

Band head spin for triaxial super-deformed bands in $^{165,167}\text{Lu}$

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Abstract. We use VMI model for the prediction of band head spin of Triaxial Super-Deformed (TSD) rotational bands. The calculated and observed transition energies are agreed well when an accurate band head spin (I_0) is predicted. The results are in good agreement with the experimentally known values of spin and transition energies. In the present paper, we have reported the band head spin of TSD bands for Lu isotope. This method brings comprehensive interpretation for spin assignment of TSD bands which could help in designing future experiments for these bands. Thus, we have reported the band head spin value of 5 TSD rotational band of Lu isotope.

1 Introduction

The study of Triaxial Super-Deformed nuclei has become an exciting field in the nuclear structure. The nuclei which has large triaxial deformation parameter γ and large quadrupole deformation ϵ_2 known as Triaxial Super-Deformed (TSD) nuclei [1, 2].

Until now, enough data for axial superdeformed nuclei are accumulated but the data for TSD nuclei are quite few like ^{163}Lu , ^{164}Lu , ^{165}Lu , ^{167}Lu , ^{171}Ta , ^{86}Zr etc. were identified experimentally. The prediction of TSD nuclei in $A \sim 160$ region has been reported in Ref. [2]. However, Bohr and Mottelson predicted the existence of the nuclear wobbling motion in rapidly rotating triaxially deformed systems [3]. The TSD nuclei has the configuration with $\epsilon_2 \sim 0.4$ and $|\gamma| \sim 20^\circ$ has been well known in a Lu isotope since long [4]. For the first time, an excited TSD band was reported in ^{163}Lu [5].

Various empirical or semi-empirical models have been proposed for the spin assignments in SD bands. Unfortunately, very few experiments could perfectly assign spins of SD bands. The theoretical modeling is only technique to calculate the spin values. Several efforts for assigning spins of SD states have been proposed in $A \sim 190$ mass region.

Further some of the methods were proposed to assign band-head spins and transition energies as a function of spin or rotational frequency ω . However, the spin assignments in these studies were not in good agreement with the experimental values. Uma et al. [6] have proposed a modified Variable Moment of Inertia (VMI) model to predict the band head spin of the superdeformed bands. In this paper a modified Variable Moment of Inertia (VMI) model is used to assign band-head spin of TSD bands directly from experimentally observed transition energy ratio which requires less computation time [7, 8]. The extension of the VMI model to an asymmetric (γ_0) rotator can only be done by an

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additional assumption. The method given is reasonable, since it reduces to the usual VMI model for ($\gamma = 0$). But the results are not changed appreciably with respect to the procedure given [9]. In this model, transition energy ratio (R) is a function of two parameters that is restoring force constant (C) and moment of inertia (\mathcal{J}_0). Both values are computed by a best-fitting method incorporating experimentally known transition energies with spin values and verified by r.m.s. deviation. It help to assign band-head spin correctly.

Thus, we organize the paper as : A brief description of theoretical and experimental efforts of superdeformed and triaxial superdeformed rotational band in nuclei is presented in the paper. In Sec 2, we briefly describe the formalism for assigning band-head spin of TSD rotational bands using modified VMI model. The results of these models are presented and compared in Sec 3 for the TSD bands. Finally Sec 4, contains a conclusion and brief remarks.

2 A brief description of the modified VMI model

Here we use modified VMI model to explain the relation of rotational energy with angular momentum (I) for each level. It is basically the sum of the potential energy term ($\mathcal{J}_I - \mathcal{J}_0$) and rotational energy term [$(\frac{\hbar^2}{2\mathcal{J}_I}).I(I + 1)$]. However, modified VMI model is a two-parameter formula which characterizes each nucleus by band-head moment of inertia (\mathcal{J}_0) and restoring force constant (C). The parameters depend on the difference in moment of inertia (\mathcal{J}_I) from that of the ground-state moment of inertia (\mathcal{J}_0). The energy level ($I_0 = 0$) of rotational band is represented as:

$$E_I(\mathcal{J}) = \frac{1}{2}C(\mathcal{J}_I - \mathcal{J}_0)^2 + \frac{1}{2} \left[\frac{I(I + 1)}{\mathcal{J}_I} \right]. \quad (1)$$

Here we assume $\hbar = 1$. While the band head energy level ($I_0 \neq 0$) of rotational band is given as:

$$E_I = E_0 + \frac{1}{2\mathcal{J}_I} [I(I + 1) - I_0(I_0 + 1)] + \frac{1}{2}C(\mathcal{J}_I - \mathcal{J}_0)^2, \quad (2)$$

where E_0 is the band-head energy of the rotational band. The variable moment of inertia \mathcal{J}_I is determined by the equilibrium condition.

$$\frac{\partial E(\mathcal{J}_I)}{\partial \mathcal{J}_I} = 0. \quad (3)$$

This is further equivalent to the cubic equation,

$$\mathcal{J}_I^3 - \mathcal{J}_I^2 \mathcal{J}_0 - ([I(I + 1) - I_0(I_0 + 1)]/2C) = 0. \quad (4)$$

The cubic equation has one real root for any finite positive value of \mathcal{J}_0 and C. Finally the transition energy for TSD bands is defined as [6]:

$$E_{\gamma(I \rightarrow I-2)} = \frac{1}{2\mathcal{J}_0} [I(I + 1) - (I - 2)(I - 1)] + \frac{1}{8C\mathcal{J}_0^4} ([I(I + 1)]^2 - [(I - 2)(I - 1)]^2). \quad (5)$$

In this equation, the parameters (\mathcal{J}_0) and (C) are determined by fitting the experimentally known transition energies using the Best-Fit Method (BFM). The band head spin is determined in terms of the observed transition energy ratio. The transition energy ratio (R) can be expressed as:

$$R = \frac{E_{\gamma(I + 2 \rightarrow I)}}{E_{\gamma(I \rightarrow I - 2)}}. \quad (6)$$

Hence the band head spin I_0 is determined from R.

The root mean square deviation (*r.m.s.*) of calculated transition energies for different I_0 values are calculated and shown in figures. The *r.m.s.* value is least for band-head spin value of a band. If I_0 value shifts away by ± 1 from the accurate value, a rapid shift in *r.m.s.* deviation (χ) can be seen. The *rms* deviation is taken as [6],

$$\chi = \left[\frac{1}{n} \sum \left| \frac{E_{\gamma}^{cal}(I_i) - E_{\gamma}^{exp}(I_i)}{E_{\gamma}^{exp}(I_i)} \right|^2 \right]^{\frac{1}{2}} \quad (7)$$

Where n is the total number of fitted transitions involved. Lowest spin I_0 is predicted from the least value of *r.m.s.*, as I_0 is known all the other spin values of the TSD band levels can be determined. It is clear that the fitting procedure is quite simple and straightforward [6]. Therefore, the experimentally observed band head spin for TSD bands are well justified with the I_0 predicted from VMI equations.

TSD Band	Stiffness Constant $C \times 10^7$ (MeV ²)	Band Head Moment of inertia (\mathcal{J}_0) (MeV ²)	Present assigned (VMI) I_0
¹⁶⁵ Lu (b1)	4.69	0.0718	12.5
¹⁶⁵ Lu (b3)	2.34	0.065	20.5
¹⁶⁷ Lu (b1)	1.63	0.0598	12.5
¹⁶⁷ Lu (b2)	5.62	0.0687	17.5
¹⁶⁷ Lu (b3)	2.32	0.0651	13.5

Table 1. The two parameters stiffness constant (C) and band head moment of inertia (\mathcal{J}_0) used in the fitting along with predicted band head spin (I_0) using modified VMI model are given.

It can be observed that a correct spin value of the lowest TSD level, I_0 , is assigned when the calculated transition energy is almost same with the observed transition energy. However, if I_0 is shifted from the correct value, even by ± 1 the *r.m.s.* deviation will no longer be least i.e. there is an increase in the *r.m.s.* deviation with the shift in band head spin. The band head spins of all the five TSD bands have been successfully predicted by this approach. Therefore, the V-shaped χ - I_0 plot gives a reasonable assignment of I_0 .

3 Results

In this section, we predicted the band head spin of Lu isotopes using modified VMI model. The fitting is successfully done for ¹⁶⁵Lu(*b1*, *b3*) and ¹⁶⁷Lu(*b1*, *b2*, *b3*) isotopes, where *b1*, *b2* and *b3* refer to band 1, band 2 and band 3 respectively. We present our results in Table 1 and 2. The experimental data are taken from Ref. [10].

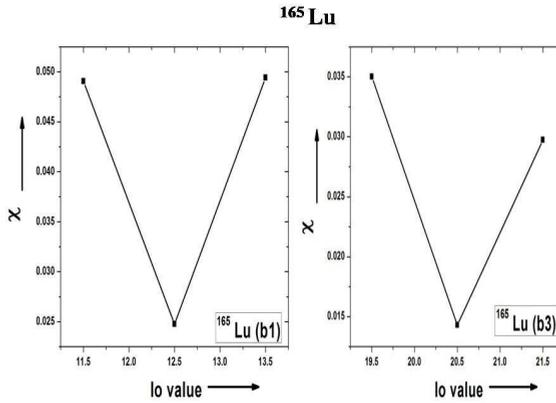


Figure 1. The r.m.s. deviation plots showing the minimum value for band head spin for ¹⁶⁵Lu

The two parameters, band head moment of inertia \mathcal{J}_0 and restoring force constant (C) are determined by fitting experimentally known transition energies using Best Fit Method (*BFM*) and the cubic equation for Lu isotopes are given in Table 1.

Using the parameters adopted in Table 1, we calculated the transition energies and with the help of calculated transition energies we are able to predict the band head spin of selected cases. The obtained gamma energies are in good agreement with the experimentally known values of energies.

TSD Band	$E_\gamma(I_4 \rightarrow I_2)$ Exp. (keV)	$E_\gamma(I_4 \rightarrow I_2)$ VMI (keV)	Band Head (I_4) Exp.	Assigned (I_4) from VMI	r.m.s. deviation (χ)
¹⁶⁵ Lu (b1)	385	392	12.5	12.5	0.0248
¹⁶⁵ Lu (b3)	661	651.4	20.5	20.5	0.0143
¹⁶⁷ Lu (b1)	471	455	12.5	12.5	0.063
¹⁶⁷ Lu (b2)	547	547.6	17.5	17.5	0.0073
¹⁶⁷ Lu (b3)	457	452.7	13.5	13.5	0.0238

Table 2. Experimental transition energies, calculated transition energies using modified VMI model along with experimental and predicted band head spin are given for 5 TSD bands for Lu

We present our results in Table 2 that gives the experimental transition energy E_γ , calculated transition energy, experimental band head spin and predicted band head spin using VMI equation. It

is also noticed that r.m.s. deviation is minimum for these bands which is shown in Figure 1 and 2 for ^{165}Lu and ^{167}Lu respectively. The predicted band head spin is compared with experimentally known spin values which provide a significant validity of this approach for TSD bands also. The results are encouraging. Overall, the obtained spin values are well justified by the r.m.s. deviation.

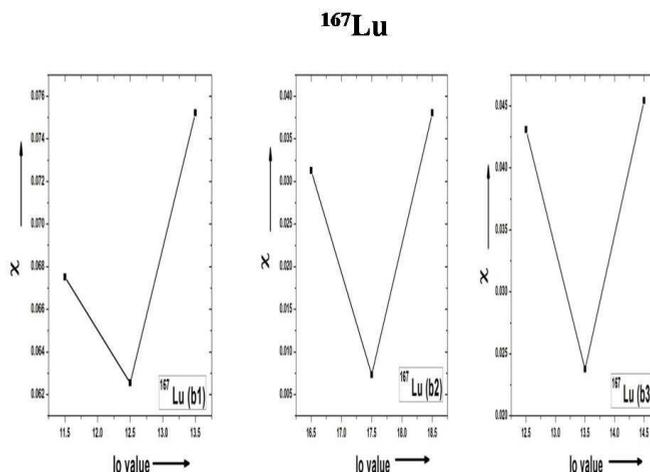


Figure 2. The r.m.s. deviation plots showing the minimum value for band head spin for ^{167}Lu

4 Conclusion

We have predicted the band head spin of 5 TSD bands in Lu isotope. Our good results for the spins and transition energies governs that the fitting parameters (C) and (\mathcal{J}_0) obtained here for these TSD bands are acceptable and the simple model employed here is appropriate to use for TSD bands. More profound analysis of TSD bands will be published in future.

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