Evidence of chiral bands in even-even nuclei

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Evidence for chiral doublet bands has been observed for the first time in the even-even nucleus $^{136}$Nd. One chiral band was firmly established. Four other candidates for chiral bands were also identified, which can contribute to the realization of the multiple pairs of chiral doublet bands ($M_D^\chi$) phenomenon. The observed bands are investigated by the constrained and tilted axis cranking covariant density functional theory (TAC-CDFT). Possible configurations have been explored. The experimental energy spectra, angular momenta, and $B(M1)/B(E2)$ values for the assigned configurations are globally reproduced by TAC-CDFT. Calculated results support the chiral interpretation of the observed bands, which correspond to shapes with maximum triaxiality induced by different multiquasiparticle configurations in $^{136}$Nd.

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FIG. 1. Partial level scheme of $^{136}$Nd showing the newly identified doublet bands and their decay toward low-lying levels. The intensities of the bands relative to the intensity of the 374-keV, $2^+ \rightarrow 0^+$ transition are also indicated below the band labels. The bands are labeled as in the complete level scheme that will be published in a forthcoming paper [34].

The $^{135}$Nd and $^{136}$Nd nuclei were the most strongly populated in the reaction, with cross sections of around 100 mb each, for a total of 480 mb. A number of $5.1 \times 10^{10}$ three-fold and higher prompt γ-ray coincidence events were accumulated using the JUROGAM II array which has an efficiency of around 4.3% at 1.3 MeV. The events were time stamped by the total data readout (TDR) data acquisition, and sorted using the GRAIN code [35]. Fully symmetrized, three-dimensional ($E_\gamma$-$E_\gamma$-$E_\gamma$) and four-dimensional ($E_\gamma$-$E_\gamma$-$E_\gamma$-$E_\gamma$) matrices of γ-ray energies $E_\gamma$ were analyzed using the RADWARE [36,37] analysis package. Spin and parity assignments were made on the basis of the measured DCO (directional correlations of oriented states) ratios ($R_{DCO}$), two-point angular correlation (anisotropy) ratio $R_{an}$ [38,39], angular distributions [40], and polarization asymmetries [41] of the transitions depopulating the states. Details of the angular correlation analysis will be given in a forthcoming paper [34].

The partial level scheme of $^{136}$Nd showing five pairs of doublet bands and their decay toward low-lying states is given in Fig. 1. An example of double-gated spectra showing the transitions connecting one of the newly identified bands, band D2-C, to band D2, is given in Fig. 2. Similar spectra have been also obtained for the other doublet bands. One of the specific features of these bands is their decay to the yrast partners via high-energy quadrupole transitions, which, due to the $E_\gamma^5$ dependence, are stronger than the in-band dipole transitions which have low energy and $E_\gamma^3$ dependence, resulting in a factor of 10 to 20 in favor of the connecting transitions. Due to this fact, many in-band dipole transitions of the new bands could not be observed with the present statistics. The $B(M1)/B(E2)$ branching ratios for some of the observed bands are shown in Fig. 3. One can see an increase at spin around 20 for the bands D1, D3, and D4, which are well correlated with an increased single-particle alignment marking a change in the configuration. The $B(M1)/B(E2)$ values of all bands have a decreasing behavior at the highest spins. For the strongest pair of chiral bands, D5 and D5-chiral, we could extract the $B(M1)/B(E2)$ values for both bands, which are nearly identical within errors in the observed spin range.

Generally speaking, for the description of chiral rotations, three-dimensional tilted axis cranking covariant density functional theory (3D TAC-CDFT) should be adopted, which has been used to investigate the chiral modes of $^{106}$Rh [42]. For the doublet bands built on the 4- and 6-quasiparticle configurations identified in $^{136}$Nd, the 3D TAC-CDFT calculations are very challenging. For simplicity, the observed five pairs of rotational bands are investigated within the framework of TAC-CDFT [43–45]. However, the chiral nature
of the assigned configurations was tested by performing 3D TAC-CDFT calculations for one chiral pair as an example. The well-known relativistic density functional PC-PK1 [46] is adopted and the Dirac equation is solved in a 3D harmonic oscillator basis in Cartesian coordinates with 10 major shells which provide convergent results in TAC-CDFT calculations [45].

By minimizing the energy with respect to the triaxial deformation \( \gamma \), both adiabatic and configuration-fixed constrained calculations similar to those in Ref. [18] were performed for various low-lying particle-hole excitations in \(^{136}\text{Nd}\). Detailed results are shown in Table I. There are three positive-parity configurations (labeled A, B, and C) and five negative-parity configurations (labeled D, D*, E, F, and H) which are candidates for the observed nearly degenerate partner bands. All these configurations with particle-hole excitations possess remarkable triaxial deformation, which is the typical feature for the chiral rotational bands. By considering the pairing correlations with separable pairing force [47,48], the changes of total energy and total angular momentum at the rotational frequency 0.2 MeV are within 0.005% and 4.5% for configuration A. Therefore, pairing correlations can be safely neglected here. In the following, the configurations listed in Table I will be justified based on the quasiparticle alignments. The energy spectra, angular momenta, as well as the \( B(M1)/B(E2) \) values of the observed doublet bands are also investigated within the TAC-CDFT framework for the assigned configurations.

Here, we take the bands D1 and D1-C as an example to show how the configurations are assigned. The quasiparticle alignments calculated by TAC-CDFT are shown in Fig. 4 in comparison with the experimental data. The experimental quasiparticle alignments are derived as in Ref. [49]. The parameters \( J_0 = 11 \hbar^2/\text{MeV} \) and \( J_1 = 20 \hbar^4/\text{MeV}^2 \) have been adopted for the Harris formula \( J = J_0 + J_1 \omega^2 \). It is clearly seen that the calculated results based on configuration A reproduce the experimental data of band D1 very well. Therefore, the configuration of band D1 is assigned as A, as shown in Table I.

The assigned configuration A is further investigated by examining the calculated excitation energies of bands D1 and D1-C using the TAC-CDFT formalism, as shown in the left panels of Fig. 5. Note that the energy references for the
TABLE I. Unpaired nucleon configurations labeled A–H and the corresponding parities, calculated by constrained CDFT. The deformation parameters $\beta$ and $\gamma$, and the excitation energies $E_4$ are also provided. It is noted that the letter G represents the ground state of $^{136}$Nd and there are no nucleons resulting from the breaking of a Cooper pair for this configuration. $\Omega$ is the projection of the spin of the $d_{5/2}$ orbital on the quantization axis, while $\pi$ and $\nu$ indicate protons and neutrons, respectively. The configurations are given in terms of spherical single-particle orbitals.

<table>
<thead>
<tr>
<th>State</th>
<th>$E_4$ (MeV)</th>
<th>Parity</th>
<th>$\beta$</th>
<th>$\gamma$</th>
<th>Unpaired nucleons</th>
</tr>
</thead>
<tbody>
<tr>
<td>G</td>
<td>0.000</td>
<td>+</td>
<td>0.24</td>
<td>27$^+$</td>
<td>$\pi(h_{11/2})^1(d_{5/2}g_{7/2})^{-1}\otimes\nu(h_{11/2})^{-1}(s_{1/2}d_{3/2})^{-1}$</td>
</tr>
<tr>
<td>A</td>
<td>0.335</td>
<td>+</td>
<td>0.21</td>
<td>21$^+$</td>
<td>$\pi(h_{11/2})^1(d_{5/2}g_{7/2})^{-1}\otimes\nu(h_{11/2})^{-1}(s_{1/2}d_{3/2})^{-1}$</td>
</tr>
<tr>
<td>B</td>
<td>3.419</td>
<td>+</td>
<td>0.22</td>
<td>19$^+$</td>
<td>$\pi(h_{11/2})^1(d_{5/2}g_{7/2})^{-1}\otimes\nu(h_{11/2})^{-1}(s_{1/2}d_{3/2})^{-1}$</td>
</tr>
<tr>
<td>C</td>
<td>3.704</td>
<td>+</td>
<td>0.26</td>
<td>23$^+$</td>
<td>$\pi(h_{11/2})^2(d_{5/2}g_{7/2})^{-2}\otimes\nu(h_{11/2})^{-1}(f_{7/2}h_{9/2})^{-1}$</td>
</tr>
<tr>
<td>D</td>
<td>1.173</td>
<td>–</td>
<td>0.22</td>
<td>19$^+$</td>
<td>$\pi(h_{11/2})^2\otimes\nu(h_{11/2})^{-1}(s_{1/2}d_{3/2})^{-1}(\Omega \sim 0\text{ or } \Omega)$</td>
</tr>
<tr>
<td>D*</td>
<td>1.346</td>
<td>–</td>
<td>0.21</td>
<td>22$^+$</td>
<td>$\pi(h_{11/2})^2\otimes\nu(h_{11/2})^{-1}(s_{1/2}d_{3/2})^{-1}(\Omega \sim -\frac{5}{2})$</td>
</tr>
<tr>
<td>E</td>
<td>1.937</td>
<td>–</td>
<td>0.21</td>
<td>23$^+$</td>
<td>$\pi(h_{11/2})^2\otimes\nu(h_{11/2})^{-1}(s_{1/2}d_{3/2})^{-1}(\Omega \sim -\frac{3}{2})$</td>
</tr>
<tr>
<td>F</td>
<td>2.778</td>
<td>–</td>
<td>0.20</td>
<td>35$^+$</td>
<td>$\pi(h_{11/2})^1(d_{5/2}g_{7/2})^{-1}\otimes\nu(h_{11/2})^{-2}$</td>
</tr>
<tr>
<td>H</td>
<td>3.494</td>
<td>–</td>
<td>0.20</td>
<td>37$^+$</td>
<td>$\pi(h_{11/2})^1(d_{5/2}g_{7/2})^{-1}\otimes\nu(h_{11/2})^{-2}$</td>
</tr>
</tbody>
</table>

positive- and negative-parity calculated bands are taken as those of band D1 at $I = 11\hbar$ and of band D4 at $I = 15\hbar$, respectively. It is found that the theoretical results based on configuration A agree with experimental data of band D1 satisfactorily, and the configuration assignment to band D1 is validated. Meanwhile, no other configuration can be found to reproduce the results of D1-C, which indicates that bands D1 and D1-C are based on the same intrinsic state. Therefore one can safely identify D1 and D1-C as good candidates for chiral partner bands.

Similarly, one can conclude that D2 and D2-C, D5 and D5-C, D3 and D3-C, as well as D4 and D4-C are candidates of chiral doublets bands built on the same intrinsic configurations B, C, D*, and D, respectively. It is noted that the back-bending phenomenon observed in bands D2, D3, and D4 can be due to crossing with 8-, 6-, and 6-qp configurations, respectively, which is beyond the current TAC-CDFT calculation.

The $B(M1)/B(E2)$ values of the five pairs of chiral bands have also been studied self-consistently, as shown in Fig. 3. The calculated $B(M1)/B(E2)$ values based on configurations A and B underestimate the experimental data of bands D1 and D2, while the theoretical $B(M1)/B(E2)$ values of the assigned configuration are in good agreement with the experimental data of band D5 over all the observed spin range. As for bands D3 and D4, a satisfactory agreement with experimental is only present in the low- and high-spin parts. An abrupt increase of $B(M1)/B(E2)$ values in bands D3 and D4 is observed near back-bending and cannot be reproduced by the calculations. As mentioned above, a back-bending also exists in band D2, which induces an abrupt increase of the $B(M1)/B(E2)$ values. Note that although the absolute $B(M1)/B(E2)$ values for bands D1 and D2 are not reproduced, the relative differences agree with the experimental data well. A better agreement over the entire observed spin range will hopefully be obtained in the near future with extensive theoretical calculations taking into account the possible change of configuration along the bands and the configuration mixing.

Finally, to give more theoretical support for the observed chiral rotation in $^{136}$Nd, we also performed challenging calculations based on 3D TAC-CDFT. The D* configuration corresponding to band D3 was calculated as an example. The orientation angle $\phi$ for the total angular momentum of band D3 calculated with 3D TAC-CDFT is shown in Fig. 6.
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FIG. 6. Evolution of the azimuth angle $\phi$ as a function of rotational frequency, for the total angular momentum of the configuration $D^*$ assigned to band D3, calculated by 3D TAC-CDFT.

It is found that the azimuth angle $\phi$ for band D3 vanishes at low rotational frequencies, providing a planar solution. Above the critical rotational frequency $0.5 \text{ MeV}$, a nonzero angle $\phi$ appears, corresponding to an aplanar solution, namely, chiral rotation, for the selected configuration, namely, $D^*$. The resulting excitation energies and total angular momenta are well reproduced by the present 3D TAC-CDFT calculations. We find aplanar solutions in band D3 which provides very strong support for the existence of chiral rotation. Moreover, the effects of 3D rotation on $B(M1)/B(E2)$ values, the total angular momentum, and excitation energies are marginal.

Summarizing, five pairs of nearly degenerate rotational bands were identified in $^{136}\text{Nd}$, one of which has a clear chiral character. It is the first time that chiral bands are observed in an even-even nucleus at high spins. The other four doublet bands, which are weakly populated, have no measured $B(M1)/B(E2)$ values, but the calculated assigned configurations show chiral geometry. This set of five nearly degenerate bands is the largest observed in a single nucleus until now. The observed bands were discussed using the TAC-CDFT, 3D TAC-CDFT, and MQ-PRM models, which reveal the chiral nature in the assigned configurations. Further theoretical calculations to investigate in detail to what extent the chiral geometry of the observed bands is realized are in progress.

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