A Study of Some Nuclear Properties of a Nuclei with Mass Numbers (A=100,102) using Interacting Boson Model

*A.R.H Subber **Mohammed A. Al Sharefi ***Ahmed M. Shuaib Department of Physics – College of Education – Basrah university Department of Physics – College of Science – Babylon university

Abstract :

In this study, we have employed the Interacting Boson Model (IBM-1 and IBM-2) to determine the most appropriate Hamiltonian for the study of a nuclei in the region (A \approx 100). Using the best fit values of parameters to construct the Hamiltonian, we have estimated energy levels and multipole mixing ratios (δ (E2/M1)) for some doubly even (A=100,102) nuclei. The results are compared with experimental data and it is observed that they are in good agreement.

الخلاصة:

. تم في هذا البحث استعمال انموذجي البوزونات المتفاعلة (الاول و الثاني) لتحديد المؤثر الهاملتوني الأكثر ملائمة لدراسة النوى في المنطقة (100≈A) باستعمال افضل تطابق للبارامترات في المؤثر الهاملتوني, وقد تم حساب مستويات الطاقة ونسب الخلط لبعض النوى ذات العدد الكتلي (A=100,102) , وتمت مقارنة النتائج مع القيم العملية وكانت النتائج في هذا البحث ذات توافق جيد .

1-Introduction :

The study of nuclear structure concerns itself with the motions of nucleons within the nucleus. In principle what is needed is the total wavefunction of the nucleus. However, for all but the lightest nuclei this becomes extremely complicated and of little practical use. As the problem becomes intractable with increasing mass, one has to resort to nuclear models which make use of simplifications and approximations [1]. These models must be simple enough to understand and calculate, but still powerful enough to predict observed phenomena. Some of the most successful nuclear models are the Collective Model of Bohr and Mottlesohn , and its group theory counterpart, the Interacting Boson Model (IBM) of Arima and Iachello . Both models have special limiting cases that form structural paradigms, and which are empirically observed in select nuclei. It is therefore important to subject these limiting cases to extreme scrutiny, as they act as benchmarks of nuclear structure.[2]

In previous years, some nuclei are calculated by (IBM) such as : Yazar et al. [3] explore the energy levels and the electric quadrupole transition probabilities $B(E2;I_i \rightarrow I_f)$ and gamma-ray E2/M1 mixing ratios for selected transitions of some even-even erbium isotopes, Zamfir et al. [4] stated that Xe isotopes of the mass region of A ~ 130 appear to evolve from U(5) to O(6)-like structure in the IBM.

2- Theoretical Basis :

2-1 (IBM-1) :

The IBM-1 Hamiltonian can be written as : [5]

$$H = \epsilon'' \hat{n}_d + a_0 P^+ P + a_1 \hat{L}^2 + a_2 Q^2 + a_3 T_3^2 + a_4 T_4^2,$$
(1)
Where :
$$P = 1/2 (d^2 - s^2)$$
$$Q = (s^+ \cdot d + d^+ \cdot s) + \chi (d^+ \cdot d)$$
$$n_d = \sqrt{5} T_o \qquad L = \sqrt{10} T$$

Journal of Kerbala University, Vol. 10 No.1 Scientific . 2012

The program code PHINT written by Scholten is used within the option of specifying the parameters in the neutron proton formalism where it take care of projecting maximum symmetry states. We have used EPS, PAIR, ELL, QQ parameters and kept OCT, HEX as zero since there contribution is very small. For these calculation the experimental data are taken from .

2-2 (IBM-2):

The microscopic picture of the IBM is very complicated. A commonly used microscopic picture is given in terms of collective pairs of nucleons. The S and D pairs of valence nucleons have angular momenta J = 0 and J = 2, respectively. These pairs correspond intuitively to the s and d bosons, respectively. The building blocks of the IBM-2 are the proton bosons s_{π} , d_{π} and the neutron bosons s_{ν} , d_{ν} . For the analysis of excitation energies in some (A=100,102) nuclei we tried to keep to a minimum number of free parameters in the Hamiltonian [6]. We thus considered equal values for the neutron and proton d-boson excitation energy, in addition to the standard quadrupole interaction and Majorana term. We only considered the dipole neutron-proton boson interaction whose strength is characterized by a single parameter $W_{\pi\nu}$. The explicit expression of the Hamiltonian adopted in the calculations is : [5,7]

$$\hat{H} = \varepsilon (\hat{n}_{d_{\pi}} + \hat{n}_{d_{\nu}}) + K_{\pi\nu} \hat{Q}_{\pi} \cdot \hat{Q}_{\nu} + K_{\pi\pi} \hat{Q}_{\pi} \cdot \hat{Q}_{\pi} + w_{\pi\nu} \hat{L}_{\pi} \cdot \hat{L}_{\nu} + \hat{M}_{\pi\nu} , \qquad (2)$$

where the indexes $n_{d\pi}$ and $n_{d\nu}$ refer to neutron and proton bosons respectively, and $M_{\pi\nu}$ is the Majorana term. Moreover,

$$\begin{split} \hat{n}_{d_{\eta}} &= (\hat{d}_{\eta}^{+} \cdot \tilde{d}_{\eta}) \,, \\ \hat{Q}_{\eta} &= [\hat{d}_{\eta}^{+} \times \hat{\tilde{s}}_{\eta} + \hat{s}_{\eta}^{+} \times \hat{\tilde{d}}_{\eta}]^{(2)} + \chi_{\eta} [\hat{d}_{\eta}^{+} \times \hat{\tilde{d}}_{\eta}]^{(2)} \,, \\ \hat{L}_{\eta} &= \sqrt{10} \, [\hat{d}_{\eta}^{+} \times \hat{d}_{\eta}]^{(1)} , \qquad \eta = \pi, \nu \,, \\ \hat{M}_{\pi\nu} &= \frac{1}{2} \xi_{2} [\hat{s}_{\nu}^{+} \times \hat{d}_{\pi}^{+} - \hat{s}_{\pi}^{+} \times \hat{d}_{\nu}^{+}]^{(2)} \cdot [\hat{s}_{\nu} \times \hat{\tilde{d}}_{\pi} - \hat{\tilde{s}}_{\pi} \times \hat{\tilde{d}}_{\nu}]^{(2)} \\ &- \sum_{k=1,3} \xi_{k} [\hat{d}_{\nu}^{+} \times \hat{d}_{\pi}^{+}]^{(k)} \cdot [\hat{\tilde{d}}_{\nu} \times \hat{\tilde{d}}_{\pi}]^{(k)} \,. \end{split}$$

The E2 transition strengths were calculated by using the operator :

$$\hat{T}(E2) = e_\pi \hat{Q}_\pi + e_\nu \hat{Q}_\nu \,,$$

where e_{π} and e_v are boson effective charges. The M1 operator is given by :[7,9]

(3)

$$\hat{T}(M1) = \sqrt{\frac{3}{4\pi}} (g_{\pi}\hat{L}_{\pi} + g_{\nu}\hat{L}_{\nu}), \qquad (4)$$

where g_{π} and g_{ν} are the proton and neutron boson g-factors, respectively. The numerical diagonalization has been carried out by using the computer code NPBOS of Otsuka.

The mixing ratio is given by :

$$\delta(E2/M1) = 0.835 E_{\gamma} \Delta(E2/M1)$$
 (5)

3- Results and Discussions :

As a starting point we used parameters extrapolated from the IBM calculations of some (A=100,102) nuclei , the fitted values of the parameters are given in Tables (1 and 2). These parameters are consistent with those of previous calculations. In general, at using IBM-2, we find the values of ε and κ decrease with increasing mass number.

Also, we can know the limit of nuclei (vibrational, rotational or γ -unstable) by the parameters of IBM versions. Also, we have estimated multipole mixing ratios (δ (E2/M1)) of some transitions for this nuclei and then compared them with some experimental results from [9,10], The mixing ratios values are given in the tables (3-5).

•~	500).								
	Nucleus	EPS	P.P.	L.L.	Q.Q.	T ₃ . T ₃	T ₄ .T ₄	CHI	SO6
	¹⁰⁰ Sr	0.000	0.000	0.0160	- 0.0091	0.000	0.000	- 1.320	1.000
	¹⁰⁰ Mo	0.3610	0.000	0.0099	0.000	0.0581	0.0112	0.000	1.000
	¹⁰² Zr	0.000	0.000	0.0200	- 0.0078	0.000	0.000	- 1.320	1.000
	¹⁰² Pd	0.5780	0.000	0.0115	- 0.020	0.000	0.000	0.000	1.000
	¹⁰² Ru	0.1057	0.000	0.0191	0.000	0.0513	0.1016	0.000	1.000
	¹⁰² Mo	0.1978	0.0850	0.0181	0.000	0.0636	0.000	0.000	1.000

Table 1 . The Hamiltonian parameters set used in the present study for the IBM-1 calculations of
some nuclei with mass numbers (A=100,102).(all in MeV except CHI
(all in MeV except CHIand SO6) ..

Table 2 . The Hamiltonian parameters set used in the present study for the IBM-2 calculations of some nuclei with mass numbers (A=100,102)

Nucleus	N_{π}	N _v	3	κ	χ_{ν}	χπ	C_{Lv} (L=0,2,4)	$C_{L\pi}(L=0,2,4)$
			(MeV)	(MeV)			(MeV)	(MeV)
¹⁰⁰ Sr	6	6	0.684	- 0.178	0.18	- 1.05	0.93,-0.67,0	0.82,0,0
¹⁰⁰ Mo	4	4	0.824	- 0.058	- 1.00	- 1.00	- 0.55,- 0.1,0.1	0,-0.5,0
¹⁰² Zr	5	6	0.523	- 0.121	0.32	- 0.26	0,- 0.48,0	0.6,0,0
¹⁰² Pd	2	3	0.750	- 0.084	- 1.10	0.60	0.3,- 0.4,0	0.2,-0.65,0
¹⁰² Ru	3	4	0.741	- 0.079	- 0.50	- 0.60	0.2, 0,0	-0.85,-0.66,0
¹⁰² Mo	4	5	0.620	- 0.060	-1.00	- 1.00	-0. 1,0.05,0.1	-0.3,-0.1,0

Table 3 . The mixing ratios values of $~^{100}\mathrm{Sr}$ and $~^{100}\mathrm{Mo}$

$I_i {\rightarrow} I_f$	¹⁰⁰ Sr		$I_i {\rightarrow} I_f$	¹⁰⁰ Mo	
	Exp.	IBM-2		Exp.	IBM-2
$2_1^+ \rightarrow 2_2^+$		6.3172	$2_1^+ \rightarrow 2_2^+$		1.2470
$2_1^+ \rightarrow 2_3^+$		0.1219	$2_1^+ \rightarrow 2_3^+$		0.4259
$4_1^+ \rightarrow 4_2^+$		14.8295	$4_1^+ \rightarrow 4_2^+$		2.9090
$3_1^+ \rightarrow 2_1^+$		6.4731	$3_1^+ \rightarrow 2_1^+$		0.2161
$1_1^+ \rightarrow 2_1^+$		0.1972	$1_1^+ \rightarrow 2_1^+$		0.2856

Journal of Kerbala University, Vol. 10 No.1 Scientific . 2012

Table 4 . The mixing ratios values of $~^{102}\mathrm{Mo}$ and $~^{102}\mathrm{Zr}$

$I_i \!$	¹⁰² Mo		$I_i \!$	¹⁰² Zr	
	Exp.	IBM-2		Exp.	IBM-2
$2_1^+ \rightarrow 2_2^+$		1.0777	$2_1^+ \rightarrow 2_2^+$		2.3913
$2_1^+ \rightarrow 2_3^+$		0.0781	$2_1^+ \rightarrow 2_3^+$		0.8948
$4_1^+ \rightarrow 4_2^+$		0.1074	$4_1^+ \rightarrow 4_2^+$		1.4056
$3_1^+ \rightarrow 2_1^+$		0.1943	$3_1^+ \rightarrow 2_1^+$		0.4474
$1_1^+ \rightarrow 2_1^+$		0.1753	$1_1^+ \rightarrow 2_1^+$		0.0059

Table 5 . The mixing ratios values of 102 Pd and 102 Ru

$I_i {\rightarrow} I_f$	¹⁰² Pd		$I_i {\rightarrow} I_f$	¹⁰² Ru	
	Exp.[11]	IBM-2		Exp. [12]	IBM-2
$2_1^+ \rightarrow 2_2^+$	10.4	6.0107	$2_1^+ \rightarrow 2_2^+$		10.5100
$2_1^+ \rightarrow 2_3^+$	8.1	0.4380	$2_1^+ \rightarrow 2_3^+$	0.25	0.2085
$4_1^+ \rightarrow 4_2^+$		4.7389	$4_1^+ \rightarrow 4_2^+$		3.3664
$3_1^+ \rightarrow 2_1^+$		0.9183	$3_1^+ \rightarrow 2_1^+$	- 8.4	0.3467
$1_1^+ \rightarrow 2_1^+$		1.9234	$1_1^+ \rightarrow 2_1^+$		0.3504

4- Conclusion :

In this work , Interacting Boson Models IBM-1and IBM-2, have been used to calculate energy levels and nuclear properties of some even-even nuclei with (A=100) and (A=102). We observed that IBM-2 is the best in study of this nuclei, also we conclude that (100 Mo , 102 Pd , 102 Ru) in U(5) limits and (100 Sr , 102 Zr) in SU(3) limit, (102 Mo) in O(6) limits. energy levels of the low lying states of these nuclei were produced, Mixing ratios $\delta(E2/M1)$ for transitions with $\Delta I = 0$, $I \neq 0$ were calculated. the results are compared with available experimental data, Satisfactory agreements were produced.

5- References :

- [1] Bangay J.C. ; Structure of 110 Cd Studies with β decay of 110 In ; MSc. Thesis, University of Guelph (2010) .
- [2] Cook N.D. ;"Models of The Atomic Nucleus" ; Philadelphia (2005) .
- [3] Yazar H.R. & Uluer I. ; Math. And Comput. Appli. ; Vol.12,No.2 (2007) .
- [4] Zamfir B., Warner D.D.; Phy.Rev. Vo.(60) (1999).
- [5] Casten R.F. & Warner D.D.; Rev. Mod. Phys. P389 (1988).
- [6] E.P.Rodal ; Rev. Mex. S.52 (2006).
- [7] Jin-Fu Z. & Al-Khudair F.; Commun. Theor. Phys.; (Beijing, China) Vol.37, No.3 (2002).

Journal of Kerbala University, Vol. 10 No.1 Scientific . 2012

- [8] Turkan N. & Maras I. ; Indian Acadm. Of Science ; Vol.68 , No.5 (2007) .
- [9] Singh B.; Nuclear Data Sheets; Vol. 109 No. 2 (2008).
- [10] Martin M.J.; Nuclear Data Sheets; Vol. 35 No.4 (1998).
- [11] Inci I. & Turkan N.; Turk J. Phys.30 (2006).

[12] Yilmaz A.H. & Kuruoglu M.; Commun. Theor. Phys.; (Beijing, China) Vol.46, No.4 (2006)

The energy levels which calculated in the present study are in the following figures : Fig.(1): Comparison between experimental,IBM-1 and IBM-2 energy levels of ¹⁰⁰Sr

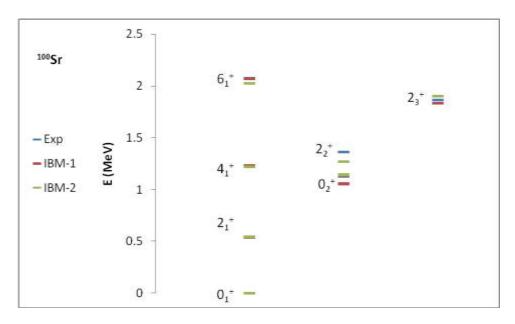


Fig.(2): Comparison between experimental, IBM-1 and IBM-2 energy levels of ¹⁰⁰Mo

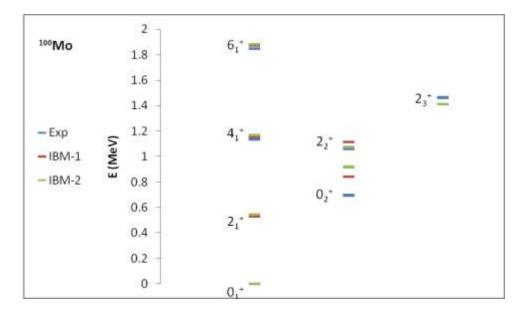


Fig.(3): Comparison between expermental,IBM-1 and IBM-2 energy levels of ¹⁰²Zr

Journal of Kerbala University, Vol. 10 No.1 Scientific. 2012

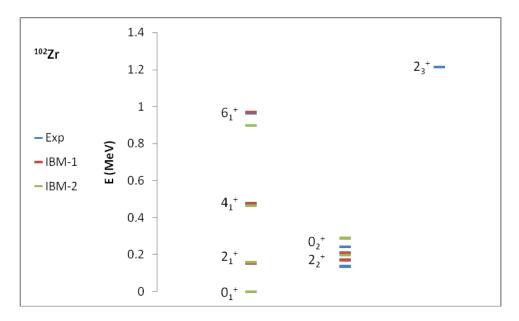
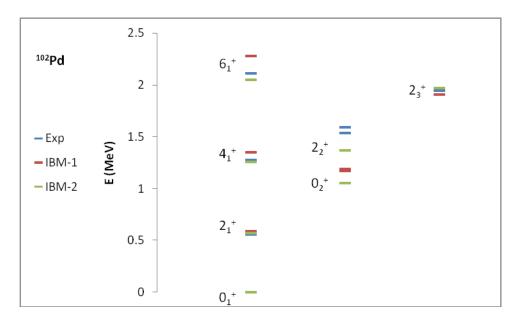
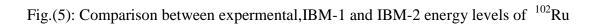


Fig.(4): Comparison between experimental,IBM-1 and IBM-2 energy levels of ¹⁰²Pd





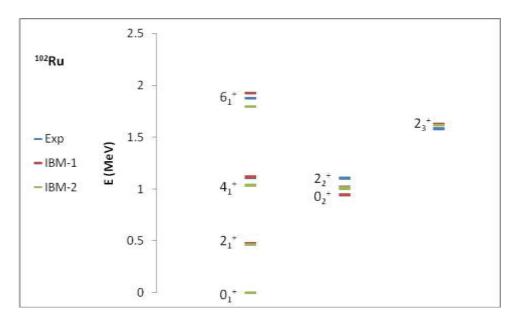


Fig.(6): Comparison between expermental,IBM-1 and IBM-2 energy levels of ¹⁰²Mo

