K-Selection rule violating E2 and M1 transition from K=6⁺ isomers to the ground band of ^{172, 174, 176}Hf

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Introduction

The well-deformed hafnium (Hf, Z = 72) isotopes contain multiquasiparticle K- isomers, with half-lives ranging from nanoseconds to years [1]. A number of strongly deformed bands have also been observed in Hf nuclei [2, 3]. Besides deformation and the associated rotational bands long lived K isomers are important features of the rare-earth and other heavy nuclei [4,5]. Here we study the rotational bands and K isomers of some neutron-rich Hf nuclei and their properties using deformed Hartree-Fock and J projection theories. Bands built on the K isomers have also been calculated.

Theoretical Framework

To this end, we have adopted in this work the Deformed Hartree-Fock (DHF) model to get the deformed singleparticle states and the deformed multinucleon confifigurations. Ground and Kisomeric intrinsic states are constructed. and for each intrinsicstate (configuration), states of good angular momenta are obtained by Angular Momentum Projection A deformed shape such as one described by slater determinant of deformed orbits $|\Phi_K\rangle$ is localized in angle and, by the uncertainty principle, is not a state of good angular momentum (J). Thus $|\Phi_K\rangle$ does not have a unique J quantum number and is a superposition of various J states [6–8],

$$|\Phi_{\rm K}\rangle = \sum_{\rm I} C_{\rm IK} |\Psi_{\rm IK}\rangle$$
.

One needs to project out states of good angular momenta from the intrinsic state Φ_K with the Angular Momentum Projection operator,

$$P_K^{IM} = \frac{2I+1}{8\pi^2} \int d\Omega D_{MK}^{I*} \Omega R(\Omega)$$
 2

Results and Discussion

The deformed HF orbits are calculated with a spherical core of 132 Sn; the model space spans the $3s_{1/2}$, $2d_{3/2}$, $2d_{5/2}$, $1g_{7/2}$, $1h_{11/2}$ and $1h_{9/2}$ orbits for protons have energies 3.654, 3.288, 0.731, 0.0, 1.705 and 6.46 MeV, and the $3p_{1/2}$, $3p_{3/2}$, $2f_{5/2}$, $2f_{7/2}$, $1h_{9/2}$ and $1i_{13/2}$ neutron states have energies 4.462, 2.974, 3.432, 0.0, 0.686 and 1.487 MeV respectively. We use a surface delta interaction (with Interaction [9] strength 0.3 MeV for $p_{-p_{1}}$ and $p_{-p_{2}}$ and $p_{-p_{3}}$ interactions) as the residual interaction among the active nucleons in these orbits. It is to be noted that the theoretical spectra shown in FIG.

1 and 2 have good K quantum numbers $(K=0^+ \text{ and } K=6^+)$. The $K=6^+ \text{ isomers are obtained by particle hole excitations over the ground state having <math>5/2^+$ and $7/2^+$ neutron intrinsic structures. The spectra are shown in Figures 1 and 2 for $^{172,176}\text{Hf}$ nuclei.

We have calculated E2 transition matrix elements for the $K=0^+$, $K=6^+$ (in-band) states as also between the K=6⁺ isomers and the ground bands. We find, in our angular momentum projection calculation, quite enhanced collective matrix element values for the in-band transitions; while the E2 transition matrix elements from K=6⁺ isomer band-heads to J=4+ of the ground band are retarded by a factor of about 10-5. Similar retarded matrix elements are also obtained for M1 transitions from the K=6⁺ isomers to the ground bands. Thus, a theoretical understanding of the retarded decay of a K-isomer to lower K bands is possible by angular momentum projection from the deformed intrinsic states.

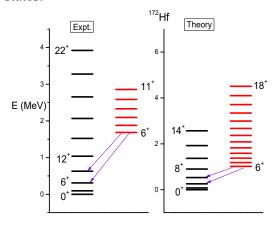


FIG. 1: Energy spectra of ¹⁷²Hf. Experimental values are taken from Ref. [10]

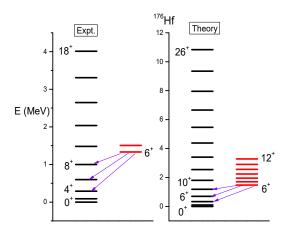


FIG. 2: Energy spectra of ¹⁷⁶Hf. Experimental values are taken from Ref. [10]

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