechanism of heavy-ion collisions, but also the structure and the composition of neutron stars [20–29]. By using the Hugenholtz–Van Hove (HVH) theorem, *L* can be decomposed in a unique way into kinetic energy, isoscalar potential, and isovector potential contributions [30,31]. As a fundamental relation for interacting

self-bound infinite Fermi systems, this theorem does not depend upon the precise nature of the interaction. While the kinetic energy and isoscalar potential contributions are relatively well constrained, the isovector potential contribution still has significant uncertainties. Owing to the existence of an isovector potential, more neutrons are pushed from the inner region of finite nuclei outwards to the surface region, and thus contribute to $R_{\rm skin}$. In this sense, *L* is related intrinsically to $R_{\rm skin}$.

The correlation between R_{skin} and L could be modified by the occurrence of α clusters [32]. This is because the α clusters may appear in the low-density region, i.e., the surface of a finite nucleus, and its impact progressively increases with the α -clustering probability [33–39]. Inside the core, α clustering is suppressed and its four nucleons are considered to move almost independently in a shell-model mean-field potential. The single-particle states are populated up to the Fermi energies of the neutrons or protons and a pairing correlation exists among the single-particle orbits. Pairing remains at high densities but the α cluster dissolves and its four nucleons (2n + 2p)turn into the single-particle motions forming the continuum of scattering states.

In this Letter, we address the question of whether α clustering at the surface of ²⁰⁸Pb and ⁴⁸Ca has a certain impact on R_{skin} and L. We use the quartetting wave-function approach (QWFA) to do so because it treats correctly both the intrinsic motion between four nucleons in the α cluster and the relative motion of the α cluster versus the core [34–37]. Strong closed-shell structure effects and complex derivative terms of the intrinsic wave function are properly taken into account in QWFA. The key quantity for clustering modification on the R_{skin} -L correlation is the α -cluster formation probability, which is quantitatively obtained by solving the coupled equations of a first-principles approach to nuclear many-body systems without adjusting any parameter.

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Intrinsic energy of an α cluster embedded in a nuclear medium. First, we simulate the α -cluster formation at the surface of a core nucleus by considering a low-density nuclear medium, in which the α -like four-nucleon correlations are described by the in-medium Schrödinger equation. The corresponding wave function of four nucleons is decomposed into a center-of-mass (c.m.) motion part $\Psi^{c.m.}$ and an intrinsic motion part φ^{intr} , which are coupled together with complex gradient terms. Such gradient terms are difficult to handle, but vanish in the case of a homogeneous nuclear medium. With the Jacobian momenta $\mathbf{p}_1 = \mathbf{P}/4 + \mathbf{k}/2 + \mathbf{k}_{12}$, $\mathbf{p}_2 = \mathbf{P}/4 + \mathbf{k}/2 - \mathbf{k}_{12}$, $\mathbf{p}_3 = \mathbf{P}/4 - \mathbf{k}/2 + \mathbf{k}_{34}$, $\mathbf{p}_4 = \mathbf{P}/4 - \mathbf{k}/2 - \mathbf{k}_{34}$, the in-medium wave equation for the intrinsic motion is reduced to [33]

$$\frac{\hbar^2}{2m} \Big[k^2 + 2k_{12}^2 + 2k_{34}^2 \Big] \varphi^{\text{intr}}(\mathbf{k}, \mathbf{k}_{12}, \mathbf{k}_{34}) \\ + \int \frac{d^3 k'}{(2\pi)^3} \frac{d^3 k'_{12}}{(2\pi)^3} \frac{d^3 k'_{34}}{(2\pi)^3} V_4 \varphi^{\text{intr}}(\mathbf{k}', \mathbf{k}'_{12}, \mathbf{k}'_{34}) \\ = (W^{\text{ext}} + W^{\text{intr}}) \varphi^{\text{intr}}(\mathbf{k}, \mathbf{k}_{12}, \mathbf{k}_{34}), \qquad (1)$$

where the centroid of the α cluster is considered to be at rest ($\mathbf{P} = 0$). V_4 is the effective in-medium interaction that contains the external mean field V_4^{ext} as well as the intrinsic nucleon-nucleon (NN) interaction modified by the Pauli blocking $V_4^{\text{intr}} = \Theta(p_1 - k_F)\Theta(p_2 - k_F)V_{\text{NN}}(\mathbf{p}_1, \mathbf{p}_2;$ $\mathbf{p}'_1, \mathbf{p}'_2)\delta(\mathbf{p}_3 - \mathbf{p}'_3)\delta(\mathbf{p}_4 - \mathbf{p}'_4) + 5$ permutations. The NN interaction is defined as a Gaussian form factor $V_{\text{NN}}(\mathbf{p}_1, \mathbf{p}_2;$ $\mathbf{p}'_1, \mathbf{p}'_2) = \lambda e^{-\frac{(\mathbf{p}_1 - \mathbf{p}_2)^2}{4\gamma^2}} e^{-\frac{(\mathbf{p}'_1 - \mathbf{p}'_2)^2}{4\gamma^2}} \delta(\mathbf{p}_1 + \mathbf{p}_2 - \mathbf{p}'_1 - \mathbf{p}'_2)$, where the potential parameters $\lambda = 1449.6$ MeV fm³ and $\gamma = 1.152 \text{ fm}^{-1}$ [33].

The minimum of the intrinsic energy W^{intr} has to be found for each density ρ with the Fermi-blocked Gaussian ansatz $\varphi^{\text{intr}}(\mathbf{p}_1, \mathbf{p}_2, \mathbf{p}_3, \mathbf{p}_4) = \frac{1}{\sqrt{N}} \varphi_{\tau_1}(\mathbf{p}_1) \varphi_{\tau_1}(\mathbf{p}_2) \varphi_{\tau_1}(\mathbf{p}_3) \varphi_{\tau_1}(\mathbf{p}_4) \delta(\mathbf{p}_1 + \mathbf{p}_2 + \mathbf{p}_3 + \mathbf{p}_4)$, where *N* is the normalization factor. The single nucleon wave function $\varphi_{\tau}(\mathbf{p})$ is given by $e^{-\frac{p^2}{2a}} \Theta[p - k_F]$ with the single variational parameter *a*. The minimum energy of a free α cluster is $W^{\text{intr}} = -28.3$ MeV at a = 0.535 fm⁻² [see Fig. 1(a)]. The intrinsic energy W^{intr} is shifted at finite density of the surrounding nuclear matter owing to the Pauli blocking. The bound state disappears and four nucleons become uncorrelated at the critical density $\rho_c = 0.03$ fm⁻³ [see Fig. 1(d)]. Note that the matter density distribution of the α cluster at the surface of a finite nucleus depends also on the surrounding density $\rho_{\alpha}(r, \rho) = 4(\frac{4a(\rho)}{3\pi})^{3/2}e^{-\frac{4a(\rho)}{3}r^2}$ by treating correctly both the energy shift and Pauli blocking effect.

Density evolution of an α cluster and formation probability in finite nuclei. The density evolution of an α cluster approaching the core nucleus is depicted in Fig. 2. The strong binding of the α cluster is gradually reduced by the energy shift due to Pauli blocking after it feels the tail of the core density. As shown in Fig. 2, the variational parameter *a* reflecting the size of the α cluster is decreased from 0.534 to 0.355 when it merges with the continuum of single-particle states. Eventually the α cluster dissolves and its four nucleons go over into single-particle states with pair correlations in the open shells on top of the core. Before that, the α cluster remains a



FIG. 1. The variation of intrinsic energies of an α cluster (a) in free space and [(b)–(d)] in homogeneous nuclear matters. (d) A critical transition occurs at $\rho_c = 0.03$ fm⁻³, where the α cluster dissolves and its four nucleons become uncorrelated.

relatively compact entity with a small extension even up to the critical density ρ_c (see Fig. 2).

The c.m. motion of an α cluster is introduced as a dynamical collective degree of freedom, which simplifies the treatment of correlated nuclear systems beyond the mean-field approximation. By separating the intrinsic motion from the c.m. motion, the c.m. wave function of the α cluster follows the equation [36,37]

$$-\frac{\hbar^{2}}{2Am}\nabla_{R}^{2}\Psi^{\text{c.m.}}(\mathbf{R}) - \frac{\hbar^{2}}{Am}\int ds_{j}\varphi^{\text{intr},*}(\mathbf{R},\mathbf{s}_{j})$$

$$\times [\nabla_{R}\varphi^{\text{intr}}(\mathbf{R},\mathbf{s}_{j})][\nabla_{R}\Psi^{\text{c.m.}}(\mathbf{R})]$$

$$-\frac{\hbar^{2}}{2Am}\int ds_{j}\varphi^{\text{intr},*}(\mathbf{R},\mathbf{s}_{j})[\nabla_{R}^{2}\varphi^{\text{intr}}(\mathbf{R},\mathbf{s}_{j})]\Psi^{\text{c.m.}}(\mathbf{R})$$

$$+\int dR'W(\mathbf{R},\mathbf{R}')\Psi^{\text{c.m.}}(\mathbf{R}') = E\Psi^{\text{c.m.}}(\mathbf{R}), \qquad (2)$$



FIG. 2. Density evolution of an α cluster approaching the core nucleus. The width parameter $a(\rho)$ and density distribution of an α cluster $\rho_{\alpha}(r, \rho)/4$ are shown for densities $\rho = 0.001, 0.01, 0.02$, and 0.03 fm⁻³, respectively.



FIG. 3. The normalized c.m. wave function of four nucleons forming the α cluster. The wave function in the range of $0 < R < R_c$ represents the c.m. wave function of four uncorrelated nucleons after dissolution. Only at the surface region with $R > R_c$ does the α cluster appear and the c.m. wave function with $R > R_c$ corresponds to the formed α cluster, as marked by the green color. The pocket region in the c.m. effective potential is denoted in the small panel.

where the second and third terms are complex derivative terms and no previous investigations of such terms have been performed. It can be strictly proved that the second term vanishes if the number of nucleons (A = 4) embedded in the medium is conserved. In contrast, the third term is nontrivial and is rather difficult to solve (ninefold integral). For the first time, we take this derivative term into account in QWFA and find that this term does affect the final α -cluster formation probability. The fourth term is the effective potential describing the c.m. motion of the α cluster under the influence of Pauli blocking with the surrounding medium. The inner c.m. effective potential $W(R < R_c)$ (R_c is the critical radius corresponding to ρ_c) is constructed from the shell-model wave functions of four nucleons forming the α cluster. Note that only states near the Fermi energy can form an α -like cluster because these shell-model states extend to the low-density regions. The inner effective potential $W(R < R_c)$ joins with the outer one $W(R > R_c) = W(R)^{\text{ext}} + W(R)^{\text{intr}}$ at $R = R_c$. An important feature of W(R) is that a pocket is formed on the surface region (see the small panel in Fig. 3), resulting from the competition between strong nuclear force attraction and repulsive Pauli blocking. The pocket plays an essential role in the formation of the α cluster at the surface of a core nucleus. As seen in Fig. 3, the normalized c.m. wave function shows a small peak around the pocket region. This is in agreement with the microscopic calculations on α clustering in Refs. [40–42]. By integrating the c.m. wave function from the critical radius R_c to infinity, the α -cluster formation probability can be microscopically obtained $P_{\alpha} = \int_{R>R_c} d^3 R |\Psi^{\text{c.m.}}R|^2 [36,37].$ We go beyond the Thomas-Fermi approximation by taking the closed-shell structure effects into account. The α -cluster formation probability is expected to vary dramatically across the major shell closures. Indeed, it is found that the α clustering in doubly magic nuclei such as ⁴⁰Ca, ¹³²Sn, and ²⁰⁸Pb is significantly hindered by shell effects (see Fig. 4). The real*istic* α -cluster formation probability in ²⁰⁸Pb is rather small,



FIG. 4. Comparison of α -cluster formation probabilities in doubly magic nuclei ⁴⁰Ca, ¹³²Sn, and ²⁰⁸Pb and in nuclei ⁴⁴Ti, ¹³⁶Te, and ²¹²Po.

 $P_{\alpha} = 9.3 \times 10^{-3}$. This is quite different from their neighbors 44 Ti, 136 Te, and 212 Po where enhanced α -cluster formation probabilities are found by using exactly the same QWFA formulism.

Impact of α clustering on R_{skin} and L. A direct relationship between L and the underlying single-nucleon potential $V_{n/p}(\rho, \delta, k) = V_0(\rho, k) \pm V_{sym}(\rho, k)\delta$ is revealed by the HVH theorem [30,31]. The advantage of this strict relationship is that it can be used to determine L in a fully transparent way. At the saturation density ρ_0 , L can be reformulated by using the effective mass m^* ,

$$L = \frac{2}{3}t(k_F^0) + \frac{3}{2}V_{\text{sym}}(\rho_0, k_F^0) + \frac{\partial V_{\text{sym}}(\rho_0, k)}{\partial k}\Big|_{k_F^0} k_F^0, \quad (3)$$

where the first term $L(1) = \frac{\hbar^2 k_F^2}{3m^*}$ denotes the contributions from kinetic energy and isoscalar potential [30]. For the nucleon effective mass, we adopt the value of $\frac{m^*}{m} = 0.70 \pm$ 0.05 widely used in the literature (see, e.g., Ref. [43]). The isovector potential $V_{\text{sym}}(\rho_0, k)$ can be deduced from the real part of global optical potentials, which is basically parametrized in the Woods-Saxon (WS) form, i.e., V(r) = $-V_0[1 \pm \kappa(\frac{N-Z}{A})]/[1 + \exp(\frac{r-R_0A^{1/3}}{a})]$ ("+" for protons and "-" for neutrons). The second term L(2) in Eq. (3) is determined by the product of the strength of the WS potential V_0 and the isovector parameter κ , i.e., $L(2) = \frac{3}{2}V_{\text{sym}}(\rho_0, k_F^0) =$ $\frac{3}{2}\kappa \cdot V_0$. The WS potential does not have an explicit energy (or momentum) dependence. From the global optical potential (GOP) constrained by nuclear reaction data [30], the energy dependence of the isovector potential is found to have a linear form $V_{\text{sym}}(\rho_0, k) = 22.75 - 0.21E(k)$ [30]. So the third term L(3) in Eq. (3) is negative because of the decreasing isovector potential with increasing energy.

We use the same "QWFA" WS global optical potential to determine (1) shell-model states and densities in ²⁰⁸Pb and ⁴⁸Ca, together with the Coulomb potential + *ls* coupling, and (2) density slope of symmetry energy *L* by using the HVH theorem. The QWFA parametrization is found to reproduce well the α -cluster decay half-lives around doubly magic nuclei ²⁰⁸Pb and ¹⁰⁰Sn [36,37]. The neutron skin thickness $R_{skin} = r_n^{rms} - r_p^{rms}$ is calculated directly from shell-model density distributions. With an explicit α -cluster degree of freedom, the rms radius is given by $r^{rms} = \{\int r^2 [\rho^{cluster}(r) + \rho^{core}(r)] d^3r\}^{1/2}$,



FIG. 5. Correlation between L and $R_{\rm skin}$ with and without α clustering in ²⁰⁸Pb. In the case of α clustering, a large amount of α -cluster formation probability is assumed ($P_{\alpha} = 1$).

where ρ^{core} is the density distribution of protons or neutrons in the core. The density distribution of two neutrons or two protons forming the α cluster is

$$\rho^{\text{cluster}}(r) = 2 \int_{R < R_c} d^3 R[|\Psi^{\text{c.m.}}(\mathbf{R})|^2 \rho(\mathbf{r})]$$

+ $\frac{1}{2} \int_{R > R_c} d^3 R[|\Psi^{\text{c.m.}}(\mathbf{R})|^2 \rho_{\alpha}(\mathbf{r} - \mathbf{R}; \mathbf{R})], \quad (4)$

where the α -cluster formation at the surface of core nucleus $(R \ge R_c)$ is taken into account in the second integral and the spatial extension of the formed α cluster is well described by $\rho_{\alpha}(\mathbf{r} - \mathbf{R}; \mathbf{R})$. We assume that the neutron skin thicknesses given by PREX-2 and CREX are all measured quantities. Figure 5 shows the correlation between L and R_{skin} with and without α clustering for ²⁰⁸Pb. L increases with the increasing R_{skin} and the *L*- R_{skin} correlation is almost linear. We have checked the L- R_{skin} correlation by using different WS parametrizations [44] and found that this behavior is general. As shown in Fig. 5, the L- R_{skin} correlation is modified significantly by assuming a large amount of formation probability $(P_{\alpha} = 1)$. However, the *realistic* α -cluster formation probability in 208 Pb is quite small, and thus its influence on L is negligible. By considering the *realistic* α -cluster formation probability in 208 Pb, the L value extracted from the PREX-2 data is 75.3 MeV (see Table I). The uncertainties of all terms contributing to this L value are considered. The uncertainty in L(1) is due to the effective mass m^* . With the $m^*/m =$ 0.70 ± 0.05 we adopted, an error bar of +2.8/-2.4 MeV is obtained. Since the R_{skin} data of PREX-2 have a large error bar, it is no surprise that there is a large error bar associated with the L(2) term. The error bar associated with the L(3) term is obtained by the world data on nucleon-nucleus scatterings, (p, n) charge exchange reactions, and single-particle energies

TABLE I. The extracted density slope parameter *L* by considering α clustering at the surface of ²⁰⁸Pb (PREX-2) and ⁴⁸Ca (CREX). *The correction of *L* due to α clustering for the lower and upper limits of $R_{\rm skin}$ is 100% and 8%, respectively.

Nuclei	<i>R</i> _{skin} (fm)	L (MeV) no α cluster	P_{α}	L (MeV) with α cluster
²⁰⁸ Pb	0.283 ± 0.071	$75.2^{+24.3}_{-24.5}$	$9.3 imes 10^{-3}$	$75.3^{+24.3}_{-24.6}$
⁴⁸ Ca	0.121 ± 0.050	$13.2^{+25.4}_{-24.9}$	$7.3 imes 10^{-2}$	$15.0^{+25.6}_{-25.0}$
*	0.071 (lower)	1.7		3.4
*	0.171 (upper)	24.8		26.8

of bound states [30]. Put all together, the error bar of *L* is approximately 24 MeV. In contrast to ²⁰⁸Pb, the *realistic* α -cluster formation probability P_{α} in ⁴⁸Ca is found to be 7.3×10^{-2} , which is much larger than that in ²⁰⁸Pb. Thus the impact of α clustering in ⁴⁸Ca on *L* cannot be ignored. The extracted *L* value with an error bar from the CREX data is $15.0^{+25.6}_{-25.0}$ MeV and the correction due to α clustering is of the order of 14%. This correction progressively increases with the α -cluster formation probability P_{α} , which could be close to unity if the contributing shell-model orbits are rather similar, especially for self-conjugate nuclei such as ⁴⁴Ti in Fig. 4.

Conclusion. α -clustering survives at the surface of heavy nuclei, which is relevant to the neutron skin thickness of heavy nuclei. The latter is a precise tool in constraining the density slope of nuclear symmetry energy. The impact of α clustering on the neutron skin thickness depends closely on the amount of the formation probability. We emphasize that the approach presented here to calculate formation probability is based on a first-principles approach to nuclear many-body systems. A proper treatment of derivative terms of intrinsic wave function has been performed and the spatial extension of the α cluster has been considered to better account for the correlation between L and R_{skin} . The present analysis shows that the L values deduced from PREX-2 and CREX experiments are not consistent with each other, even with the α -clustering effect included. We expect that a better account of the model dependence in extracting R_{skin} from parity-violating asymmetry $A_{\rm PV}$ will further improve the estimation of L. Moreover, state-of-art approaches can be applied to describe shell-model states and nuclear densities and the Gaussian ansatz used in the variational calculations can be improved. An exact solution of the four-nucleon correlation near the critical density by including both self-energy corrections and Pauli blocking should be tackled in the future.

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