X(5) Symmetry to ¹⁵²Sm

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The excited positive and negative parity states, potential energy surfaces, $V(\beta, \gamma)$, electromagnetic transition probabilities, B(E1), B(E2), electric monopole strength X(E0/E2) and staggering effect, $\Delta I = 1$, were calculated successfully using the interacting boson approximation model *IBA*-1. The calculated values are compared to the available experimental data and show reasonable agreement. The energy ratios and contour plot of the potential energy surfaces show that ¹⁵²Sm is an X(5) candidate.

1 Introduction

Phase transition is one of the very interesting topic in nuclear structure physics. The even-even samarium series of isotopes have encouraged many authors to study that area extensively experimentally and theoretically.

Experimentally, authors studied levels energy with their half-lives, transition probabilities, decay schemes, multipole mixing ratios, internal conversion coefficients, angular correlations and nuclear orientation of γ -rays[1-4].

Theoretically, different theoretical models have been applied to that chain of isotopes. One of the very interesting models is the interacting boson approximation model *IBA* [5-10]. Iachello [11,12] has made an important contribution by introducing the new dynamical symmetries E(5) and X(5).

E(5) is the critical point symmetry of phase transition between U(5) and O(6) while X(5) is between U(5) and SU(3)nuclei. The aim of the present work is to calculate:

- 1. The potential energy surfaces, $V(\beta, \gamma)$;
- 2. The levels energy, electromagnetic transition rates *B*(*E*1) and *B*(*E*2);
- 3. The staggering effect, and
- 4. The electric monopole strength X(E0/E2).

2 IBA-1 model

2.1 Levels energy

The *IBA*-1 Hamiltonian [13-16] employed on 152 Sm in the present calculation is:

$$H = EPS \cdot n_d + PAIR \cdot (P \cdot P)$$

+ $\frac{1}{2} ELL \cdot (L \cdot L) + \frac{1}{2} QQ \cdot (Q \cdot Q)$ (1)

$$+5 OCT \cdot (T_3 \cdot T_3) + 5 HEX \cdot (T_4 \cdot T_4),$$

where

170

$$P \cdot p = \frac{1}{2} \begin{bmatrix} \left\{ (s^{\dagger} s^{\dagger})_{0}^{(0)} - \sqrt{5} (d^{\dagger} d^{\dagger})_{0}^{(0)} \right\} x \\ \left\{ (ss)_{0}^{(0)} - \sqrt{5} (\tilde{d} \tilde{d})_{0}^{(0)} \right\} \end{bmatrix}_{0}^{(0)},$$
(2)

$$L \cdot L = -10 \sqrt{3} \left[(d^{\dagger} \tilde{d})^{(1)} x (d^{\dagger} \tilde{d})^{(1)} \right]_{0}^{(0)}, \qquad (3)$$

$$Q \cdot Q = \sqrt{5} \left[\begin{cases} (S^{\dagger} \tilde{d} + d^{\dagger} s)^{(2)} - \frac{\sqrt{7}}{2} (d^{\dagger} \tilde{d})^{(2)} \\ \\ \left\{ (s^{\dagger} \tilde{d} + + \tilde{d} s)^{(2)} - \frac{\sqrt{7}}{2} (d^{\dagger} \tilde{d})^{(2)} \\ \end{cases} \right]_{0}^{(0)}, \quad (4)$$

$$T_3 \cdot T_3 = -\sqrt{7} \left[(d^{\dagger} \tilde{d})^{(2)} x (d^{\dagger} \tilde{d})^{(2)} \right]_0^{(0)} , \qquad (5)$$

$$T_4 \cdot T_4 = 3 \left[(d^{\dagger} \tilde{d})^{(4)} x (d^{\dagger} \tilde{d})^{(4)} \right]_0^{(0)} .$$
(6)

In the previous formulas, n_d is the number of bosons; $P \cdot P$, $L \cdot L$, $Q \cdot Q$, $T_3 \cdot T_3$ and $T_4 \cdot T_4$ represent pairing, angular momentum, quadrupole, octupole and hexadecupole interactions respectively between the bosons; *EPS* is the boson energy; and *PAIR*, *ELL*, *QQ*, *OCT*, *HEX* are the strengths of the pairing, angular momentum, quadrupole, octupole and hexadecupole interactions respectively (see Table 1).

2.2 Transition rates

The electric quadrupole transition operator employed is:

$$\Gamma^{(E2)} = E2SD \cdot (s^{\dagger}\tilde{d} + d^{\dagger}s)^{(2)} + \frac{1}{\sqrt{5}}E2DD \cdot (d^{\dagger}\tilde{d})^{(2)}.$$
 (7)

E2SD and E2DD are adjustable parameters.

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

$$B(E_2, I_i - I_f) = \frac{[\langle I_f || T^{(E_2)} || I_i \rangle]^2}{2I_i + 1}.$$
(8)

3 Results and discussion

In this section we review and discuss the results.

Salah A. Eid and Sohair M. Diab. X(5) Symmetry to ¹⁵²Sm

nucleus	EPS	PAIR	ELL	QQ	OCT	HEX	E2SD(eb)	E2DD(eb)
¹⁵² Sm	0.3840	0.000	0.0084	-0.0244	0.0000	0.0000	0.1450	-0.4289

Table 1: Parameters used in IBA-1 Hamiltonian (all in MeV).

3.1 The potential energy surfaces

The potential energy surfaces [17], $V(\beta, \gamma)$, as a function of the deformation parameters β and γ are calculated using:

$$E_{N_{\Pi}N_{\nu}}(\beta,\gamma) = \langle N_{\pi}N_{\nu};\beta\gamma | H_{\pi\nu} | N_{\pi}N_{\nu};\beta\gamma\rangle =$$

$$= \zeta_{d}(N_{\nu}N_{\pi})\beta^{2}(1+\beta^{2}) + \beta^{2}(1+\beta^{2})^{-2}\times$$

$$\times \left\{ kN_{\nu}N_{\pi}[4-(\bar{X}_{\pi}\bar{X}_{\nu})\beta\cos 3\gamma] \right\} +$$

$$+ \left\{ [\bar{X}_{\pi}\bar{X}_{\nu}\beta^{2}] + N_{\nu}(N_{\nu}-1)\left(\frac{1}{10}c_{0}+\frac{1}{7}c_{2}\right)\beta^{2} \right\},$$
(9)

where

$$\bar{X}_{\rho} = \left(\frac{2}{7}\right)^{0.5} X_{\rho} \rho = \pi \text{ or } \upsilon.$$
 (10)

The calculated potential energy surfaces, $V(\beta, \gamma)$, are presented in Figures 1, 2, 3. ¹⁵²Sm lies between ¹⁵⁰Sm which is a vibrational like nucleus, U(5), Fig. 1, while ¹⁵⁴Sm is a rotational like, SU(3), nucleus, Fig. 3. So, ¹⁵⁰Sm can be an X(5)candidate where levels energy, transition probability ratios as well as the potential energy surfaces are supporting that assumption (see Table 2).

3.2 Energy spectra and electric transition rates

The energy of the positive and negative parity states of ¹⁵²Sm isotope are calculated using computer code PHINT [19]. A comparison between the experimental spectra [18] and our calculations, using values of the model parameters given in Table 1 for the ground state, $\beta 1$, $\beta 2$ and γ bands are illustrated in Fig. 4. The agreement between the calculated levels energy and their corresponding experimental values are fair, but they are slightly higher especially for the higher excited states in $\beta 1$, $\beta 2$ and γ bands. We believe this is due to the change of the projection of the angular momentum which is due mainly to band crossing. Fig. 5 shows the position of X(5) and E(5) between the other types of nuclei.

Unfortunately there are no available measurements of electromagnetic transition rates B(E1) for ¹⁵²Sm nucleus, Table 3, while some of B(E2) are measured. The measured $B(E2, 2_1^+ \rightarrow 0_1^+)$ is presented, in Table 4, for comparison with the calculated values [20]. The parameters E2SD and E2DDdisplayed in Table 1 are used in the computer code NPBEM [19] for calculating the electromagnetic transition rates and the calculated values are normalized to $B(E2, 2_1^+ \rightarrow 0_1^+)$. No new parameters are introduced for calculating electromagnetic transition rates B(E1) and B(E2) of intraband and interband.



Fig. 1: Potential energy surfaces for ¹⁵⁰Sm .



Fig. 2: Potential energy surfaces for ¹⁵²Sm .



Fig. 3: Potential energy surfaces for ¹⁵⁴Sm .

nucleus	$E_{4_1^+}/E_{2_1^+}$	$E_{6_1^+}/E_{2_1^+}$	$E_{8_1^+}/E_{2_1^+}$	$E_{0_2^+}/E_{2_1^+}$	$E_{6_1^+}/E_{0_2^+}$	$E_{0_3^+}/E_{2_1^+}$	$BE2(4_1^+ - 2_1^+)/BE2(2_1^+ - 0_1^+)$
¹⁵² Sm	3.02	5.83	9.29	5.66	1.03	8.92	1.53
X(5)	3.02	5.83	9.29	5.65	1.53	6.03	1.58

Table 2: Energy and transition probability ratios.



Fig. 4: Experimental[18] and calculated levels energy



Fig. 5: Triangle showing the position of X(5) and E(5).



Fig. 6: Staggering effect on ¹⁵²Sm.

$I_i^- I_f^+$	<i>B</i> (<i>E</i> 1)Exp.	<i>B</i> (<i>E</i> 1)IBA-1
$1_1 0_1$		0.0979
$1_1 0_2$		0.0814
31 21		0.2338
31 22		0.0766
31 23		0.0106
$3_2 2_1$		0.0269
$3_2 2_2$		0.0291
$3_2 2_3$		0.0434
51 41		0.3579
5 ₁ 4 ₂		0.0672
5 ₁ 4 ₃	<u> </u>	0.0050
71 61		0.4815
71 62		0.0574
9 ₁ 8 ₁		0.6075
91 82		0.0490
111 101		0.7367
111 102		0.0413

Table 3: Calculated B(E1) in ¹⁵²Sm.

3.3 Staggering effect

The presence of (+ve) and (-ve) parity states has encouraged us to study the staggering effect [21-23] for ¹⁵²Sm isotope using staggering function equations (11, 12) with the help of the available experimental data [18].

$$St(I) = 6\Delta E(I) - 4\Delta E(I-1) - 4\Delta E(I+1) + + \Delta E(I+2) + \Delta E(I-2),$$
(11)

with

$$\Delta E(I) = E(I+1) - E(I).$$
 (12)

The calculated staggering patterns are illustrated in Fig. 6 and show an interaction between the (+ve) and (-ve) parity states for the ground state band of ¹⁵²Sm.

3.4 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. The strength of the electric monopole transition, $X_{if'f}(E0/E2)$, [24] can be calculated using equations (13, 14) and presented in Table 5.

$$X_{if'f}(E0/E2) = \frac{B(E0, I_i - I_f)}{B(E2, I_i - I_{'f})},$$
(13)

Salah A. Eid and Sohair M. Diab. X(5) Symmetry to ¹⁵²Sm

$I_i^+ I_f^+$	<i>B</i> (<i>E</i> 2)Exp*.	<i>B</i> (<i>E</i> 2)IBA-1
$2_1 0_1$	0.670(15)	0.6529
31 21		0.0168
41 21	0.1.017(4)	1.0014
61 41	1.179(33)	1.1304
02 21	0.176(1)	0.3363
$2_2 2_1$	0.0258(26)	0.0610
22 41	0.091(11)	0.1057
42 21	0.0035(35)	0.0003
42 41	0.037(23)	0.0458
$2_3 0_1$	0.0163(11)	0.0141
23 21	0.0417(42)	0.0125
23 41	0.0416(32)	0.0296
43 21	0.0035(13)	0.0038
43 41	0.037(13)	0.0084
43 42		0.1235
43 22		0.0070
43 23		0.3110
42 22		0.6418
81 61		1.1681
81 62		0.0376
101 81		1.1421

Table 4: Calculated B(E2) in ¹⁵²Sm (* from Ref.[20])

$I_i^+ I_f^+$	X(E0/E2)Exp*.	X(E0/E2)IBA-1
02 01	0.7(0.1)	0.85
$0_3 0_2$		3.68
$0_3 0_1$		0.72
04 03		4.39
04 02		0.64
04 01		1.27
$2_2 2_1$	4.5(0.5)	3.52
$2_3 2_1$		12.23
$2_3 2_2$		11.19
43 41		1.76
43 42		1.40
44 41		0.44
44 42		3.15
42 41	6.6(2.10)	2.02
62 61		1.46
8 ₂ 8 ₁		1.20
102 101		1.07

Table 5: $X_{if'f}$ (E0/E2) ratios in ¹⁵²Sm (* from Ref [20]).

Salah A. Eid and Sohair M. Diab. X(5) Symmetry to ¹⁵²Sm

where
$$I_i = I_f = 0$$
, $I_{f} = 2$ and $I_i = I_f \neq 0$, $I_f = I_{f}$

$$X_{if'f}(E0/E2) = (2.54 \times 10^9) A^{3/4} \times \frac{E_{\gamma}^5 (\text{MeV})}{\Omega_{KL}} \alpha(E2) \frac{T_e(E0, I_i - I_f)}{T_e(E_2, I_i - I_f)}.$$
 (14)

where:

A : mass number;

 I_i : spin of the initial state where E0 and E2 transitions 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 [25];

 $\alpha(E2)$: 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

 $T_e(E_2, I_i - I_f)$: absolute transition probability of the *E*2 transition between I_i and I_f states.

3.5 Conclusions

The *IBA*-1 model has been applied successfully to the 152 Sm isotope and:

- 1. Levels energy are successfully reproduced;
- 2. Potential energy surfaces are calculated and show X(5) characters to ¹⁵²Sm;
- 3. Electromagnetic transition rates *B*(*E*1) and *B*(*E*2) are calculated;
- 4. Staggering effect has been calculated and beat pattern observed which show an interaction between the (-ve) and (+ve) parity states, and
- 5. Strength of the electric monopole transitions $X_{if'f}(E0/E2)$ are calculated.

Submitted on January 10, 2016 / Accepted on January 12, 2016

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