

E0 Transitions in ^{146}Sm

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Abstract: The potential energy surfaces $V(\beta, \gamma)$ for ^{146}Sm , levels energy belonging to the ground state (g.s), beta (β), gamma (γ) bands, electromagnetic transition probabilities $B(E2)$'s, mixing ratios $\delta(E2/M1)$ and strength of the electric monopole, (E0), transitions $X(E0/E2)$ are calculated within the frame work of the interacting boson approximation model (IBA-2). The results of these calculations are compared to the available experimental data and reasonable agreement has obtained.

Keywords: Interacting boson approximation model, Levels energy, Transition probabilities, Mixing ratios, Electric monopole transitions

1. Introduction

The even-even samarium isotopes are of considerable interest because of the phase transition from spherical to deformed shape. This transition has attracted many authors to study that series of isotopes theoretically and experimentally.

The theoretical works that discussed spectral statistics of the rare-earth nuclei either using quantum mechanical effects to see the effects of the half-life of α -decay of spherical isotopes [1] or shell model configuration effects and compare to the available experimental data[2]. Also, interacting boson approximation Hamiltonian [3-6] and finite amplitude method [7] are used in calculating the energy density and the electric monopole strength in the chain of samarium isotopes.

Experimentally the lifetime of levels in ^{146}Sm has measured [8] using (^{11}B , 4n) reaction at 54 MeV on natural La target evaporated on gold. The results are compared to the calculations of the IBA model and cluster – vibrator model. The calculated values of the two models are not inconsistent with the observed transition probabilities up to $6^+ \rightarrow 4^+$ transition and failed to reproduce that for $8^+ \rightarrow 6^+$ transition.

Unfortunately scarce data are available about ^{146}Sm and the interacting boson approximation model IBA-2 has been used in calculating levels energy, transition probabilities, potential energy surfaces $V(\beta, \gamma)$, mixing ratios $\delta(E2/M1)$ and strength of the electric monopole, (E0), transitions $X(E0/E2)$.

2. Interacting boson approximation model (IBA-2)

2.1 Levels energy:

IBA-2 model [9-11] has been applied to the positive-parity of the low - lying states in ^{146}Sm . The Hamiltonian employed for the present calculation is given [12] as:

$$H = \varepsilon_{\pi} (d^{\dagger} \times d)_{\pi} + \varepsilon_{\nu} (d^{\dagger} \times d)_{\nu} + V_{\pi\pi} + V_{\nu\nu} + KQ_{\pi} \cdot Q_{\nu} + M_{\pi\nu} \quad (1)$$

Where ε_{π} , ε_{ν} are the, (proton), π and (neutron), ν energies respectively and they are assumed equal

$$\varepsilon_{\nu} = \varepsilon_{\pi} = \varepsilon$$

The third and fourth term of equation. (1) represent the π - ν interactions operator which is given by :

$$V_{\rho\rho} = \sum_{L=0,2,4} \frac{1}{2} C_{L\rho} (2L+1)^2 \left[(d^{\dagger} \times d^{\dagger})_{\rho}^L \cdot (d \times d)_{\rho}^L \right]^{(0)} \quad \rho = \pi \text{ or } \nu \quad (2)$$

The fifth term is the quadrupole operator which is given by the usual expression, (K is the strength of the π and ν bosons quadrupole interactions),:

$$Q_{\rho} = \left[(s^{\dagger} \times d) + (d^{\dagger} \times s) \right]_{\rho} + x_{\rho} (d^{\dagger} \times d)_{\rho} \quad \rho = \pi \text{ or } \nu \quad (3)$$

The sixth term is the Majorana operator and is given by :

$$M_{\Pi\nu} = -2 \sum_{k=1,3} \zeta_k (d_{\Pi}^{\dagger} x d_{\nu}^{\dagger}) \cdot (d_{\Pi} x d_{\nu}) + \zeta_2 (d_{\Pi}^{\dagger} x s_{\nu}^{\Pi} - s_{\Pi}^{\dagger} x d_{\nu}^{\dagger})^{(2)} \cdot (d_{\Pi} x s_{\nu} - s_{\Pi} x d_{\nu}) \quad (4)$$

2.2 Reduced transition probabilities, B (E2)'s:

The electric quadrupole transition operator [9] employed in this study is defined as :

$$T^{(E2)} = e_{\pi} Q_{\pi} + e_{\nu} Q_{\nu} \quad (5)$$

Where:

$T^{(E2)}$: absolute transition probability of the electric quadrupole (E2) transition,

e_{π} and e_{ν} : the π and ν effective charges, and

Q_{ρ} : the quadrupole operators which is the same as that in equation.(3)

$\rho = \pi$ or ν .

The reduced electric quadrupole transition rates between $I_i \rightarrow I_f$ states is given by

(6)

$$B(E2, I_i - I_f) = \left[\langle I_f || T^{(E2)} || I_i \rangle \right]^2 / (2I_i + 1) \quad \text{where}$$

I_i : the initial state of the electric quadrupole transition, and
 I_f : the final state of the electric quadrupole transition.

The proton and neutron boson numbers N_{π} and N_{ν} respectively can be treated as parameters. They are fixed to be half the number of valence fermions and counted from the nearest closed shell.

The effective charges e_{π} and e_{ν} depend on the total number of bosons N , number of ν bosons N_{ν} , number of π bosons N_{π} and the experimental value of $B(E2, 2_1^+ \rightarrow 0_1^+)$, where

$$N = N_{\pi} + N_{\nu} \quad (7)$$

The effective charges of the vibrational limit are calculated [13] using:

$$(N_{\pi})^{-1} [(N)(N+3)^{-1} B(E2, 2_1^+ \rightarrow 0_1^+)]^{1/2} = e_{\pi} + e_{\nu} N_{\nu} / N_{\pi} \quad (8)$$

3. Results and discussion

3.1 The potential energy surfaces:

The potential energy surfaces, $V(\beta, \gamma)$, for ^{146}Sm nucleus as a function of the deformation parameters β and γ has been calculated using [14] equation:

$$E_{N_{\nu} N_{\pi}}(\beta, \gamma) = \langle N_{\pi} N_{\nu}; \beta\gamma | H_{np} | N_{\pi} N_{\nu}; \beta\gamma \rangle = \varepsilon_d (N_{\nu} + N_{\pi}) \beta^2 (1 + \beta^2)^{-1} + \beta^2 (1 + \beta^2)^{-2} \{ k N_{\pi} N_{\nu} [4 - (\bar{X}_{\pi} + \bar{X}_{\nu}) \beta \cos 3\gamma + \bar{X}_{\pi} \bar{X}_{\nu} \beta^2] + N_{\nu}(N_{\nu}-1) \left(\frac{C_o}{10} + \frac{C_2}{7} \right) \beta^2 \} \quad (9)$$

where

$$\bar{X}_{\rho} = (2/7)^{0.5} X_{\rho} \quad \rho = \pi \text{ or } \nu \quad (10)$$

The contour plot of the potential energy surfaces is presented in Figure 1, while the plot for the potential energy surfaces versus the deformation parameters β and γ for $\gamma=0^\circ$ (prolate) and $\gamma=60^\circ$ (oblate) is presented in Figure 2. It is clear from the graphs that ^{146}Sm nucleus has vibrational character since vibrational character require:

- (a) Minimum surface energy at deformation $\beta=0$, and
- (b) Circular contours around $\beta=0$.

The two requirements are exist for the nucleus as shown in figures 1,2. The previous arguments support the assumption made by other authors that ^{146}Sm is a vibrational-like nucleus.

The value of the deformation parameter β which is corresponding to the minimum value of the potential energy implies a decrease of the moment of inertia, and in turn to an increase of the spacing between levels in the ground state and β bands.

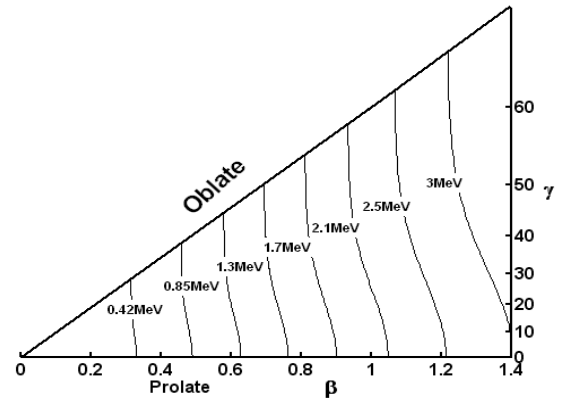


Figure 1.: Contour plot of the potential energy surfaces

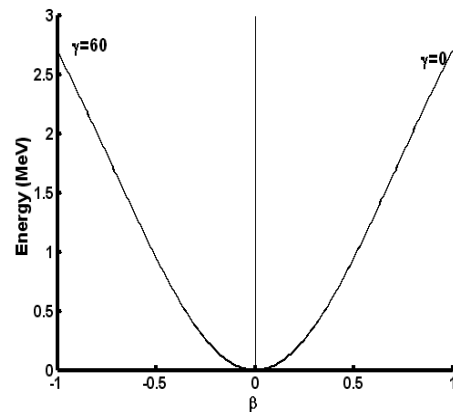


Figure 2.: The potential deformation energy (MeV) versus the deformation β . ($\gamma=0^\circ$ prolate and $\gamma=60^\circ$ oblate)

3.2 Levels energy:

The computer codes NPBOS and NPBEM [15] of the IBA-2 have been used in calculating energy eigenvalues, wave functions and E2 transition matrix elements. The Hamiltonian parameters of equation (1) used in the present calculation are:

$$\begin{aligned} \varepsilon &= 0.890 & K &= -0.080 & X_{\parallel} &= -1.300 & X_{\nu} &= -1.400 \\ \varepsilon_1 = \varepsilon_3 &= 0.090 & \varepsilon_2 &= 0.100 & & & & \\ C_{L\nu} &= 0.000, 0.050, 0.000 & & & & & & \\ C_{L\pi} &= 0.000, 0.100, 0.000 & & & & & & \text{(all in MeV)}. \end{aligned}$$

The results of the calculation of energy levels together with the experimental data are displayed in **Figure 3**.

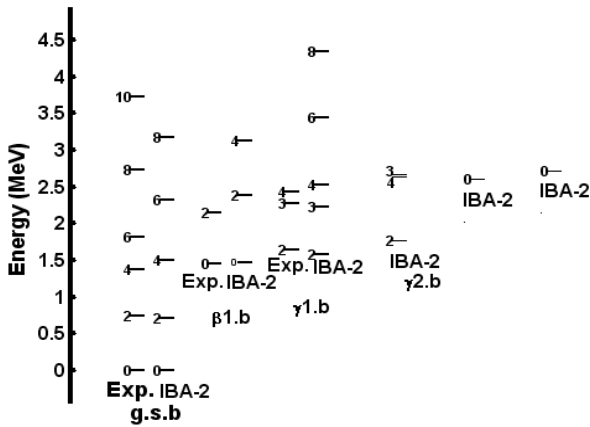
The calculation is limited to states of spin- parity up to I=8⁺ and the energies of seventeen levels were reproduced and in general, reasonable agreement has achieved between the experimentally measured [16] and the calculated ones .

3.3 Reduced transition probabilities, B (E2)'s :

The π and ν effective charges e_{\parallel} and e_{ν} were calculated using the vibrational limit [13]. The calculated values are :

$$e_{\parallel} = 0.231 \text{ e.b and } e_{\nu} = 0.531 \text{ e.b}$$

The calculated and the experimental B(E2) values are listed and compared in **Table 1**. There is only one experimental value available [17] which is B(E2,0₁⁺-2₁⁺)= 0.48(8) e²b². It **Figure 4**: Comparison between the experimental and



theoretical energy levels

is based on the life time measurement of 747.1 keV level which is < 7.2 ps.

The quasi ground state , β and γ – bands are presented in spectra columns as in Figure 4. The calculated transition probability of the internal gamma rays in any band have greater values than the gamma transitions in inter bands have, **Table 1**.

These energy and ratios of different energies to those of the first excited 2₁⁺ state are listed in Table 2 and show vibrational character to ¹⁴⁶Sm nucleus.

Table 1 : Reduced transition probability values , B (E2) (in e²b²)

I _i ⁺	I _f ⁺	B(E2) Exp*	B(E2) IBA-2	I _i ⁺	I _f ⁺	B(E2) Exp*	B(E2) IBA-2
2 ₁	0 ₁	0.090	0.1065	3 ₂	4 ₁		0.0082
2 ₁	0 ₂		0.0321	3 ₂	4 ₂		0.0515
2 ₁	0 ₃		0.0012	3 ₂	4 ₃		0.0019
2 ₁	0 ₄		0.0001	3 ₂	4 ₄		0.0088
2 ₂	0 ₁		0.0015	4 ₁	2 ₁		0.1718
2 ₂	0 ₂		0.0109	4 ₁	2 ₂		0.0008
2 ₂	0 ₃		0.0067	4 ₁	2 ₃		0.0233
2 ₂	0 ₄		0.0095	4 ₁	2 ₄		0.0247
2 ₃	0 ₁		0.0083	4 ₂	2 ₁		0.0003
2 ₃	0 ₂		0.0383	4 ₂	2 ₂		0.0649
2 ₃	0 ₃		0.0000	4 ₂	2 ₃		0.0270
2 ₃	0 ₄		0.0118	4 ₂	2 ₄		0.0130
2 ₄	0 ₁		0.0003	4 ₃	2 ₁		0.0099
2 ₄	0 ₂		0.0650	4 ₃	2 ₂		0.0005
2 ₄	0 ₃		0.0058	4 ₃	2 ₃		0.0327
2 ₄	0 ₄		0.0132	4 ₃	2 ₄		0.0160
2 ₁	2 ₂		0.1058	4 ₄	2 ₁		0.0001
2 ₁	2 ₃		0.0111	4 ₄	2 ₂		0.0021
2 ₁	2 ₄		0.0000	4 ₄	2 ₃		0.0006
2 ₂	2 ₁		0.1058	4 ₄	2 ₄		0.0187
2 ₂	2 ₃		0.0243	4 ₁	3 ₁		0.0311
2 ₂	2 ₄		0.0319	4 ₁	3 ₂		0.0063
2 ₃	2 ₁		0.0111	4 ₂	3 ₁		0.0140
2 ₃	2 ₂		0.0243	4 ₂	3 ₂		0.0401
2 ₃	2 ₄		0.0134	4 ₃	3 ₁		0.00 83
2 ₄	2 ₁		0.0000	4 ₃	3 ₂		0.0015
2 ₄	2 ₂		0.0319	4 ₄	3 ₁		0.0607
2 ₄	2 ₃		0.0134	4 ₄	3 ₂		0.0068
3 ₁	2 ₁		0.0039	4 ₂	4 ₁		0.0589
3 ₁	2 ₂		0.1462	4 ₃	4 ₁		0.0000
3 ₁	2 ₃		0.0074	4 ₃	4 ₂		0.0034
3 ₁	2 ₄		0.0056	4 ₄	4 ₃		0.0250
3 ₂	2 ₁		0.0034	4 ₄	4 ₂		0.0483
3 ₂	2 ₂		0.0161	4 ₄	4 ₁		0.0006
3 ₂	2 ₃		0.0545	6 ₁	4 ₁		0.2080
3 ₂	2 ₄		0.0585	6 ₁	4 ₃		0.0247
3 ₁	4 ₂		0.0180	6 ₁	4 ₄		0.0095
3 ₁	4 ₃		0.0107	6 ₂	4 ₁		0.0006

[17]

Table 2: The levels energy ratios

Quantity	Exp*	IBA-2	SU(5)	O(6)	SU(3)
E ₂ ⁺ /E ₂ ¹⁺	2.20	2.21	2.00	2.50	3.33
E ₄ ¹⁺ /E ₂ ¹⁺	1.84	2.09	2.00	2.50	3.33
E ₆ ¹⁺ /E ₂ ¹⁺	2.42	3.24	3.00	4.50	7.00
E ₄ ²⁺ /E ₂ ¹⁺	3.05	3.53	3.00	4.50	7.00
E ₃ ¹⁺ /E ₂ ¹⁺	3.03	3.11	3.00	4.50	7.00
E ₈ ¹⁺ /E ₂ ¹⁺	3.66	4.45	4.00	7.00	12.00
E ₄ ³⁺ /E ₂ ¹⁺	-	3.70	4.00	7.00	12.00

[18]

SU (5) = vibrational limit,
O (6) = gamma unstable limit, and
SU (3) = rotational limit.

3.4 Multipole mixing ratio $\delta(E 2/M 1)$:

The reduced electric E2 and magnetic M1 matrix elements of different gamma ray transitions were calculated and used in calculating multipole mixing ratios using [19]

$$\delta(E2/M1) = 0.835E_\gamma(MeV) \frac{\langle I_f || T^{E2} || I_i \rangle}{\langle I_f || T^{M1} || I_i \rangle} \quad (11)$$

The calculated values are presented in Table 3 in comparison to the available experimental ones and reasonable agreement has achieved.

Table 3: Mixing ratios (δ) for ^{146}Sm

I_i^+	I_f^+	Exp.(*)	IBA-2
2 ₂	3 ₁	0.33±0.04	0.47
2 ₃	3 ₁		-2.42
3 ₁	2 ₂		-1.47
3 ₂	2 ₃		2.82
3 ₂	3 ₁		1.87
3 ₂	4 ₁	-0.10 ⁺²⁰ ₋₂₆	0.95
4 ₄	4 ₂	0.79 ⁺²⁹ ₋₂₄	1.27

*[16]

3.5 Strength of the electric monopole transitions :

The electric monopole transitions, E0, are normally occurring between two states of the same spin and parity by transferring energy and zero unit of angular momentum. It is a pure penetration effect where it caused by the electromagnetic interaction between the nuclear charge and the atomic electron which penetrates the nucleus. The E0 strength can be considered as the ratio between the reduced transition probability of competing E0 and electric quadrupole, E2, transitions de-populating the same level. The strength of the electric monopole transition, (E0 / E2), can be determined by [20]:

$$X_{if} (E0/E2) = \frac{B(E0, I_i - I_f)}{B(E2, I_i - I_f)} \quad (12)$$

$$X_{if} (E0/E2) = 2.54 \times 10^9 \chi A^{4/3} \frac{(E_\gamma MeV)^5}{\Omega_{K,L}} \quad (13)$$

$$\alpha(E2) = \frac{T_e(E0, I_i - I_f)}{T_e(E2, I_i - I_f)}$$

where

- B(E0, I_i - I_f) : reduced transition probability of the E0 transition,
- B(E2, I_i - I_f) : reduced transition probability of the E2 transition.
- A : mass number ,
- I_i : spin of the initial state where E0 & E2 transition are depopulating it,
- I_f : spin of the final state of E0 transition,
- I_f' : spin of the final state of E2 transition,
- E_γ : gamma ray energy,
- Ω_{kl} : electronic factor for K,L shells [21],
- α(E2) : conversion coefficient of the E2 transition,

- T_e(E0, I_i-I_f): absolute transition probability of the E0 transition between I_i and I_f states, and
- T_e(E2, I_i-I_f) : absolute transition probability of the E2 transition between I_i and I_f states.

The calculated strength of ten electric monopole transitions for ΔI[±]=0 , I_i[±]=0, I_f[±]=0 and I_f[±] =2 are presented in Table 4. The large value for (0₃⁺ - 0₂⁺ - 2₃⁺) may be due to the vanishing B(E2) value rather than the large B(E0) value. Unfortunately there is no experimental data for comparison.

Table 4. X_{iff'}(E0/E2) ratios for E0 transitions

I_i^+	I_f^+	$I_{f'}^+$	Exp	IBA-2	I_i^+	I_f^+	$I_{f'}^+$	Exp	IBA-2
0 ₂	0 ₁	2 ₁	0.0057		0 ₄	0 ₃	2 ₂		0.0010
0 ₃	0 ₂	2 ₃	2.9032		0 ₄	0 ₃	2 ₁		0.1081
0 ₃	0 ₂	2 ₂	0.0431		0 ₄	0 ₂	2 ₄		0.0034
0 ₃	0 ₂	2 ₁	0.2439		0 ₄	0 ₂	2 ₃		0.0038
0 ₃	0 ₁	2 ₃	1.3548		0 ₄	0 ₂	2 ₂		0.0047
0 ₃	0 ₁	2 ₂	0.0201		0 ₄	0 ₂	2 ₁		0.3783
0 ₃	0 ₁	2 ₁	0.1138		0 ₄	0 ₁	2 ₄		0.0012
0 ₄	0 ₃	2 ₄	0.0009		0 ₄	0 ₁	2 ₃		0.0013
0 ₄	0 ₃	2 ₃	0.0011		0 ₄	0 ₁	2 ₂		0.0016
0 ₄	0 ₃	2 ₂	0.0013		0 ₄	0 ₁	2 ₁		0.1351

4. Conclusions

1. The contour plot of the potential energy surfaces V(β, γ) shows that ^{146}Sm is a spherical nucleus and has a vibrational character.
2. The vibrational limit of the IBA-2 has been used and energies of seventeen-levels were calculated.
3. The reduced transition probability of gamma transitions were calculated and normalized to B(E2, 0₁⁺→2₁⁺) = 0.48(8).
4. The mixing ratios and the strength of twenty electric monopole (E0) transitions was calculated for the first time.

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