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Collective Hamiltonian for chiral modes

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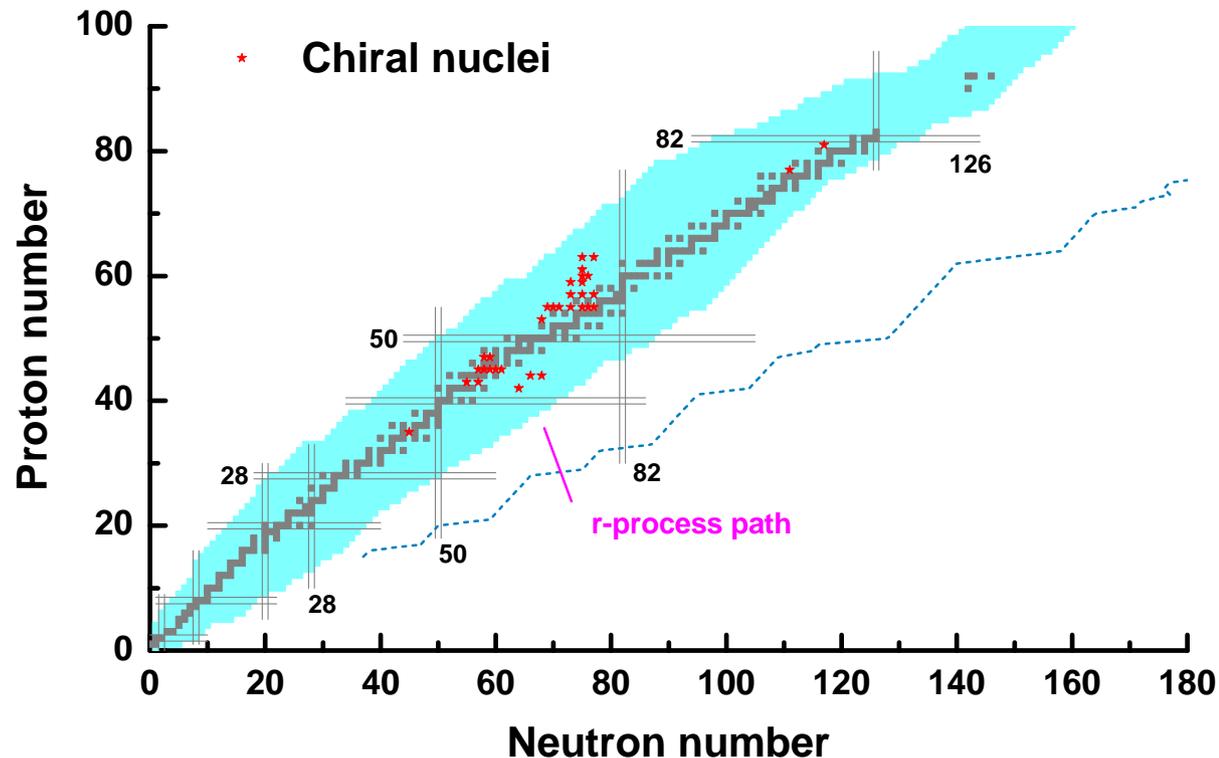
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Introduction

- Chirality is a subject of general interest in molecular physics, elementary particle physics, and optical physics.
- In nuclear physics, the occurrence of chirality was originally suggested in 1997. [Frauendorf and Meng1997NPA](#)
- The chirality was firstly observed in the four $N = 75$ isotones: ^{130}Cs , ^{132}La , ^{134}Pr , and ^{136}Pm in 2001. [Starosta2001PRL](#)
- So far, more than 30 candidate chiral nuclei have been reported experimentally in the $A \sim 80, 100, 130,$ and 190 mass regions.



Introduction

- Chiral doublet bands were firstly predicted by tilted axis cranking (TAC) approach and particle rotor model (PRM). [Frauendorf and Meng1997NPA](#)
- Numerous efforts have been devoted to the development of both the TAC methods and PRM models.
- PRM: [Peng2003PRC](#), [Koike2004PRL](#), [Zhang2007PRC](#), [Droste2009EPJA](#), [Qi2009PLB](#), [Lawrie2010PLB](#)
 - ✓ quantal model: the total angular momentum is a good quantum number; energies and transition probabilities are treated fully quantally.
 - ✗ deformation parameters have to be assumed at the very beginning.
- TAC: [Dimitrov2000PRL](#), [Olbratowski2004PRL](#), [Olbratowski2006PRC](#), [Meng2013Front.Phys.](#)
 - ✓ permits the calculation for the orientation of the density distribution relative to the angular momentum vector; easily extended to the multi-quasiparticle case.
 - ✗ cannot give the quantum tunneling of chiral doublet bands.
- TAC+RPA: [Mukhopadhyay2007PRL](#), [Almehed2011PRC](#)
 - ✓ go beyond the mean-field approximation.
 - ✗ restricted in the description of the chiral vibration.

Motivation

- **It is imperative to search a unified method for studying both chiral rotation and vibration in the framework of TAC.**
- **In the present work, the collective Hamiltonian for chiral modes have been constructed.**

Theoretical framework

- **Determine the potential term:** starting from the tilted axis cranking model

$$\begin{aligned}\hat{h}' &= \hat{h}_{\text{def}} - \vec{\omega} \cdot \hat{j}, \\ \vec{\omega} &= (\omega \sin \theta \cos \varphi, \omega \sin \theta \sin \varphi, \omega \cos \theta).\end{aligned}\quad (1)$$

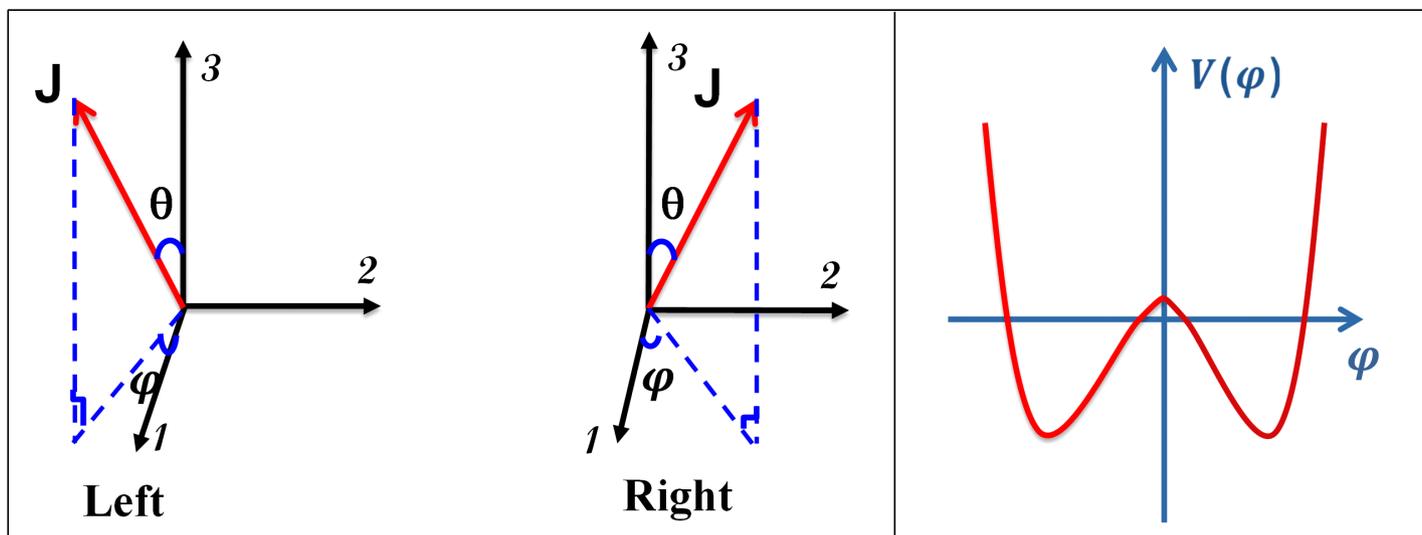
where $\hat{j} = \hat{j}_\pi + \hat{j}_\nu$, and $\hat{h}_{\text{def}} = \hat{h}_{\text{def}}^\pi + \hat{h}_{\text{def}}^\nu$ is the single- j shell Hamiltonian with

$$\hat{h}_{\text{def}}^{\pi(\nu)} = \pm \frac{1}{2} C \left\{ (\hat{j}_3^2 - \frac{j(j+1)}{3}) \cos \gamma + \frac{1}{2\sqrt{3}} (\hat{j}_+^2 + \hat{j}_-^2) \sin \gamma \right\}.\quad (2)$$

Then minimize the total Routhian

$$E'(\theta, \varphi) = \langle h' \rangle - \frac{1}{2} \sum_{k=1}^3 \mathcal{J}_k \omega_k^2, \quad \mathcal{J}_k = \mathcal{J}_0 \sin^2(\gamma - \frac{2\pi}{3}k).\quad (3)$$

with respect to θ for given φ to obtain the collective potential $V(\varphi)$.



Theoretical framework

- **Determine the kinetic term:** the kinetic term is $T_{\text{kin}} = \frac{1}{2}B\dot{\varphi}^2$ with mass parameter B , which is obtained from the cranking approximation as

$$B = 2\hbar^2 \sum_{l \neq 0} \frac{(E_l - E_0) \left| \frac{\partial \vec{\omega}}{\partial \varphi} \langle l | \hat{j} | 0 \rangle \right|^2}{[(E_l - E_0)^2 - \hbar^2 \Omega^2]^2}, \quad (4)$$

in which $|l\rangle$, $|0\rangle$ and E_l , E_0 are respective the eigen states of the cranking Hamiltonian and the corresponding eigenvalues. The Ω is chiral vibrational frequency and can be taken as $\Omega = 0$ for the case of chiral rotation.

- **Collective Hamiltonian:**

✓ Classical form:

$$H_{\text{coll}} = T_{\text{kin}}(\varphi) + V(\varphi) = \frac{1}{2}B\dot{\varphi}^2 + V(\varphi), \quad (5)$$

where $V(\varphi)$ and B have been obtained from TAC calculation.

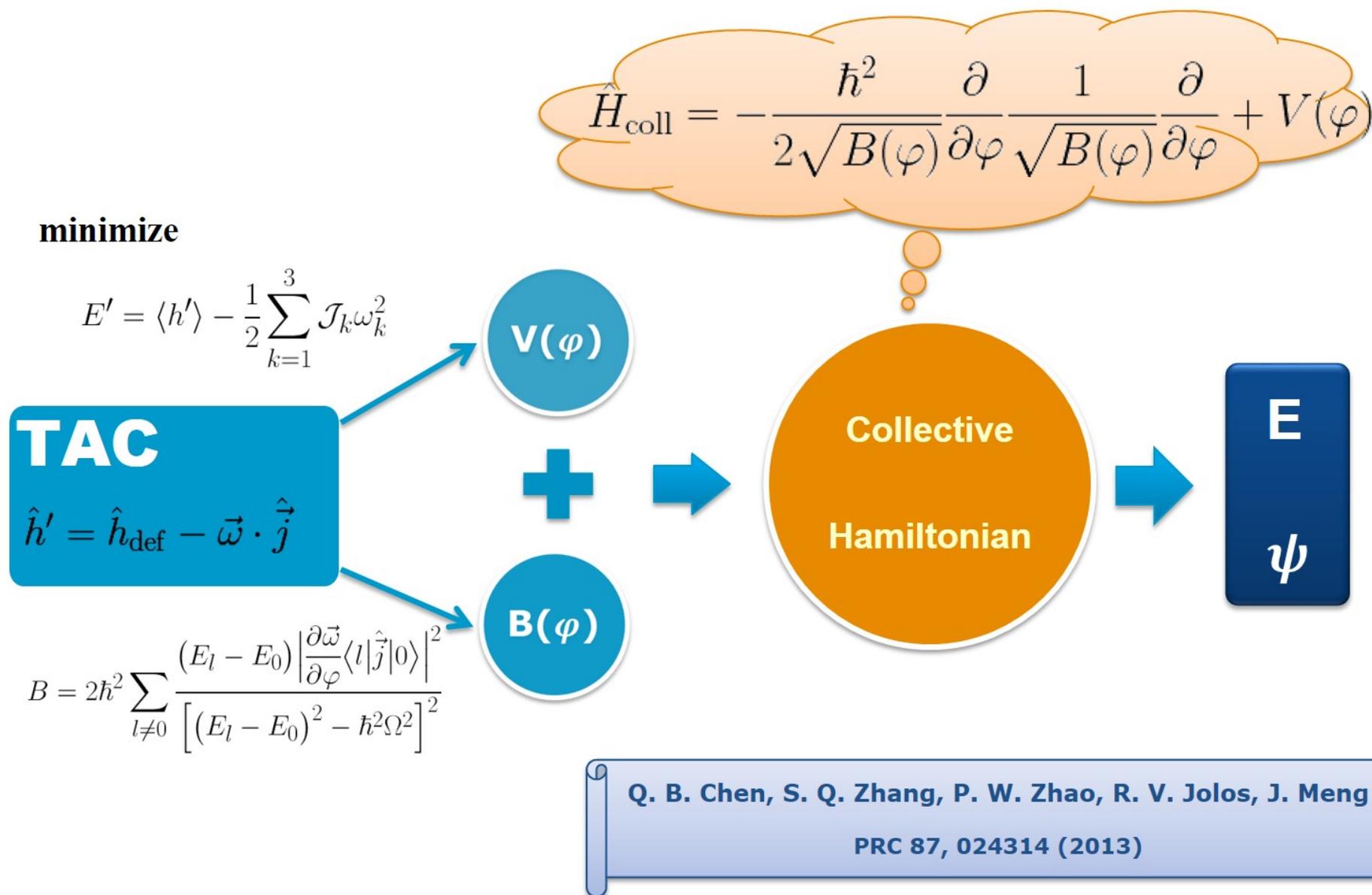
✓ Quantal form:

$$\hat{H}_{\text{coll}} = -\frac{\hbar^2}{2\sqrt{B(\varphi)}} \frac{\partial}{\partial \varphi} \frac{1}{\sqrt{B(\varphi)}} \frac{\partial}{\partial \varphi} + V(\varphi), \quad (6)$$

obtained according to the Pauli prescription:

$$\hat{H}_{\text{kin}} = -\frac{\hbar^2}{2} \frac{1}{\sqrt{\det B}} \sum_{ij} \frac{\partial}{\partial q_i} \sqrt{\det B} (B^{-1})_{ij} \frac{\partial}{\partial q_j}. \quad (7)$$

Theoretical framework



Numerical details

- Angular moments of valence nucleons: $j_\pi = j_\nu = 11/2 \hbar$;
- Single- j shell Hamiltonian coefficients: $C_\pi = 0.25 \text{ MeV}$, $C_\nu = -0.25 \text{ MeV}$;
- Triaxial deformation: $\gamma = -30^\circ$;
- Moment of inertia: $\mathcal{J}_0 = 40 \hbar^2/\text{MeV}$;
- Potential energy surface mesh points (θ_i, φ_j) is represented as

$$\begin{aligned}\theta_i &= (i - 1) \times 1^\circ, (i = 1, \dots, 91), \\ \varphi_j &= (j - 91) \times 1^\circ, (j = 1, \dots, 181).\end{aligned}$$

Total Routhian

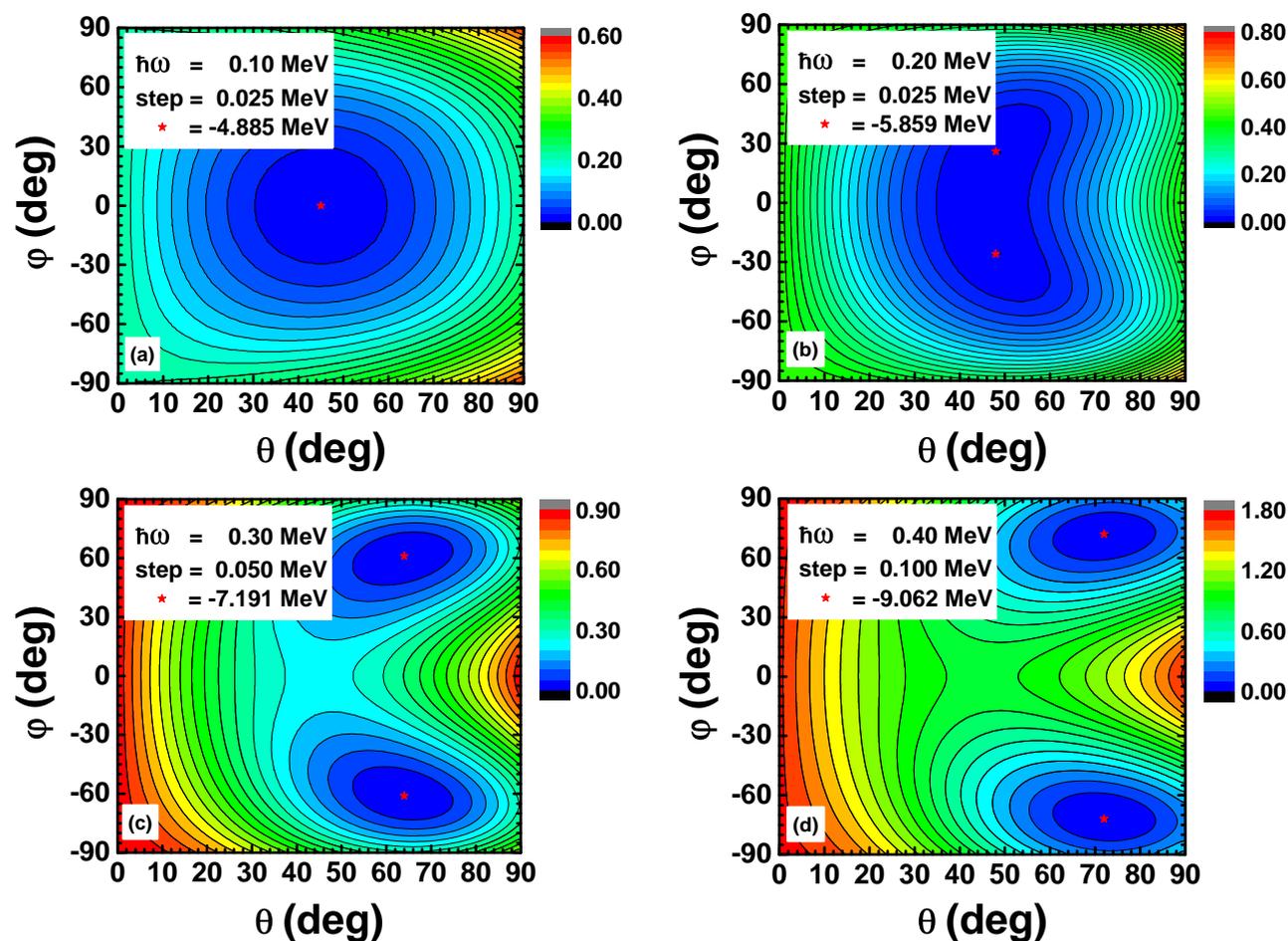


Figure: Total Routhian surface calculations for the $h_{11/2}$ proton particle and the $h_{11/2}$ neutron hole coupled to a triaxial rotor with $\gamma = -30^\circ$ at the frequencies $\hbar\omega = 0.1, 0.2, 0.3, 0.4$ MeV.

Remarks

- Potential energy surfaces are symmetrical with the $\varphi = 0^\circ$.
- The minimal points change from $\varphi = 0^\circ$ to $\varphi \neq 0^\circ$.

Collective potential

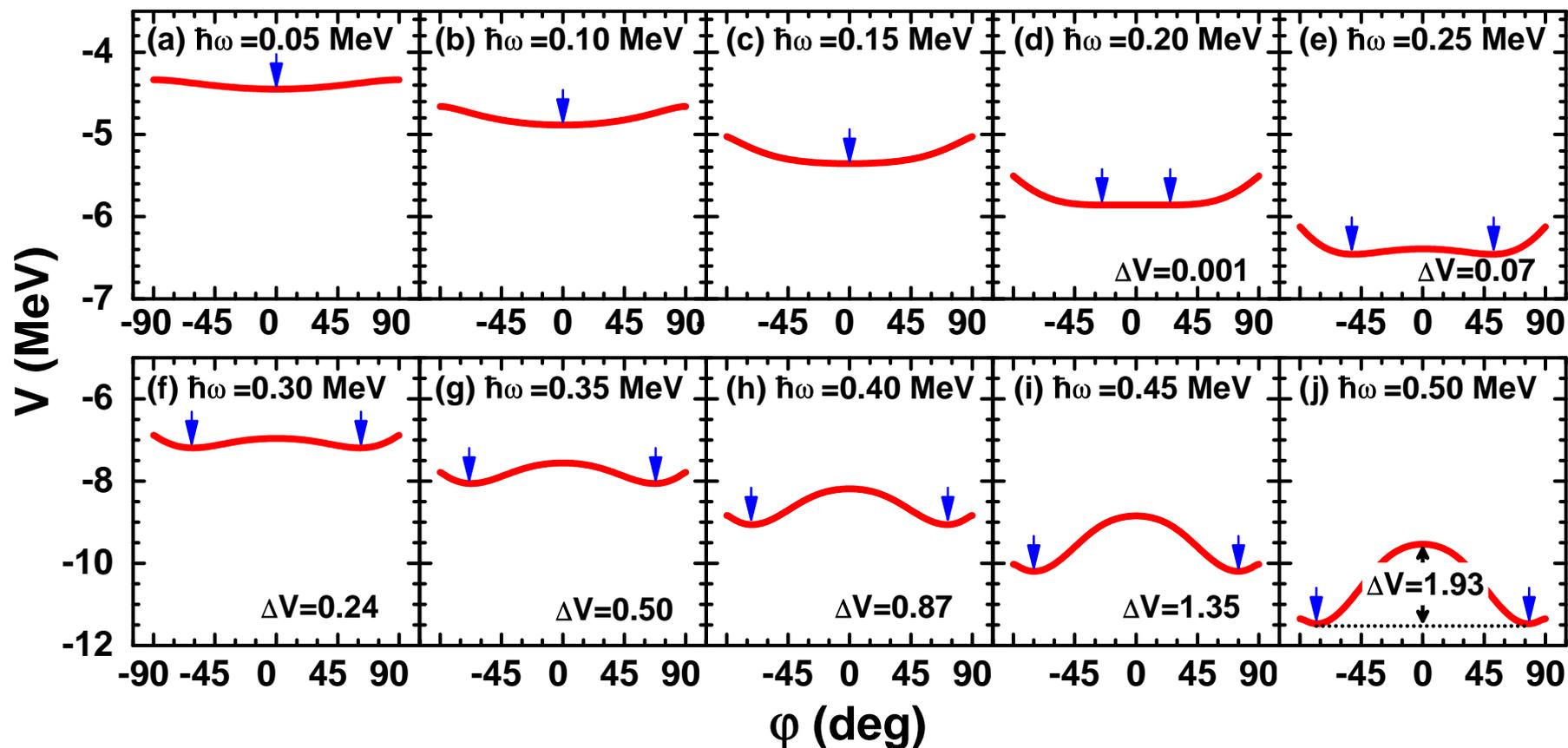


Figure: The potential energy $V(\varphi)$ as a function of φ extracted from the total Routhian surface calculations. The arrow labels the position of the potential minimum V_{\min} .

Remarks

- The shape of potential change from harmonic oscillator type to double well type.
- The potential barrier increases from 1 keV at $\hbar\omega = 0.2$ MeV to about 2 MeV at $\hbar\omega = 0.50$ MeV.

Mass parameter

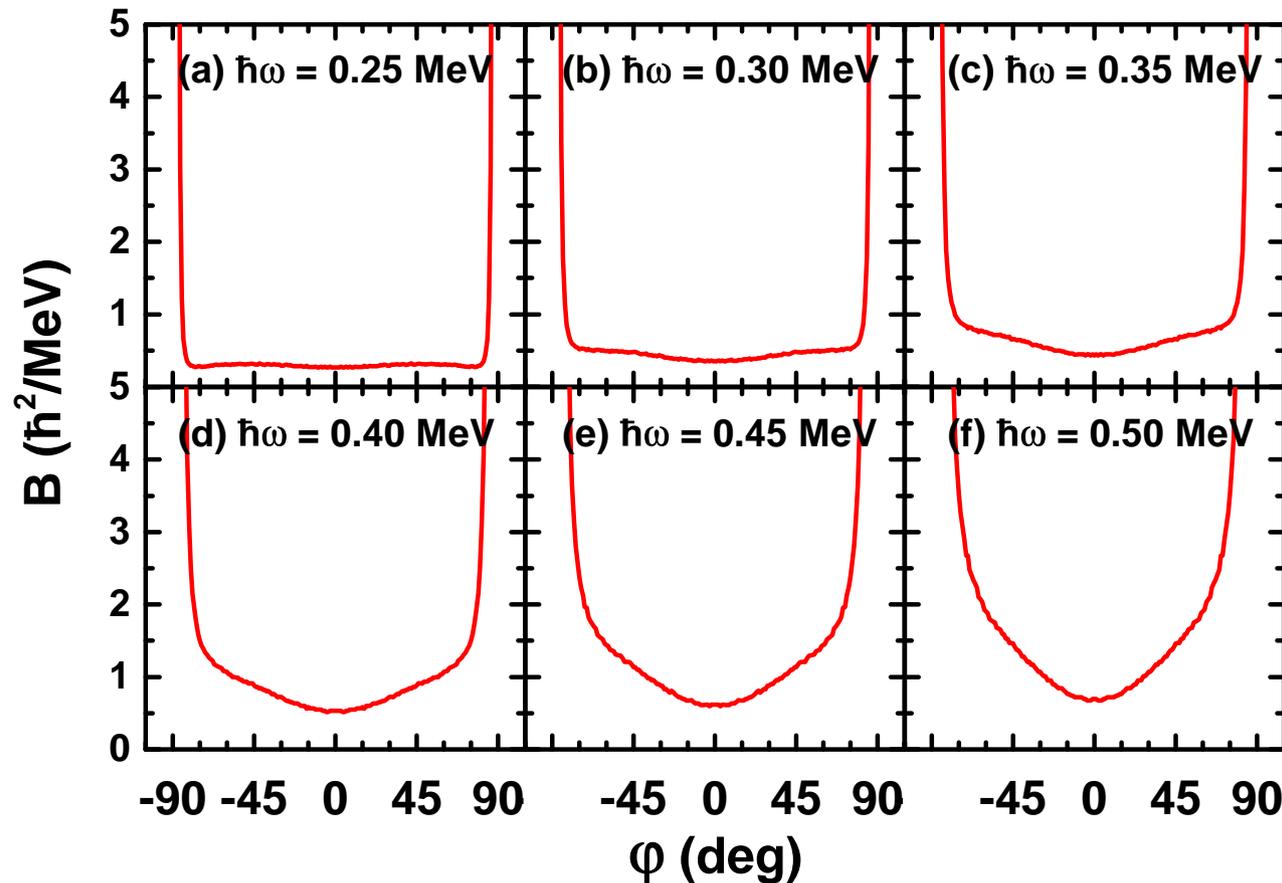


Figure: The mass parameter $B(\varphi)$ as a function of φ for the chiral rotation cases obtained based on TAC.

Remarks

- $B(\varphi)$ is symmetric with respect to $\varphi = 0^\circ$ and increase dramatically when φ approach to $\pm 90^\circ$.
- In the interior part, $B(\varphi)$ is increased remarkably with $|\varphi|$ for $\hbar\omega \geq 0.35 \text{ MeV}$, while its dependence on φ is weak for $\hbar\omega = 0.25$ and 0.35 MeV .

Energy levels

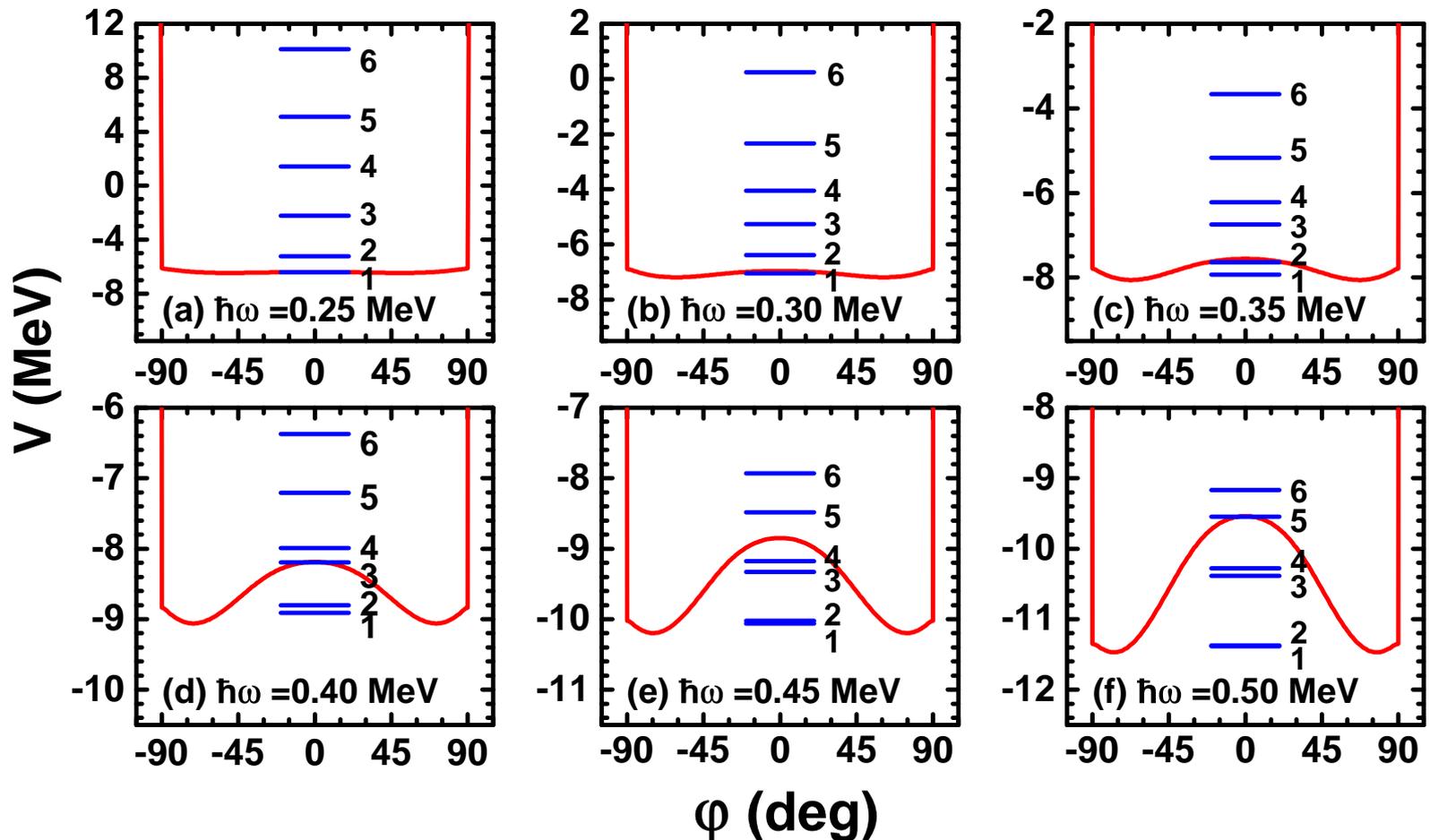


Figure: The six lowest energy levels, labeled as 1-6, obtained from the collective Hamiltonian.

Remarks

- The three pairs of energy levels become close to each other.
- Are the levels 2-4 corresponding to the **excited chiral doublet bands**? [Droste2009EPJA](#), [Chen2010PRC](#), [Hamamoto2013PRC](#)

Wave function

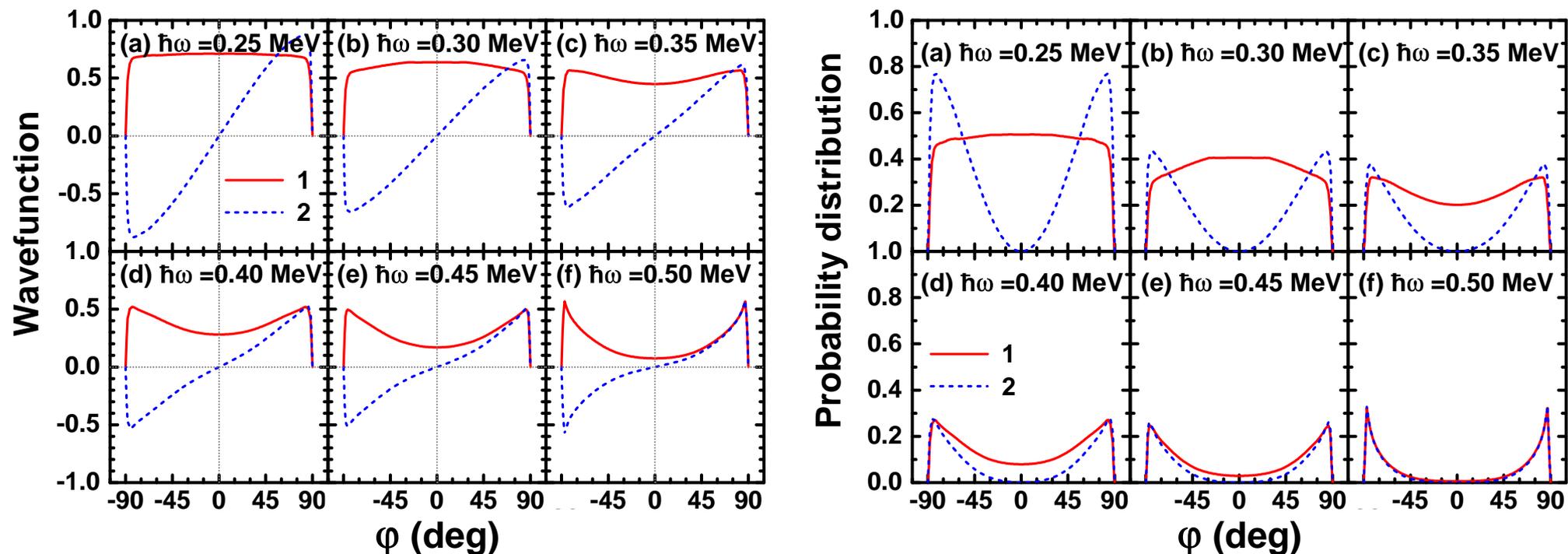


Figure: Wave functions $\psi(\varphi)$ and probability distributions $|\psi(\varphi)|^2$ for the lowest two levels 1 and 2 obtained from collective Hamiltonian.

Remarks

- The wave functions are symmetric for level 1 and antisymmetric for level 2 with respect to $\varphi \rightarrow -\varphi$.
- The wave functions of level 1 tends to show similar pattern with level 2.

Comparison with exact solutions

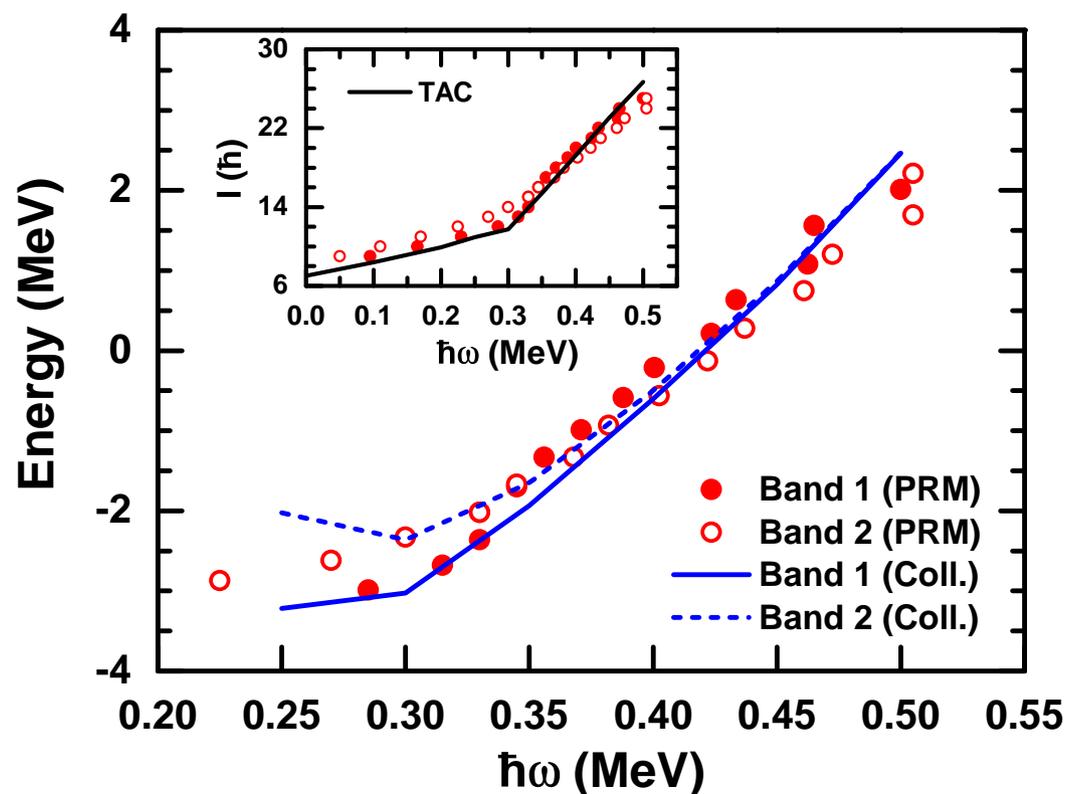


Figure: The energy spectra of the doublet bands obtained from the collective Hamiltonian in comparison with the exact solutions by the PRM.

Remarks

- Apart from the agreement of collective Hamiltonian and PRM results for the yrast band, the partner band of PRM can also be reasonably reproduced by the collective Hamiltonian.
- The second chiral vibration character obtained by PRM is not taken into account by the present collective Hamiltonian investigation.

Summary and perspective

Summary:

- A collective model which is able to describe the chiral rotation and vibration is proposed and applied to a system of one $h_{11/2}$ proton particle and one $h_{11/2}$ neutron hole coupled to a triaxial rigid rotor.
- Based on the tilted axis cranking approach, both the potential energy and mass parameter as functions of φ are obtained and included in the collective Hamiltonian.
- It is found that for chiral rotation, the partner states become more degenerate with the increase of the cranking frequency, and the wave function of levels 1 and 2 tend to show similar pattern.

Perspective:

- The collective Hamiltonian is expected to describe the chiral doublet bands in more realistic nuclei, such as tilted axis cranking relativistic mean-field theory, to obtain a more microscopic collective potential.
- The fluctuation of the potential energy with θ will be taken into account.
- The adiabatic self-consistent collective coordinate (**ASCC**) method will be introduced to calculate the mass parameter. [Marumori1980PTP](#), [Matsuo2000PTP](#), [Hinohara2010PRC](#)

Thank you!