Searching for $4\alpha$ linear-chain structure in excited states of $^{16}\text{O}$ with a covariant density functional theory

J. M. Yao, 1, 2 N. Itagaki, 3 and J. Meng 4, 5, 6

1Department of Physics, Tohoku University, Sendai 980-8578, Japan
2School of Physical Science and Technology, Southwest University, Chongqing 400715, China
3Yukawa Institute for Theoretical Physics, Kyoto University, Kyoto 606-8502, Japan
4State Key Laboratory of Nuclear Physics and Technology, School of Physics, Peking University, Beijing 100871, China
5School of Physics and Nuclear Energy Engineering, Beihang University, Beijing 100191, China
6Department of Physics, University of Stellenbosch, Stellenbosch, South Africa

A study of $4\alpha$ linear-chain structure (LCS) in high-lying collective excitation states of $^{16}\text{O}$ with a covariant density functional theory is presented. The low-spin states are obtained by configuration mixing of particle-number and angular-momentum projected quadrupole deformed mean-field states with generator coordinate method (GCM). The high-spin states are determined by cranking calculations. These two calculations are based on the same energy density functional PC-PK1. We have found the LCS candidate in both high-lying low-spin GCM states and cranking high-spin states with similar moment of inertia and band-head energy, which are estimated to be around 0.11 MeV and 30 MeV, respectively. The intrinsic configuration is considered to be the one that $4\alpha$ clusters stay along a common axis and nucleons occupy the $(s)^4(p)^4(d)^4(f)^4$ configurations in a nonlocal way. Moreover, our results indicate that the spin and orbital angular momenta of all nucleons are parallel in the LCS states but the sum of spin-orbit splitting energies turns out to be much smaller than that of shell-like state. Besides, our fully microscopic GCM calculation has reproduced the excitation energies and $B(E2)$ values rather well for the rotational band built on the second $0^+$ state which has been previously considered to have $^{12}\text{C}+\alpha$ structure, and the dominant configuration turns out to be four $\alpha$ clusters with “kite”-like shape.

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I. INTRODUCTION

The excited states close to particle emission threshold in doubly-magic $^{16}\text{O}$ have been of much interest due to the formation of $\alpha$ clustering structure, which is different from the spherical ground state with shell-like picture. In particular, the possible existence of $4\alpha$ linear-chain structure (LCS) in highly excited states of $^{16}\text{O}$ has been under intensive discussion. About half-century ago, Chevallier et al. observed resonant $2^+, 4^+$, and $6^+$ states in the reaction $^{12}\text{C}(\alpha, \text{Be})\text{Be}$ and proposed that these states may correspond to the rotating states with $4\alpha$ LCS. The moment of inertia (MOI) was estimated as $h^2/(2J) = 64$ keV with band-head excitation energy of 16.8 MeV [1]. This proposal was supported by the analysis of decay widths for the states [2]. Later, Freer et al. performed the $^{12}\text{C}(^{16}\text{O}, 4\alpha)^{12}\text{C}$ reaction, and obtained a slightly smaller MOI ($96 \pm 20$ keV) with bandhead excitation energy of $17.0 \pm 0.7$ MeV [3]. The calculation by Bauhoff et al.[4] using the Brink’s $\alpha$ cluster model supported the existence of $4\alpha$ LCS state in low-spin states of $^{16}\text{O}$, which predicted the MOI to be 64 keV and the band-head excitation energy of 16.3 MeV, in excellent agreement with the data [1]. Most recently, a new measurement of the $^{12}\text{C}(\alpha, \text{Be})\text{Be}$ excitation function was carried out. Unfortunately, this experiment did not provide any evidence to support the existence of $4\alpha$ LCS in $^{16}\text{O}$ [5].

The formation of cluster structure in finite quantum many-body systems in itself is an interesting phenomenon. In nuclear physics, the cluster structure is essential to understand many problems of nuclear structure and reactions. A fully microscopic understanding of cluster formation necessitates the treatment of the individual nucleons as the fundamental ingredients, with the clusters and their properties emerging automatically from the calculations, instead of assuming a priori certain geometrical arrangements for the clusters. Self-consistent mean-field approaches with all nucleons treated on the same footing provide us good tools for understanding these phenomena from this aspect. The previous studies in this context have shown that the nucleons are prone to form cluster structure in the nuclear system with either high excitation energy and high spin with large deformation [6, 7], deep confining nuclear potential [8, 9], or expansion with low density [10].

In the recent decade, several studies based on selfconsistent mean-field approaches have been carried out for the $4\alpha$ LCS in $^{16}\text{O}$. The relativistic mean-field (RMF) calculations with a constraint on nuclear quadrupole moment demonstrated the existence of $4\alpha$ LCS in hyperdeformed mean-field state with deformation parameter $\beta \approx 3.6$ and excitation energy of 37 MeV [11, 12]. It can be understood that the degeneracy of single-nucleon levels is reduced by deformation since the spherical symmetry is lost, but can be recovered at some specific deformations, for example, with the long-short axis ratio around 2:1 (superdeformed shape) or 3:1 (hyperdeformed shape), which favors the formation of clusters [13]. The stability...
of this state against quadrupole shape fluctuation in \(^{16}\text{O}\) has been studied with a generator coordinate method implemented with particle-number and angular-momentum projection (GCM+PNAMP) based on a non-relativistic Skyrme-Hartree-Fock (SHF)+BCS calculation with SLy4 force \([14]\). It was found that the \(0^+\) state with dominated \(8p-8h\) character and excitation energy of 32 MeV is close to the \(4\alpha\) LCS. A more recent calculation by Ichikawa \textit{et al.} \([7]\) using a cranking SHF method suggested that a rapidly rotating \(^{16}\text{O}\) can deform into the \(4\alpha\) LCS in the region of angular momentum \((13 - 18h)\). The MOI was estimated to be around 60 – 80 keV with a very high band-head energy 38 MeV. The formation of LCS in high-spin state was regarded as a result of the competition between nuclear attractive and centrifugal forces due to rapid rotation. It means that even without taking into account the orthogonality condition to the low-lying state, which can enhance the stability of LCS and it was considered in Refs. \([15, 16]\), one finds a region of angular momenta for the formation of LCS in high-spin states.

In recent years, the structure of the proposed \(4\alpha\) LCS rotational states in Ref. \([1]\) has been mainly interpreted in two different ways. A study of the structure and scattering of \(\alpha+^{12}\text{C}\) system using a double folding model in the coupled-channel method suggests that the rotational states have the \(^{12}\text{C}(0^+_{\alpha})+\alpha\) cluster structure \([17]\). This picture is also supported by the most recent study using the GCM+AMP method based on an extended \(^{12}\text{C}(AMD)+\alpha\) cluster model, which proposes the \(^{12}\text{C}(0^+_{\alpha})+\alpha\) state \((J^\pi = 0^+)\) to be a \(4\alpha\) condensed gas-like state with excitation energy of 17 MeV and rms charge radius of 3.8 fm \([18]\). On the other hand, most recently, the \(4\alpha\) LCS in \(^{16}\text{O}\) was investigated with the Brink model wave function by Suhara \textit{et al.} \([19]\), suggesting that the \(4\alpha\) LCS has the one-dimensional \(\alpha\) condensate character, where the \(\alpha\) clusters are trapped into a one-dimensional potential in a nonlocal manner, like a gas.

The aim of this work is to search for the \(4\alpha\) LCS in \(^{16}\text{O}\) with a covariant density functional theory (CDFT), which has already achieved great success in describing variety aspects in nuclear physics \([20-23]\). On one hand, it has been pointed out in Ref. \([8]\) that relativistic energy density functionals (EDF) are characterized by deep single-nucleon potentials and therefore are prone to predict the occurrence of much more pronounced cluster structures in nuclear ground state than the non-relativistic EDF. On the other hand, most clustering states appear in light nuclei and are deformed in the intrinsic frame. The effects from shape fluctuations and restoration of rotational symmetry need to be examined. Moreover, the cluster structure in nuclear excited states has not been studied with a relativistic functional. Therefore, the search for the \(4\alpha\) LCS in both low-spin and high-spin excited states of \(^{16}\text{O}\) and the investigation of the structure properties of these exotic states based on a relativistic EDF are very interesting. To this end, both the GCM+PNAMP and cranking methods on top of the CDFT are adopted.

The paper is organized as follows. In Sec. II, we introduce the basic formulae of the CDFT and its extension for nuclear collective excitation states briefly. The results from both GCM calculation for nuclear low-spin states and cranking calculation for nuclear high-spin states based on the CDFT are presented and discussed in Sec. III. Finally, a summary and outlook are given in Sec. IV.

II. THEORETICAL FRAMEWORK

The starting point of the point-coupling type of CDFT is an EDF which has the following form \([24, 25]\),

\[
E_{\text{DF}}[\rho_i, \nabla \rho_i, j_i^\mu, \nabla j_i^\mu] = \text{Tr}[\rho \cdot p + \beta m] + \int dr \left( \frac{\alpha_S}{2} \frac{\rho^2}{\rho S} + \frac{\beta_S}{3} \rho^3 + \frac{\gamma_S}{4} \rho^4 + \frac{\delta_S}{5} \rho S \nabla \rho S \right) + \frac{\alpha_V}{2} j_i^\mu j_i^\mu + \frac{\gamma}{4} (j_i^\mu j_i^\mu)^2 + \frac{\delta_V}{2} j_i^\mu j_i^\mu \nabla (j_i^\mu j_i^\mu) + \frac{1}{4} \hat{F}_{\mu \nu} F^{\mu \nu} - F_{0 \mu} \partial_0 A_\mu + e \left( 1 - \frac{3}{r_0} \frac{\gamma}{2} j_i^\mu A_\mu \right),
\]

where densities \(\rho_i\) and currents \(j_i^\mu\) are bilinear combinations of Dirac spinors, namely \(\psi \Gamma_i \psi\) with \(i = S, V, TV\) representing the symmetry of the coupling. The subscript \(S\) stands for isoscalar-scalar (\(\Gamma_S = 1\)), \(V\) for isoscalar-vector (\(\Gamma_V = \gamma\)), and \(TV\) for isovector-vector (\(\Gamma_{TV} = \gamma^\mu t_3\)) type of coupling characterized by their transformation properties in isospin and in space-time. \(A^\mu\) is the four-component electromagnetic field. The coupling constants \(\alpha, \beta, \gamma, \delta\) are determined in the optimization of the EDF for the properties of several finite nuclei and/or nuclear matter \([24, 25]\).

In the following, we will introduce the extensions of the CDFT for low-spin and high-spin states separately in a brief way. More detailed description can be found in Refs. \([26-30]\) and Refs. \([31, 32]\) respectively.

A. The GCM calculation for low-spin states

The wave function of nuclear low-spin state is given by the superposition of a set of both particle-number and angular-momentum projected (PNAMP) quadrupole deformed mean-field states in the framework of GCM \([30]\),

\[
|JNZ; \alpha\rangle = \sum_{q,K} f^{JNZK}_\alpha (q) \hat{P}_{MN} \hat{P}_{MK} \hat{P}^N \hat{P}^Z |q(\beta, \gamma)\rangle,
\]

where \(\alpha = 1, 2, \ldots\) distinguishes different collective states with the same angular momentum \(J\). The operators \(\hat{P}^N\), \(\hat{P}^Z\), and \(\hat{P}_{MK}\) project onto good neutron and proton numbers and onto good angular momentum.
mean-field states $|q(\beta, \gamma)\rangle$ are Slater determinants of single-(quasi)particle states from the RMF+BCS calculation with constraints on the mass quadrupole moments

$$Q_{20} = \sqrt{\frac{5}{16\pi}}(2z^2 - x^2 - y^2)$$

and

$$Q_{22} = \sqrt{\frac{15}{32\pi}}(x^2 - y^2),$$

where the deformation parameters $\beta, \gamma$ are related to the quadrupole moments by $\beta = \frac{4\pi}{3AR^2}Q_{20}, \gamma = \tan^{-1}\left(\sqrt{\frac{Q_{22}}{Q_{20}}}\right)$, respectively, with $R = 1.2A^{1/3}$ and $A$ being the mass number. For simplicity, the mean-field states $|q(\beta, \gamma)\rangle$ are restricted to axially deformed, namely, $\gamma = 0^\circ$ (prolate) and $180^\circ$ (oblate). In this case, $K = 0$, and the $f_{\alpha JNK}^{(sN)}$ is replaced with $f_{\alpha JNK}^{(s)}$. Moreover, $q(\beta, \gamma)$ is abbreviated with $\beta$ subsequently.

Minimization of nuclear total energy with respect to the coefficient $f_{\alpha JNK}^{(s)}$ leads to the Hill-Wheeler-Griffin (HWG) equation [33],

$$\sum_{\beta'} [\mathcal{H}^{(\beta, \beta')}_a] - E^{(\beta, \beta')}_a f_{\alpha JNK}^{(\beta', \beta)}(\beta') = 0, \quad (3)$$

where $\mathcal{N}^{(\beta, \beta')}_a$ and $\mathcal{H}^{(\beta, \beta')}_a$ are the norm kernel and the energy kernel, respectively. The solution of HWG equation provides the energy spectrum and all the information needed for calculating the electric multipole transition strengths in low-spin excited states.

\section*{B. The cranking RMF calculation for high-spin states}

The GCM+PNAMP method can provide nuclear excited states with good quantum numbers, which are essential for spectroscopic study. However, it is currently limited to low-spin states due to the computation difficulty. To obtain nuclear high-spin states, we are restricted to the semiclassical cranking method on top of the CDFT, in which the nucleus is cranked along $x$-axis with a constant rotational frequency $\omega$. The wave function of single-particle (s.p.) state is the solution of the Dirac equation in body-fixed rotating frame, which turns out to have the following form [23, 34],

$$(h_0 - \omega J_x)\psi_k(r, s, t) = \varepsilon_k \psi_k(r, s, t), \quad (4)$$

where the single-particle Hamiltonian in non-rotating frame $h_0$ is

$$h_0 = \alpha \cdot [-i\nabla - \mathbf{V}(r)] + \beta (m + S(r)) + V_0(r). \quad (5)$$

The $\varepsilon_k$ and $\psi_k$ are the energy and wave function of single-particle (s.p.) state in rotating frame, respectively. $m$ is bare nucleon mass. The $S(r), V_0(r)$ represent the scalar potential and vector potentials. The $\mathbf{V}(r)$ is the space-like (time-odd) component of vector potential, which is often called nuclear magnetism [20] and is nonzero only in time-reversal invariance violated systems. In the cranking RMF approach for a rapidly rotating nucleus, the Coriolis term $\omega J_z$ violates the time-reversal invariance in the intrinsic frame and therefore generates the nonzero nuclear magnetism term $\mathbf{V}(r)$.

\section*{C. Expansion of Dirac spinor on a basis}

The Dirac equation in both deformation constrained RMF+BCS and cranking RMF calculations is solved by expanding the Dirac spinor $\psi_k$ on a three-dimensional harmonic oscillator (HO) basis in Cartesian coordinates,

$$\psi_k(r, s, t) = \left(\begin{array}{c} f_k(r, s) \\ ig_k(r, s) \end{array} \right) \chi_k(t), \quad (6)$$

where $\chi_k(t)$ is the isospin part and

$$f_k(r, s) = \sum_\alpha f_{\alpha k}(\alpha) + \sum_{\bar{\alpha}} f_{\bar{\alpha}k}(\bar{\alpha}), \quad (7a)$$

$$g_k(r, s) = \sum_\alpha g_{\alpha k}(\alpha) + \sum_{\bar{\alpha}} g_{\bar{\alpha}k}(\bar{\alpha}). \quad (7b)$$

The HO basis $\{\alpha, \bar{\alpha}\}$ are chosen as eigenstates of the $x$-simplex operator $\hat{S}_x = \hat{P}e^{-i\pi J_z}$ with positive and negative eigenvalues

$$\begin{cases} |\alpha\rangle = |n_x n_y n_z\rangle \frac{i^{n_x}}{\sqrt{2}} \left(1 + (-1)^{n_x+1}ight), \\ |\bar{\alpha}\rangle = |n_x n_y n_z\rangle \frac{(-1)^{n_x+n_y+1}}{\sqrt{2}} \left(-1 - (-1)^{n_x}\right), \end{cases} \quad (8)$$

where $|n_x n_y n_z\rangle$ is the space part of the three-dimensional HO wave function. The phase factor $i^{n_x}$ is chosen in order to have a real matrix elements for the Dirac equation [35].

In the deformation constrained RMF+BCS calculation for generating mean-field reference states as inputs of GCM calculation, the symmetries associated with parity, $x$-simplex, and time-reversal invariance are imposed. In this case, one has to expand the Dirac spinor only in half of the full basis with positive or negative simplex eigenvalue. Therefore, we expand the large $f_k$ and small $g_k$ component in Eq. (7) only on the basis with positive and negative $x$-simplex values, respectively. In the cranking RMF calculation for high-spin states, we adopt the tilde-cranking code developed in Refs. [23, 31, 32] and restrict the rotation along $x$-axis. In this code, only the symmetries associated with parity ($P$) and the combination of time reversal ($T$) and space reflection with respect to $x-z$ plane ($\hat{P}y : y \rightarrow -y$) are imposed. Since the time-reversal invariance is violated by the coriolis term, the time-reversal partner states are not degenerate in energy. The Dirac equation has to be solved in the full basis with both positive and negative $x$-simplex values, cf. Eq. (7). More details are also introduced in the review paper [23].

\section*{III. RESULTS AND DISCUSSIONS}

In both the deformation constrained RMF+BCS and cranking RMF calculations, 12 major HO shells are
adopted, which turns out to give difference in total energy within 1 MeV compared with the value by 14 major shells for the configurations with deformation parameter $\beta$ up to 4.0. We note that the GCM+PNAMP calculation with 14 major HO shells for the expansion of Dirac spinor is very time-consuming. In particular, as we will see in the results, this energy difference is marginal and will not have much influence on our conclusions, compared with the excitation energy of LCS candidate states. There is no parameters in our study other than those in the EDF, for which the relativistic point-coupling parametrization PC-PK1 [25] is used throughout this work. In the deformation constrained RMF+BCS calculation, a density-independent $\delta$ force implemented with an energy-dependent smooth cutoff factor [36] is adopted in the same way as the PC-PK1 was parameterized. In the cranking RMF calculation, pairing correlations between nucleons are neglected due to the anti-pairing effect of the Coriolis term.

A. 4$\alpha$ LCS in low-spin states

Figure 1(b) displays the mean-field and both particle-number and angular-momentum (PNAMP) projected energy as a function of the intrinsic quadrupole deformation $\beta$ in $^{16}$O. To search for the LCS in the mean-field state, we also plot the density profiles for some configurations on the curve. It is shown that there is a shoulder on the energy curve around $\beta = 3.6\alpha$, at which an evident 4$\alpha$ LCS is found. The excitation energy of this state is 39.1 MeV. After restoration of rotational symmetry, this energy is reduced to 31.5 MeV. The configuration mixing calculation predicts the excitation energy of $0^+$ state with the 4$\alpha$ LCS to be 29.6 MeV. The result is consistent with the value of 32.0 MeV obtained by the previous GCM+PN1DAMP calculation using the SLy4 force [14], which suggests a 8p-8h character for this state. However, these values are still much higher than the suggested value of 16.8 MeV in Ref. [1]. Besides the rotational band with 4$\alpha$ LCS (with $\beta \in [2.8, 4.0]$), an oblate deformed rotational band (with $\beta \in [-1.0, -0.5]$), together with a highly prolate deformed rotational band (with $\beta \in [1.0, 2.0]$), which is similarity to the “kite” structure found in Refs. [4, 11], is displayed in Fig. 1. Moreover, when the deformation parameter $\beta$ is increased up to $\beta = 4.8\alpha$, a structure of $^8$Be+$^8$Be appears in $^{16}$O, which is beyond the scope of the present study.

It is basically considered that for the $\alpha$ gas-like state, all of the $\alpha$ clusters occupy the same $S$ orbital, which has a spatially extended distribution [37]. This picture is well described by the THSR (Tohsaki, Horiuchi, Schuck and Röpke) wave function [38]. If $\alpha$ clusters are formed, the spin-orbit interaction energy should be exactly zero. From this point of view, the spin-orbit interaction energy may provide us a way to distinguish the shell-like and cluster-like structure. Fig. 1(a) displays the spin-orbit interaction energy

$$SO1 = -\sum_k v_{k}^{2} \langle k | V_{s.o.} | k \rangle,$$

and the sum of the matrix elements with its absolute

\begin{figure}[h]
\centering
\includegraphics[width=\textwidth]{fig1.png}
\caption{(Color online) (a) Spin-orbit interaction energy $SO1$ (9) and sum of spin-orbit splitting energy $SO2$ (10) and (b) total energy of the mean-field and PNAMP states (normalized to the energy of $0^+_1$ state) in $^{16}$O as a function of the intrinsic quadrupole deformation $\beta$. The horizontal short lines with bullets are the GCM solutions. Only the states with similar dominated configurations (oblate, “kite”, and LCS respectively) are plotted and placed at their “average” deformation $\beta = \frac{\sum^4_{\beta} | g_{\beta}^{O} |^{2} \beta }{4}$, cf.(12). The insets are the contours of intrinsic total density on $y$-$z$ plane (at $x = 0.3$ fm) for some typical configurations along the curve.}
\end{figure}

\begin{figure}[h]
\centering
\includegraphics[width=\textwidth]{fig2.png}
\caption{(Color online) Comparison of the spin-orbit interaction matrix element $\langle k | V_{s.o.} | k \rangle$, cf. (11), for the $k$-th occupied s.p. state in the mean-field states of $^{16}$O with $\beta = 0.0$ and $\beta = 3.6\alpha$.}
\end{figure}
values
\[ SO_2 = \sum_k v_k^2 |\langle k|V_{s.o.}|k \rangle|, \tag{10} \]
for each mean-field state \(|\beta\rangle\), where a “−” sign is introduced in defining \(SO_1\) to have positive values because the spin-orbit interaction is attractive, \(v_k^2\) is the occupation probability of the \(k\)-th s.p. state in the mean-field state \(|\beta\rangle\) and the matrix element is given by [20, 39]
\[ \langle k|V_{s.o.}|k \rangle = \langle k|\frac{1}{4m^2}(\nabla V_{\text{fs}}) \cdot (p \times \sigma)|k \rangle, \tag{11} \]
with \(V_{\text{fs}} = -\frac{m}{m_{\text{eff}}} (V_0 - S)(r)\), and \(m_{\text{eff}} = m - \frac{1}{2}(V_0 - S)(r)\). As in Ref. [40], only the large component in the Dirac spinor \(\psi_k\) is used in the calculation of quantity \(\langle k|V_{s.o.}|k \rangle\). One finds that for the spherical state the \(SO_1\) is very small due to shell closure (not exact zero because of the elimination of small component in the Dirac spinor), while the \(SO_2\) is very large, which reflects the size of spin-orbit splitting energies. Moreover, it is shown that the \(SO_2\) decreases with the deformation \(|\beta\rangle\) in both oblate and prolate sides, which is found to be a common feature in atomic nuclei. In particular, the values of \(SO_1\) and \(SO_2\) become close to each other in the configurations with \(|\beta\rangle > 1.0\), which is a sequence of population of nucleons from the upward spherical (spin-orbit antiparallel) \(p_{1/2}\) orbital to the downward spherical \(d_{3/2}\) orbital with the increasing of quadrupole deformation \(\beta\). The \(SO_1\) value turns out to be about 3.8 MeV for the configuration at \(\beta = 3.6\) with 4\(a\) LCS. Figure 2 shows the comparison of matrix element \(\langle k|V_{s.o.}|k \rangle\) for each occupied s.p. state in the mean-field states with \(\beta = 0.0\) and \(\beta = 3.6\). We note that the \(\langle k|V_{s.o.}|k \rangle\) values at \(\beta = 3.6\) are much smaller than those at the spherical shell-like state. In particular, the values of the matrix elements for \(\beta = 3.6\) are always negative, which indicates that the spin and orbital angular momenta of all nucleons are parallel to each other. We note that the spin-orbit interaction energy \(SO_1\) is not exactly zero, which may indicate that there exists a small mixture of non-\(\alpha\)-cluster components in the configuration.

Figure 3 displays the calculated low-spin spectra of \(^{16}\text{O}\) in comparison with experimental data. Following Ref. [4], the level sequences with similar structure are connected with dotted lines. Even though the calculated excitation energies are systematically higher than the data, three rotational bands based on different configurations are shown. The excitation energies of “kite”-like rotational states are in qualitative agreement with the measured levels connected with dotted lines, the band-head state of which \((0^+_2)\) has been a mysterious state and was suggested to have \(^{12}\text{C}(0^+_2) + \alpha\) structure in Refs. [18, 42, 44–46]. Table I presents the detailed properties of these “kite”-like states. Our calculated \(B(E2)\) values are in rather good agreement with the data [43]. The present calculation provides a full microscopic description of these states and suggests these states to have “kite”-like structure.

![FIG. 3: (Color online) (a) Experimental and (b) calculated low-spin spectra for \(^{16}\text{O}\). The data are taken from Ref. [41].](image)

TABLE I: The rms charge radius \(R_{\text{c}}^l\) (fm), spectroscopic quadrupole moment \(Q_s\) (e·fm\(^2\)) and \(E2\) transition strength \(B(E2)\) (e\(^2\)·fm\(^4\)) for the low-spin states with “kite”-like structure from the GCM+PN1DAMP calculation using the PC-PK1 force. The results are compared with other calculations [14, 18, 42] and the experimental data [43].

| \(J^+\) | \(R_{\text{c}}^l\) ($) | \(Q_s\) (e·fm\(^2\)) | Present Ref. [14] | Ref. [18] | Ref. [42] | Exp. [43] |
|------|----------|-----------------|------------------|------------|----------|----------|
| 0\(^+\) | 3.01     | 0               | 97.9             | 48.2       | 177      | 60.1     | 65 ± 7   |
| 2\(^+\) | 3.05     | -23.0           | 192.1            | 290        | 96.2     | 156 ± 14 |
| 4\(^+\) | 3.11     | -34.5           | 226.9            | 301        | 81.6     | 9.6      |
TABLE II: The excitation energy $E_x$ (MeV), rms charge radius $R_c^2$ (fm), spectroscopic quadrupole moment $Q_s$ (e fm$^2$) and $E2$ transition strength $B(E2 \downarrow)$ (e$^2$fm$^4$) for the low-spin states with dominated $4\alpha$ LCS from the GCM+PN1DAMP calculation using the PC-PK1 force. The calculated rms charge radius for the ground state is $2.76$ fm. The ratio $R_c^2/R_{c^*}^2$ is around 1.4.

| $J^+$ | $E_x$  | $R_c^2$ | $Q_s$  | $B(E2 \downarrow)$ |
|------|--------|---------|--------|-------------------|
| 0^+  | 29.6   | 3.88    | 0      | 561               |
| 2^+  | 30.4   | 3.95    | -50.4  | 874               |
| 4^+  | 32.0   | 3.99    | -66.5  | 874               |
| 6^+  | 34.8   | 4.02    | -74.4  | 663               |

with oblate, “kite”-like and LCS characters, where the $g^J(\beta)$s are related to the weight function $f^{JNZ}_\alpha$ in Eq. (2) by the following relation,

$$g^J(\beta) = \sum_{\beta'} (\mathcal{N}^J(\beta, \beta'))^{1/2} f^{JNZ}_\alpha(\beta'),$$  \hspace{1cm} (12)

and are orthonormal to each other. It is shown in Fig. 4 that the collective wave functions of $0^+, 2^+, 4^+$ LCS candidate states are very similar with a sharp peak at $\beta = 3.2$, while the $6^+$ state is fragmented. Table II presents the properties of the low-spin states with $4\alpha$ LCS. The rms charge radii $R_c^2$ of these states are around $3.9$ fm, which is close to the value of $3.8$ fm in the recent GCM calculation based on the AMD model [18], but much smaller than the value of $5.0$ fm by the $4\alpha$ orthogonality condition model [47]. The ratio to the rms charge radii of ground state $R_c^2/R_{c^*}^2$ is around $1.4$, in surprised agreement with the theoretical threshold value of $\sim 1.45$ for $\alpha$ formation [10].

Figure 5 displays the density of the dominated mean-field configurations in the $4\alpha$ LCS candidate states with deformation parameter $\beta = 2.8, 3.2, 3.6$ and $4.0$. The $4\alpha$ LCS is shown clearly in these configurations, especially at $\beta = 3.6$. To understand the nature of the $4\alpha$ LCS, the total intrinsic density of mean-field state at $\beta = 3.6$ is decomposed into densities of the four lowest occupied s.p. states, as shown in Fig. 6. Our main findings are as follows:

i) Integration of the densities in panel (a), (b), (c) and

FIG. 6: (Color online) The density distribution (in fm$^{-3}$) of each s.p. state corresponding to the mean-field configuration with $\beta = 3.6$ in $^{16}$O. The density at $x = y = 0.3$ fm as a function of $z$ is plotted with gray curves. The s.p. states are labeled with the quantum numbers $[n_x n_y n_z]$ of the largest component in the HO basis for the large component of Dirac spinor, cf. Eq. (7).

FIG. 7: (Color online) The weight $|f_{\alpha k}|^2$ [cf.(7)] of the most dominated components $[n_x n_y n_z]$ in the HO basis for the large component of Dirac spinors in the mean-field configuration with $\beta = 3.6$ in $^{16}$O as a function of quantum number $n_z$. 

FIG. 5: (Color online) The total density distribution (in fm$^{-3}$) of intrinsic states with deformation parameter $\beta = 2.8$ (a), 3.2 (b), 3.6 (c), and 4.0 (d). The rms radii of long and short axis are 3.7 fm and 1.1 fm, respectively, for the mean-field state with $\beta = 3.6$. 

FIG. 7: (Color online) The weight $|f_{\alpha k}|^2$ [cf.(7)] of the most dominated components $[n_x n_y n_z]$ in the HO basis for the large component of Dirac spinors in the mean-field configuration with $\beta = 3.6$ in $^{16}$O as a function of quantum number $n_z$. 

FIG. 6: (Color online) The density distribution (in fm$^{-3}$) of each s.p. state corresponding to the mean-field configuration with $\beta = 3.6$ in $^{16}$O. The density at $x = y = 0.3$ fm as a function of $z$ is plotted with gray curves. The s.p. states are labeled with the quantum numbers $[n_x n_y n_z]$ of the largest component in the HO basis for the large component of Dirac spinor, cf. Eq. (7).

FIG. 5: (Color online) The total density distribution (in fm$^{-3}$) of intrinsic states with deformation parameter $\beta = 2.8$ (a), 3.2 (b), 3.6 (c), and 4.0 (d). The rms radii of long and short axis are 3.7 fm and 1.1 fm, respectively, for the mean-field state with $\beta = 3.6$. 

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(d) of Fig. 6 over the coordinate gives 4.00, 3.99, 4.00, 3.51 respectively. Due to pairing correlation, the missing ~ 0.5 particles (~ 0.25 neutrons and ~ 0.25 protons) are scattered to a higher neutron and proton negative-parity state with the largest HO component [010]. The levels with dominated [010] component is high in energy because of level crossings by the large deformation.

ii) The gray curves in Fig. 6 suggests that the 4α clusters stay along a common axis and nucleons occupy the (s)⁴(p)⁴(d)⁴(f)⁴ configurations. To study the localization of nucleons in these s.p. states, following Ref. [9] we calculate the localization parameter $\alpha = \sqrt{\langle r^2 \rangle - \langle r \rangle^2} / \langle r \rangle$, where the average internucleon equilibrium distance is chosen as $\bar{r} = 0.9$. We obtain 1.4 for Fig. 6(a) and Fig. 6(c) and 2.1 for Fig. 6(b) and Fig. 6(d), all of which are much larger than the typical value ~ 1.0 of localized clustering state. It means that the four α clusters in the 4α LCS candidate states of $^{16}$O are delocalized and close to the quantum liquid/gas phase of nucleonic matter, in agreement with the conclusion found in the recent Brink model study [19].

iii) The weight of the most dominated HO components $|nx,np,nz\rangle$ in the four lowest s.p. states [cf. Fig. 6] is displayed in Fig. 7, where $nx, np$ turn out to be zero in all the dominant HO components. We note that the s.p. states labeled with the [004] and [001] in Fig. 6(b) and (d) actually have other competing HO components. The composition of s.p. wave function indicates the one-dimensional character of the 4α LCS in the mean-field state at $\beta = 3.6$.

**B. 4α LCS in high-spin states**

To search for the 4α LCS in the high-spin states, we perform cranked RMF calculations with various rotational frequencies. We choose the $x$-axis as the cranking axis and start the calculations from a triaxially deformed Woods-saxon potential [39]. Table III lists the properties of convergent LCS candidate solutions with the rotational frequencies $\omega$ in between 2.75 and 4.00 MeV, which is higher than the value in between 1.9 (2.0) and 2.2 (2.1) MeV by the cranking SHF calculation using SkI4 (SLy6) force [7]. The angular momentum for the state with LCS ranges from 12.6$b$ to 18.0$b$, which is almost the same as that by the cranking SHF calculation [7] as a consequence of its quantum nature. According to the definition of angular momentum $\sqrt{J_{\text{cra}}(J_{\text{cra}}+1)} = \langle J_z \rangle^2$, where $\langle J_z \rangle = \sum_A A \langle k' j' k | J_z | k \rangle$, one finds that the angular momentum $J$ is defined by the spin direction of all nucleons, different from the classic quantities (such as radius and MOI) which are much dependent on the nuclear density distribution. The difference in the cranking frequency comes from the different MOI of the states. The

| $\langle \hbar \omega \rangle$ | $J_{\text{cra}}$ | $J_{\text{rid}}$ | $E_{\text{tot}}$ | $(\beta, \gamma)$ | $R_{\text{ch}}$ | SO1 | SO2 |
|-----------------|------------|----------|----------------|----------------|----------|-----|-----|
| 2.75            | 12.63      | 12.73    | −78.3 (2.13, 35′) | 3.55           | 6.9      | 6.9 |
| 3.00            | 13.55      | 13.89    | −75.6 (2.16, 35′) | 3.57           | 6.4      | 6.5 |
| 3.25            | 14.46      | 15.05    | −72.8 (2.20, 35′) | 3.61           | 5.9      | 6.0 |
| 3.50            | 15.39      | 16.20    | −69.7 (2.27, 35′) | 3.64           | 5.5      | 5.6 |
| 3.75            | 16.43      | 17.36    | −65.9 (2.39, 35′) | 3.70           | 5.0      | 5.1 |
| 4.00            | 17.96      | 18.52    | −60.1 (2.65, 35′) | 3.82           | 4.3      | 4.6 |

**TABLE III:** The rotational frequency $\hbar \omega$ (MeV), angular momentum $J(\hbar)$, total energy $E_{\text{tot}}$ (MeV), deformations $(\beta, \gamma)$, charge radius (fm), and the spin-orbit energies SO1 and SO2 (MeV). The angular momentum with the rigid-body MOI is calculated as $J_{\text{rid}} = J \omega$, where the MOI $1/(2J) = 0.11$ (MeVh$^{-2}$) is obtained from the parameterization of the energy from the cranking RMF calculations to the rotational formula $E_{\text{tot}}(J_{\text{cra}}) = J_{\text{cra}}(J_{\text{cra}} + 1)/(2J)$.

**FIG. 8:** (Color online) The total density distribution (in fm$^{-3}$) at $x = 0.3$ fm in $^{16}$O corresponding to LCS by the cranking RMF calculation with rotational frequency $\omega = 3.0, 3.5, 4.0$ MeV, respectively. The rms radii of (long, mediate, short) axis are (a) (3.0, 3.1, 3.1) fm, (b) (3.1, 1.4, 1.1) fm, and (c) (3.3, 1.4, 1.1) fm, respectively.
configurations illustrated in Fig. 5 and that by the cranking HF calculation based on the non-relativistic Skyrme EDF [7]. The length of the longest $z$-axis with total nucleon density $\rho \leq 0.02$ fm$^{-3}$ is $\sim 13$ fm, much shorter than that ($\sim 16$ fm) in the non-relativistic Skyrme calculation [7]. It provides a simple explanation for the smaller MOI $\mathcal{J}$ (by a factor of 1.5 – 1.8) in the present calculation.

Figure 9 displays the energy of occupied s.p. states in the LCS candidate states by the cranking RMF calculation as a function of rotational frequency $\omega$. Each state is labeled with the quantum numbers of the largest HO component $[n_x n_y n_z]$ in the large component of Dirac spinor, cf. (7).

i) The excitation energy at low-spin region from the pure AMP calculation for the mean-field state at $\beta = 3.2$ is very close to that of LCS candidate states by the GCM+PN1DAMP calculation. It can be understood from Fig. 4 that the GCM states are dominated by the single configuration at $\beta = 3.2$. The MOI of the low-spin states from the GCM calculation turns out to be $\hbar^2/(2\mathcal{J}) = 0.12$ MeV.

ii) The band-head excitation energy by the extrapolation from the cranking RMF results is 30.2 MeV at $J = 0$, very close to the energy 29.6 MeV of $0^+$ state by the GCM+PN1DAMP calculation. The MOI is estimated to be $\hbar^2/(2\mathcal{J}) = 0.11$ MeV, larger than the value 0.06–0.08 MeV obtained from the cranking SHF calculation. In other words, we obtain a smaller MOI $\mathcal{J}$, which is multiplied with a higher rotational frequencies $\omega$ resulting the same angular momentum as the cranking SHF calculation for LCS candidate states.

IV. SUMMARY

We have presented a detailed study of the $4\alpha$ linear-chain structure in both low-spin and high-spin excited states of $^{16}$O within a covariant density functional theory. The low-spin states have been calculated by configuration mixing of particle-number and angular-momentum projected quadrupole deformed mean-field states from deformation constrained RMF+BCS calculation. The high-spin states have been determined by the cranking RMF calculation. Our conclusions for the linear-chain
structure in high-lying low-spin and high-spin states are summarized as follows:

i) For the low-spin candidate states, an evident $4\alpha$ LCS has been shown in the dominated intrinsic configurations. The analysis of the intrinsic configuration suggests that the $4\alpha$ clusters stay along a common axis and nucleons occupy the $(s)^4(p)^2(d)^4(f)^4$ configurations in a nonlocal way. Moreover, our results indicate that the spin and orbital angular momenta are parallel in the LCS states but the sum of spin-orbit splitting energies turns out to be much smaller than that of shell-like state, which is consistent with the picture of formation of $\alpha$ clusters. The dynamical correlation effects from restoration of rotational symmetry and configuration mixing play an important role in lowering ($\sim 9$ MeV) the excitation energy of these exotic states.

Besides, the rotational band built the second $0^+$ state has been extensively discussed for many years and this band has been considered to have $^{12}$C+$\alpha$ structure. We have reproduced the energies and $B(E2)$ values of this band rather well in a fully microscopic way, and the dominant configuration turns out to be four $\alpha$ clusters with “kite”-like shape.

ii) For the rotational high-spin candidate states with $4\alpha$ LCS, the MOI and band-head ($J = 0$) excitation energy are estimated to be around 0.11 MeV and 30 MeV, respectively, which are slightly different from the results (0.06 – 0.08 MeV and 38 MeV) of previous cranking SHF calculations. Moreover, we have found that the $4\alpha$ LCS in high-spin excited states is less pronounced (less stretched along the symmetric axis) than that by the cranking calculation based on the non-relativistic Skyrme EDFs. However, the LCS states are found at the same angular momenta region, i.e. $13 – 18\hbar$.

Finally, we point out that the present study demonstrates the ability of the projected GCM based on the CDFT for the cluster structures in nuclear low-spin states. The extension of the present study to the cluster structures in excited states of other light nuclei is also very interesting, such as the stability of clusters found in superdeformed states of Ar isotopes against the shape fluctuation [49]. For some other exotic cluster states, however, the other shape degrees of freedom, such as octupole deformation might be required. Work along this direction is in progress. Moreover, we note that for the high-spin states the cranking solutions pertain to the intrinsic frame and therefore cannot directly be compared to experiment. The implementation of AMP technique for time-reversal violated system is necessary to carry out a detailed spectroscopic study for the high-spin states. Some efforts have recently been devoted along this direction based on the non-relativistic Skyrme EDF [50, 51]. The application of such methods for the clusters in nuclear high-spin states will also be very interesting.

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