Structure of odd-odd Ga and As nuclei, and dynamical and supersymmetries

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The structure of 66 Ga, 68 Ga, 70 Ga, 70 As, 72 As, 73 As, 74 As, and 76 As is studied via $(p, n\gamma)$ [and in some cases $(\alpha, n\gamma)$] reactions. In-beam γ -ray, two-dimensional $\gamma\gamma$ -coincidence, internalconversion electron, and γ -ray angular-distribution spectra, as well as $\sigma(E_{\text{LEV}}, E_p)$ relative reaction cross sections are measured with Ge(HP), Ge(Li) γ and combined superconducting-magnetic plus Si(Li) electron spectrometers at different bombarding-particle energies. The proposed new level schemes contain 300 (among them 70 new) levels. Level spins, parities, and γ -ray branching and mixing ratios are deduced. Energy spectra, electromagnetic moments, reduced transition probabilities, y-ray branching ratios, and one-nucleon transfer reaction spectroscopic factors are calculated in the framework of the interacting boson model (IBM), the interacting boson-fermion model (IBFM), and the interacting boson-fermion-fermion model (IBFFM) for about 20 nuclei in the Ga-As region (⁶⁴⁻⁶⁷Zn, ⁶⁵⁻⁶⁸Ga, ⁶⁸⁻⁷³Ge, ⁷⁰⁻⁷⁴As). The odd-odd nuclei are described on the basis of a consistent parametrization deduced from the even-even core and the two neighboring odd-A nuclei. A reasonable description of the experimental data is obtained. The energy splitting of proton-neutron multiplets in odd-odd Ga and As nuclei is discussed by the use of the parabolic rule associated with the IBFFM in low perturbation order. The energy spectra of ${}_{34}^{74}\text{Se}_{40}, {}_{33}^{75}\text{Se}_{40}, {}_{33}^{74}\text{As}_{40}, {}_{33}^{74}\text{As}_{41}$ supermultiplet nuclei are calculated on the basis of the $U_{\pi}(6/12) \otimes U_{\nu}(6/12)$ supersymmetry (SUSY) theory, using a simple closed energy formula; 44 states of these four different nuclei are reasonably described with only seven fitted parameters. The existence of supersymmetry is supported also by one-nucleon transfer reaction data, electromagnetic properties, and IBFM, IBFFM and SUSY model wave functions of the levels considered. © 1995 American Institute of Physics.

1. INTRODUCTION

The main intention of the present work was to make a detailed in-beam γ - and electron-spectroscopic study of the odd-odd Ga and As isotopes and to give a consistent description of the structure of the Zn, Ga, Ge, and As nuclei in the framework of the interacting boson (boson-fermionfermion) model [IB(FF)M]. The nuclei investigated in this program are shown in Fig. 1.

The experimental work was motivated by the fact that the level systems of the investigated odd-odd nuclei (and especially the level spin and parity values) were known very scantily. For example, in ⁶⁸Ga unambiguous spin-parity data had been determined only for three levels before our measurements. The beams of the Debrecen isochronous cyclotron enabled excitation of both particle and collective states, and the high-resolution, high-efficiency Ge(HP) detectors and unique superconducting magnetic electron spectrometer, constructed at the Institute of Nuclear Research, assured a good possibility for complex γ - and electron-spectroscopic in-beam studies.

A theoretical interpretation of the ⁶⁶Ga and ⁶⁸Ga level schemes was formerly completely missing. In the case of ^{70,72,74}As we have performed calculations in the interacting boson-fermion-fermion model for the first time. These calculations give a consistent and detailed description of the energy spectra and electromagnetic properties of the corresponding even-even, odd-A and odd-odd Zn, Ga, Ge, and As nuclei.

The new experimental data obtained on ⁷³As and ⁷⁴As offered a possibility of checking the supersymmetry predictions for ⁷⁴Se, ⁷⁵Se, ⁷³As, and ⁷⁴As nuclei.

2. EXPERIMENTAL METHODS AND RESULTS

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We have studied the excited states of the Ga and As nuclei in the proton and α -particle beams of the Debrecen

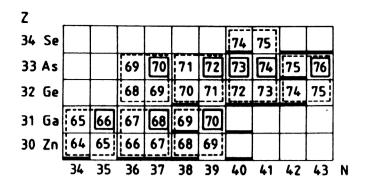


FIG. 1. Nuclei investigated in this program experimentally are indicated by eight mass numbers in solid frames. Mass numbers in dotted frames: theoretical calculations. Z: atomic; N: neutron numbers; mass number underlined with a thick line: stable nucleus.

103-cm (and in some experiments of the Jyväskylä 90-cm) isochronous cyclotrons via $(p, n\gamma)$ and $(\alpha, n\gamma)$ reactions. The targets were prepared from enriched Cu, Zn, and Ge isotopes. The description of the spectroscopic channels of the Debrecen cyclotron can be found in Refs. 1 and 2.

High-resolution Ge (HP and LEPS) detectors were used for a γ and superconducting magnetic lens plus Si(Li) spectrometer (SMLS) for electron-spectroscopic measurements. The SML spectrometer, which was described in detail in Ref. 3, has high transmission, good energy resolution, and low background.

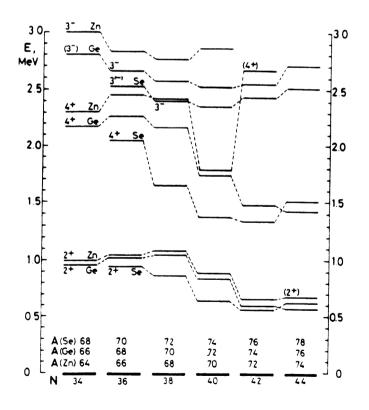


FIG. 2. Energies of 2_1^+ , 4_1^+ , and 3_1^- states of the even–even Zn, Ge, and Se nuclei. Data were taken from the corresponding Nuclear Data Sheets.

In order to obtain "complete" spectroscopic information, γ -ray (E_{γ}, I_{γ}) , $\gamma\gamma$ -coincidence, internal-conversion electron, and γ -ray angular-distribution spectra, as well as the relative reaction cross section $\sigma_{rel}(E_{LEV}, E_p)$ were measured at different bombarding-particle energies. Level schemes, spin-parity values, and y-branching and y-mixing ratios have been deduced. Great attention was paid to the reliability and consistency of the experimental data. For example, the level spins have been determined with three different methods: a) from the decay properties and internalconversion coefficients of transitions; b) from a Hauser-Feshbach analysis of the (p,n) reaction cross sections; c) from γ -ray angular distributions. The configuration of levels has been determined from all available data: from nucleontransfer reaction studies, $\log ft$ values of β decay, electromagnetic moments, transition probabilities, γ-branching ratios, predictions of parabolic-rule calculations, etc.

As a result of experimental work about 880 (among them 440 new) γ rays have been identified, and 280 (including 240 new) α_K internal-conversion coefficients and \sim 300 (among them 70 new) levels have been determined in 66,68,70 Ga and 70,72,73,74,76 As nuclei.

The results have been published in the following papers: ⁶⁶Ga (Refs. 4 and 13), ⁶⁸Ga (Refs. 5, 6, and 13), ⁷⁰Ga (Ref.

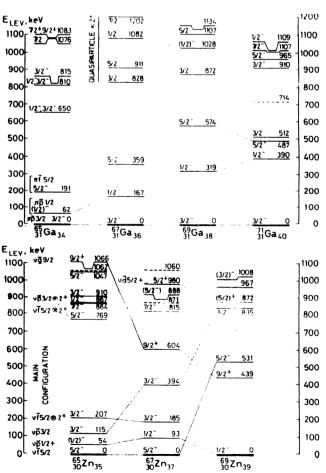


FIG. 3. The low-lying levels of odd-A ^{65–69}Zn and ^{65–71}Ga nuclei. Data were taken from the corresponding Nuclear Data Sheets and original papers.

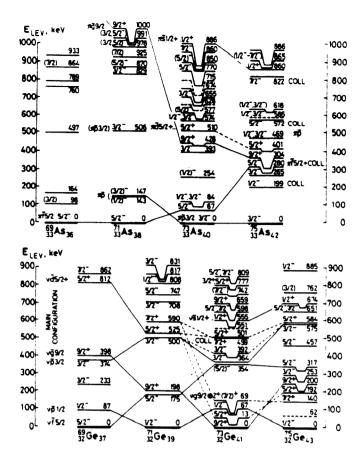


FIG. 4. The low-lying levels of odd- A^{69-75} Ge and $^{69-75}$ As nuclei. Data were taken from the corresponding Nuclear Data Sheets and original papers.

7), 70 As (Ref. 8), 72 As (Refs. 9 and 13), 73 As (Ref. 10), 74 As (Refs. 11–13), and 76 As (Ref. 14). These papers contain detailed descriptions of the experimental methods and results, deduced new level schemes, and their discussion. The level spectra were based mainly on $\gamma\gamma$ -coincidence relations, as well as on the energy and intensity balance of the transitions.

3. SYSTEMATICS OF THE EVEN-EVEN AND SINGLE-ODD NUCLEI IN THE Zn-Se REGION

If we want to describe the structure of odd-odd nuclei, we need to know the energy levels of the neighboring eveneven and odd-A nuclei. The systematics of the 0_1^+ , 2_1^+ , 4_1^+ , and 3_1^- states of the even-even Zn, Ge, and Se nuclei is shown in Fig. 2. The low-lying levels of the odd-A Zn, Ga and Ge, As nuclei are presented in Figs. 3 and 4, respectively. The main configurations of states are also indicated in Figs. 3 and 4 on the basis of the one-nucleon transfer reaction and other available data.

The systematics of neutron quasiparticle energies and occupation probabilities in the Ni–Se region are shown in Fig. 5.

In order to get preliminary estimates for the configurations of the low-lying levels of odd-odd Ga and As nuclei, we have performed parabolic-rule¹⁶ calculations. These calculations proved very useful for the prediction of the energy splitting of different proton-neutron multiplets in odd-odd

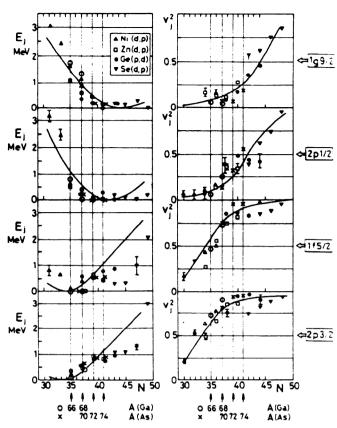


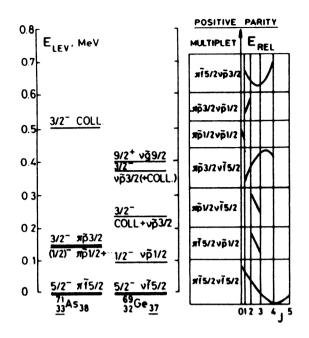
FIG. 5. Experimental quasiparticle energies (E_j) and occupation probabilities as a function of the neutron number (N) (points) and the corresponding BCS theoretical results (curves). Data were taken from Fournier *et al.*¹⁵ Values used in the present calculations are shown by circles (\bigcirc) for Ga and crosses (\times) for As odd-odd nuclei.

In (Ref. 1) and Sb (Ref. 2) nuclei. In the Ga and As isotopes rather strong configuration mixing is expected among the close-lying identical spin-parity states, which may imply a limitation on the applicability of the parabolic rule. Nevertheless, it can be used for the first orientation, as indeed the IBFFM calculations showed later.

As the parabolic rule and its applications to In and Sb nuclei were described in detail in Refs. 1, 2, and 16, here we show only the results obtained on the relative energy splitting of proton-neutron multiplets in ⁷⁰As (Fig. 6) and ^{66,68,70}Ga, ^{70,72,74,76}As nuclei (Fig. 7). For these predictions the same proton and neutron occupation probabilities were used, as later in the IBFFM calculations (see Sec. 4).

4. INTERACTING BOSON AND BOSON-FERMION-FERMION MODEL CALCULATIONS

In order to get deeper and more detailed insight into the structure of the low-lying odd-odd Ga and As states, we have calculated the level energies and electromagnetic properties on the basis of the interacting boson (boson-fermion-fermion) model.



ENERGY SPLITTING OF DIFFERENT p-n MULTIPLETS IN 33 As 37 ACCORDING TO THE PARABOLIC RULE

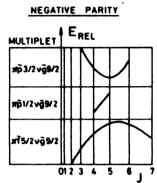


FIG. 6. Approximate relative energy splitting of different proton-neutron multiplets in ⁷⁰As, as predicted by the parabolic rule. The abscissa is scaled according to J(J+1), where J is the spin of the

4.1. Hamilton operator

The Hamiltonian of the model is 17-20

$$H_{\rm IBFFM} = H_{\rm IBFM}(\pi) + H_{\rm IBFM}(\nu) - H_{\rm IBM} + H_{\rm RES}(\pi\nu), \qquad (1)$$

where $H_{\rm IBM}$ denotes the IBM Hamiltonian for the even–even core nucleus, 21,22 $H_{\rm IBFM}(\pi)$ and $H_{\rm IBFM}(\nu)$ are the IBFM Hamiltonians for the neighboring odd-even nuclei with an odd proton and odd neutron, respectively, $^{23-25}$ and $H_{RES}(\pi\nu)$ is the Hamiltonian of the residual interaction.

The Hamiltonian of the core has the form²⁶⁻²⁸

$$H_{\text{IBM}} = h_1 \hat{N} + h_2 \{ (d^+ d^+)_0 [(N - \hat{N})(N - \hat{N} - 1)]^{1/2} + \text{H.c.} \}$$

$$+ h_3 [(d^+ d^+ \tilde{d})_0 (N - \hat{N})^{1/2} + \text{H.c.}]$$

$$+ \sum_{L=0,2,4} h_{4L} [(d^+ d^+)_L (\tilde{d}\tilde{d})_L]_0, \qquad (2)$$

where \hat{N} is the d-boson number operator and N is the total number of s and d bosons. The relation of the $\{h\}$ parameters to the parameters defined in Ref. 21 is as follows:

MULTIPLET	PA - RITY		56Ga ₃₅	68Ga ₃₇	70Ga 39	⁷⁰ AS ₃₇	77 AS 39	74AS ₄₁	76AS ₄₃
#₽̃3/2∨g̃9/2	-		U	U	U	Ü	Ü	Ü	Ü
ภ¤ี3/2∨¤1/2	+		/	/	/	/	/	/	/
≢p̃ 3/2 vī̃ 5/2	+		I	P	P	P	P	P	P
я́р 3/2 v́р 3/2	+		J	0	0	0	3	1	J
я₹5/2 ∨5 9/2	-		0	P	0	0	O	0	Ü
Я₹5/2∨₿1/2	+	Ξ	\	\	\	\	\	\	\
ภ์สี 5/2 v ี 5/2	+	(RELATIVE	J	v	U	U	v	V	U
я75/2vβ3/2	+	8	IJ	U	U	U	U	U	U
яў 1/2 v ў 9/2	-	اخ	/	/	/	/	/	/	1
яр 1/2 vр 1/2	+	ENERGY	\	\	\	\	\	\	\
яр 1/2 v ў 5/2	+	EN	\		\	\	\		\
я́р 1/2 v р 3/2	+		/	/	/	/	/	/	/
яğ9/2 v ğ 9/2	+		3	Ü	J	Ü	Ü	Ü	J
лф9/2 үр 1/2	-		/	/	/	/	/	/	/
яў9/2 v7 5/2	-		P	U	U	U	U	U	U
#g9/2∨p3/2	-		<u>U</u>	U	U	Ü	Ù	U	U
		ı						1(1+	.1)

FIG. 7. Approximate relative energy splitting of different proton-neutron multiplets in some odd-odd Ga and As nuclei as a function of J(J+1), where J is the spin of the state. The predictions are based on the parabolic rule.

TABLE I. Parameters of the IBFFM calculations.

Paramet	ers	6/Ga ₃₅	MGa ₃₇	70As ₃ ,	72A1,9	74As ₄₁
Core, MeV	۸,	0.8	1 039	09	0.88	0.68
	h,	-02	0	-0.15	-0.3	-0.25
	۸,	0	0	0.06	0 12	0.1
	h _{en}	01	0.147	0	0.2	0
	h ₄₂	-0.15	-0.2292	-0.5	-0.4	-0.3
_	h44	0.3	0.5595	-0.08	-0.08	-0.08
Total boson N ⁰ ,	N	4	5	3	4 (or 7)	4
	z	-0.5	-0.5	-1.323*	-1.323*	-1.323*
e	€VIB	1 35	1.35	0.8	0.8	0.8
Quasi-proton	E(14 3/2)	0	0	0.3	0	0
energy, MeV	E(# 5/2)	0.75	0.75	0	0.41	0.41
	E(xp 1/2)	0.615	0.615	0.3	0.74	0.74
	E(xg 9/2)			1.3	2.2	1.55
	E(#4 5/2)				5.2	4.55
Оссир.	V 2(Rp 3/2)	0.6	0.6	0.607	0.607	0.579
probability	V 2(#/ 5/2)	0.07	0 07	0.309	0.309	0.341
	V 2(Rp 1/2)	0 09	0.09	0.131	0.131	0.118
	V2(Rg 9/2)			0.07	0 07	0.06
	$V^2(nd.5/2)$				0 01	0.01
Quasi-neutron	E(vp 3/2)	0.09	0 49	0.60	0.95	0 95
energy, MeV	E(vf 5/2)	0	0	0	0 53	0 53
	E(vp 1/2)	0.83	0 03	0.2	0	0
	E(vg 9/2)	1 69	1.36	0.9	02	0 2
	E(vd 5/2)				3 2	3.2
Occup	V ² (vp 3/2)	0.73	0.89	0.80	0.941	0.957
probability	V 2(vf 5/2)	0.48	0.73	0.75	0.949	0.962
	V2(vp 1/2)	0.11	0.25	0.16	0.327	0.624
	V2(vg 9/2)	0.08	0.04	0.09	0.09	0.217
	V 2(vd 5/2)				0 01	0.01
Strengths of	A ₀ *	0 05	0.05	0 05	0.05	0.05
boson-fermion	Γ <u>ξ</u>	0.48	0.4	0.4	0.5	0.5
interaction, MeV		14	0.5	0.5	0 5x +	0.5
	۸'n	'	"	0.5		0.5
	A ₀ ^v	0	0	1 0	1.5x -	0
		0.02	0.15			
	Ľ	""	013	0.2	0 55R +	0.55
	AV	1 63			0 45x -	
	۸,	105	1.3	13	1 30x +	1.3π+
					5.50π -	5.5x ~
Effective gyromagn. ratios	8,*	0.4g, E./ 0	0.4g _z ^{g_f}	$0.40g_{z}^{R,f}\pi +$	0.7g, E.f	0.7g,x.
ay ronningst. Lattics				0.65g, R. f x -		
	E team	1.56	1.59 0	0	3.340	3.372
•	s,*	0.9g, v/ V	0.5g," x+	0.48,7	0.78,	0.78,
	-1	"'	0.9g, v/ x -			J. 78,
	٠٠ ••	-1.07	-1.09 A	o	-4.40	4.40
	8 tens					-4.48
Strengths of residual	v_{8}	-04	-0.4	0 R +	0π+	0 x +
interaction, MeV				-04π−	-0.6 π –	-06π-
	V _{ee}	0.4	0.4	0.4 π +	0	0
	V tens			03x-		
		-001	0 015	0.015	0 0 1 5	

 e^{π} = 1.5 e, e^{ν} = 0.5 e, g_l^{π} = 1, g_l^{ν} = 0, g_R = Z/A in all cases. *1.323 = $\sqrt{7/2}$.

^{*1.225=} $\sqrt{t/2}$. π^+ and π^- stand for positive- and negative-parity states, respectively.

**The tensor gyromagnetic ratio, $g_{\text{tens}}^{\pi,\nu}$ (Refs. 59 and 60) is similar to the one used in Ref. 61. $\diamond \frac{1}{50}(r^2)g_x^{\pi}$ (frec) = 1.59; $\diamond \frac{1}{50}(r^2)g_x^{\nu}$ (free) = -1.09. $\Box g_x^{\pi,f} = g_x^{\pi}$ (free) = 5.5857. $\nabla g_x^{\nu,f} = g_x^{\nu}$ (free) = -3.8263.

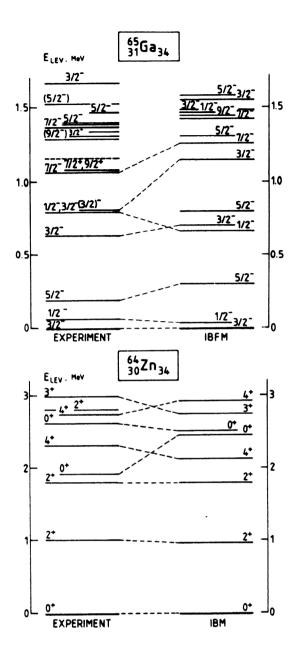


FIG. 8. Experimental energy spectra of 64 Zn (Ref. 30) and 65 Ga (Ref. 31) and the corresponding theoretical IBM and IBFM results. 4

$$h_1 = \epsilon_d - \epsilon_s + \left(\frac{1}{\sqrt{5}}u_2 - u_0\right)(N-1),$$

$$h_2 = \frac{1}{\sqrt{2}} \tilde{v}_0, \quad h_3 = \tilde{v}_2,$$

$$h_{4L} = \sqrt{2L+1} \left(\frac{1}{2} c_L - \frac{1}{\sqrt{5}} u_2 + \frac{1}{2} u_0 \right). \tag{3}$$

The IBFM Hamiltonian employed here is of the form²⁷

$$H_{\rm IBFM}(\alpha) = H_{\rm IBM} + \sum_{i} \tilde{\epsilon}_{i}(\alpha) + H_{\rm BF}(\alpha),$$
 (4)

where α stands for an odd proton $(\alpha = \pi)$ or neutron $(\alpha = \nu)$. The second and third terms are the quasiparticle and boson–fermion (particle-vibration) interaction Hamiltonians, respectively, and

$$\begin{split} H_{BF}(\alpha) &= \sum_{j} A_{j} \{ (d^{+}\tilde{d})_{0} [c_{j}^{+}(\alpha)\tilde{c}_{j}(\alpha)]_{0} \}_{0} \\ &+ \sum_{j_{1}j_{2}} \Gamma_{j_{1}j_{2}} \{ Q_{2} [c_{j_{1}}^{+}(\alpha)\tilde{c}_{j_{2}}(\alpha)]_{2} \}_{0} \\ &+ \sum_{j_{1}j_{2}j_{3}} \Lambda_{j_{1}j_{2}j_{3}} : \{ [c_{j_{1}}^{+}(\alpha)\tilde{d}]_{j_{3}} [\tilde{c}_{j_{2}}(\alpha)d^{+}]_{j_{3}} \}_{0} :, \quad (5) \end{split}$$

with

$$A_{i} = A_{0}\sqrt{5}(2j+1),$$
 (6)

$$\Gamma_{j_1 j_2} = \Gamma_0 \sqrt{5} (u_{j_1} u_{j_2} - v_{j_1} v_{j_2}) \langle j_1 || Y_2 || j_2 \rangle, \tag{7}$$

$$\Lambda_{j_1 j_2 j_3} = -2\Lambda_0 \frac{\sqrt{5}}{\sqrt{2j_3 + 1}} (u_{j_1} v_{j_3} + v_{j_1} u_{j_3})$$

$$\times (u_{j_2}v_{j_2} + v_{j_2}u_{j_2}) \cdot \langle j_3 || Y_2 || j_1 \rangle \langle j_3 || Y_2 || j_2 \rangle,$$
 (8)

$$Q_{2\mu} = d_{\mu}^{+} \sqrt{N - \hat{N}} + \sqrt{N - \hat{N}} \tilde{d}_{\mu} + \chi (d^{+} \tilde{d})_{2\mu}. \tag{9}$$

The Hamiltonian of the residual interaction was taken in the form

$$H_{\text{RES}}(\pi\nu) = 4\pi V_{\delta} \delta(\mathbf{r}_{\pi} - \mathbf{r}_{\nu}) \, \delta(r_{\pi} - R_{0}) - \sqrt{3} V_{\sigma\sigma}(\boldsymbol{\sigma}_{\pi} \cdot \boldsymbol{\sigma}_{\nu})$$

$$+ V_{\text{tens}} \left[\frac{3(\boldsymbol{\sigma}_{\pi} \cdot \mathbf{r}_{\pi\nu})(\boldsymbol{\sigma}_{\nu} \cdot \mathbf{r}_{\pi\nu})}{r_{\pi\nu}^{2}} - (\boldsymbol{\sigma}_{\pi} \cdot \boldsymbol{\sigma}_{\nu}) \right], \quad (10)$$

where $\mathbf{r}_{\pi\nu} = \mathbf{r}_{\pi} - \mathbf{r}_{\nu}$, $R_0 = 1.2\sqrt[3]{A}$ fm; $H_{\text{RES}}(\pi\nu)$ includes surface delta, spin-spin, and tensor interactions.

The Hamiltonian (1) was diagonalized in the proton-neutron-boson basis: $|(j_{\pi},j_{\nu})j_{\pi\nu},n_dI;J\rangle$, where j_{π} and j_{ν} stand for the proton and neutron angular momenta coupled to $j_{\pi\nu}$, n_d is the number of d bosons, I is their angular momentum, and J is the spin of the state. The computer codes used in the calculations were written by Brant, Paar and Vretenar.²⁹

4.2. Method of calculation. Parametrization

In the first stage of the calculations we fitted the $\{h_i\}$ parameters of the $H_{\rm IBM}$ Hamiltonian (2) to the energy spectrum of the corresponding core nuclei (⁶⁴Zn, ⁶⁶Zn, ⁶⁸Ge, ⁷⁰Ge, and ⁷²Ge).

The total boson number is given by the number of valence shell pairs. For example, in $^{66}_{30}$ Zn₃₆ (the core of $^{68}_{31}$ Ga₃₇) we have N=5, since in this nucleus there are one proton and four neutron bosons, relative to the Z=N=28 doubly magic nucleus. In the case of $^{70}_{32}$ Ge₃₈ (the core of $^{72}_{33}$ As₃₉) and $^{71}_{32}$ Ge₃₉ the calculations were performed with both N=7 and N=4. The two calculations, with renormalization of the other parameters, gave similar results for the energy spectra and electromagnetic properties. Thus, in the further calculations for 70 As, 72 As, 74 As, and nine neighboring nuclei we have used reduced total boson numbers, by renormalizing the IBM pa-

TABLE II. Main components (\geq 4%) in the IBFFM wave functions of some low-lying states in ⁶⁶Ga. The basis states are $|(j_{\pi}j_{\nu})j_{\pi\nu},n_{d}I;J\rangle$ (see the text). The last column displays the corresponding amplitudes in the wave functions.

J ^z	$(j_{\mathbf{x}}, j_{\mathbf{v}})$	j _{πv} ; n _d I	A	J *	$(j_{\mathbf{x}},j_{\mathbf{v}})$	$j_{\pi v}$; $n_d I$	A
0+	(3/2,5/2)	2:12	0.62	2+	(3/2,5/2)	2;00	0.35
	(3/2,3/2)	0;00	0.23		(3/2,3/2)	2;00	0.45
	(3/2,3/2)	2;12	0.43		(3/2,5/2)	3;12	-0.24
	(1/2,5/2)	2;12	-0.26		(3/2,3/2)	3;12	-0.41
	(1/2,3/2)	2;12	0.30	02	(3/2,3/2)	0;00	0.36
	(3/2,5/2)	2;32	-0.22		(1/2,1/2)	0:00	-0.25
11	(3/2,5/2)	1;00	0.54		(5/2,5/2)	0;00	-0.53
	(3/2,5/2)	1;20	-0.24		(5/2,5/2)	0;20	0.28
	(3/2,3/2)	2;12	0.31		(3/2,5/2)	2;12	-0.43
	(1/2,5/2)	2;12	0.24		(1/2,3/2)	2;12	0.21
	(3/2,5/2)	3;12	0.41	4;	(1/2,5/2)	3;12	-0.21
	(3/2,3/2)	3;12	-0.23		(3/2,5/2)	4;00	0.62
2†	(3/2,5/2)	2;00	0.62		(5/2,5/2)	4:00	0.37
	(1/2,5/2)	2;00	-0.25		(3/2,5/2)	4;20	-0.26
	(3/2,5/2)	2;12	-0.23		(5/2,5/2)	5;12	0.39
	(3/2,5/2)	2:20	-0.27	5-	(1/2,9/2)	5;00	0.63
	(1/2,5/2)	3;12	-0.27		(5/2,9/2)	5;00	-0.38
	(5/2,5/2)	4;12	-0.26		(1/2,9/2)	5;20	-0.30
1 2	(3/2,5/2)	1;00	0.51		(5/2,9/2)	7;12	0.35
	(3/2,5/2)	1;20	-0.21	7-	(1/2,9/2)	5;12	0.32
	(3/2,5/2)	2;12	0.30		(5/2,9/2)	7;00	0.74
	(3/2,3/2)	2;12	-0.24		(5/2,9/2)	7;12	-0.26
	(1/2,3/2)	2;12	-0.26		(5/2,9/2)	7;20	-0.39
	(3/2,3/2)	3;12	0.33				
31	(3/2,5/2)	3;00	-0.46				
	(3/2,3/2)	3;00	0.20				
	(1/2,5/2)	3;00	0.50				
	(1/2,5/2)	3;20	-0 23				
	(5/2,5/2)	5;12	0.43				

rameters. This strongly reduced the volume of computations for odd-odd nuclei, without a substantial effect on the properties of the low-lying states.

The χ and vibrational charge (e^{VIB}) parameters were fitted to the electromagnetic moments and reduced B(E2) transition probabilities of the core nuclei.

The nucleus 66 Zn exhibits a level scheme which is characteristic of a spherical vibrator. In this case the parameters h_2 and h_3 were chosen to be equal to zero. In other nuclei we employed a parametrization which corresponds to a transition between the U(5) and O(6) dynamical symmetries, but somewhat closer to the U(5) character.

The core parameters employed in the calculations are given in the first part of Table I.

In the second stage of the calculations we adjusted the parameters of $H_{\rm IBFM}(\pi)$ [Eqs. (4)–(8)] to the experimental data of the corresponding odd-Z, even-N nuclei. The proton quasiparticle energies and occupation probabilities were

taken mostly from proton-transfer reaction data and/or pairing-force (BCS) calculations.

The A_0^{π} monopole, Γ_0^{π} dynamical quadrupole, and Λ_0^{π} exchange boson-fermion interaction strengths were fitted to the low-energy spectra of ⁶⁵Ga, ⁶⁷Ga, ⁷¹As, and ⁷³As, respectively. (The level spectrum of ⁶⁹As was scarcely known.) The effective gyromagnetic ratios g_s^{π} and g_{tens}^{π} were determined by fitting to the electromagnetic moments and reduced B(E2), B(M1) transition probabilities (γ -branching ratios).

In the third stage of the calculations we adjusted the parameters A_0^{ν} , Γ_0^{ν} , and Λ_0^{ν} of $H_{\rm IBFM}(\nu)$ [Eqs. (4)–(8)] to the low-energy spectra of the corresponding even-Z, odd-N nuclei: ⁶⁵Zn, ⁶⁷Zn, ⁶⁹Ge, ⁷¹Ge, and ⁷³Ge. The proton quasiparticle energies and occupation probabilities were close to the calculated BCS values of Ref. 15 and to the systematics of the experimental data (Fig. 5). The parameters g_x^{ν} and $g_{\rm tens}^{\nu}$ were fitted to the electromagnetic properties.

Finally, the parameters V_{δ} , V_{qq} , and V_{tens} of the residual

TABLE III. Wave functions of some low-lying states of 70 As. Only the strongest components are given. The basis states are $|(j_{\pi}j_{\nu})j_{\pi\nu}, n_d I; J\rangle$ (see the text).

J*	$(j_{\mathbf{x}},j_{\mathbf{v}})$	$j_{\pi v}; n_d^{-1}$	M	Jπ	(j_{π}, j_{ν})	$j_{\pi v}; n_d I$	И
0+	(5/2,5/2)	0;00	0.83	34	(3/2,5/2)	3.00	0.79
0_2^+	(1/2,1/2)	00;0	0.70	4†	(5/2,5/2)	4:00	0.77
1+	(3/2,5/2)	1;00	0.73	42	(3/2,5/2)	4;00	0.56
1+2	(5/2,5/2)	1;00	0.66		(5/2,5/2)	5;12	0.48
1 3	(1/2,1/2)	1;00	0.66	2-	(5/2,9/2)	2;00	0.76
21	(5/2,5/2)	2;00	0.51	3-	(5/2,9/2)	3;00	0.75
	(1/2,5/2)	2;00	0.45	4-	(1/2,9/2)	4;00	0.52
2+	(5/2,5/2)	2;00	0.56		(5/2,9/2)	4:00	0.41
	(3/2,5/2)	2;00	0.41	4-2	(5/2,9/2)	4;00	0.65
2*	(5/2,1/2)	2;00	0.63	5-	(5/2,9/2)	5;00	0.56
31	(1/2,5/2)	3;00	0.51		(1/2,9/2)	5;00	0.49
	(5/2,5/2)	3;00	0.48	6-	(5/2,9/2)	6;00	0.70
3 ⁺ ₂	(5/2,5/2)	3;00	0.63	7-	(5/2,9/2)	7:00	0.75
3*	(5/2,1/2)	3;00	0.78				

interaction (10) were fitted to the experimental data on odd-odd nuclei. The core parameters and occupation probabilities remained unaltered in all cases. We remark that in 70 As, 72 As, and 74 As readjustment of the quasiparticle energies, boson-fermion interaction strengths, and effective gyromagnetic ratios was needed in order to obtain better agreement with the experimental data. Such a renormalization seems to be in accord with a general observation in the region of soft nuclei: the dynamical deformation can be sizable if one nucleon is added. Consequently, the parameters may change noticeably from those derived from the neighboring nuclides. A similar feature was found, for example, in the A = 130 region.

The parameters applied for the IBFFM description of the properties of the odd-odd nuclei are summarized in Table I.

4.3. Results. Discussion

The calculated energy spectra are compared with the experimental data in Figs. 8 (⁶⁴Zn, ⁶⁵Ga), 9 (⁶⁵Zn, ⁶⁶Ga), 10 (⁶⁶Zn, ⁶⁷Ga), 11 (⁶⁷Zn, ⁶⁸Ga), 12 (⁶⁸Ge, ⁶⁹Ge, ⁷⁰As), 13 (⁷⁰Ge, ⁷¹As), 14 (⁷¹Ge, ⁷²As), 15 (⁷²Ge, ⁷³As), and 16 (⁷³Ge, ⁷⁴As). Although the level schemes are usually very complicated (especially in odd–odd nuclei), reasonable agreement has been obtained between experiment and theory. The calculated states are usually assigned to the experimental levels on the basis of the energy, spin, parity, one-nucleon transfer reaction data (if they exist), decay properties, and wave functions.

The main components of the wave functions of some low-lying states in ⁶⁶Ga and ⁷⁰As are shown in Tables II and III, respectively. Usually, the wave functions are very complex. There are states which have more than 600 components. Nevertheless, in some cases the states are dominated

by only one proton—neutron multiplet. Thus, it is worth comparing the experimental and IBFFM energy spectra also with the predictions of the parabolic rule (Figs. 6 and 7).

The $\pi \tilde{p}_{3/2} \nu \tilde{f}_{5/2}$ multiplet. As an example, let us discuss the energy splitting of the $\pi \tilde{p}_{3/2} \nu \tilde{f}_{5/2}$ multiplet in ^{66,68,70}Ga and ^{70,72}As. The parabolic rule predicts an open-up parabola for the energy splitting in ⁶⁶Ga and an open-down one in ⁶⁸Ga (Fig. 7). This is in accord with the experimental facts, because the 1^+ , 2^+ , 3^+ , and 4^+ states of the multiplet can be identified with the 44- and 109-keV $1_1^+ + 1_2^+$, 66-keV 2_1^+ , 162-keV 3_1^+ , and 415-keV 4_1^+ states in ⁶⁶Ga and 0-keV 1_1^+ , 175-keV 2_1^+ , 376- and 676-keV $3_1^+ + 3_2^+$, and 496-keV 4_1^+ states in ⁶⁸Ga (Figs. 9 and 11). The IBFFM calculations are in accord with the classification of the parabolic rule, because the $1_1^+ + 1_2^+$, 2_1^+ , 3_1^+ , and 4_1^+ states in ⁶⁶Ga, as well as the 1_1^+ , 2_1^+ , $3_1^+ + 3_2^+$, and 4_1^+ states in ⁶⁸Ga have dominant (or at least very strong) $\pi \tilde{p}_{3/2} \nu \tilde{f}_{5/2}$ components. The inversion of the parabola is connected with the fact that the $\tilde{f}_{5/2}$ neutron is particle-like in $\frac{66}{31}$ Ga₃₇ ($V^2 > 0.5$).

The $\pi \tilde{p}_{3/2} \nu \tilde{f}_{5/2}$ multiplet is also seen in ⁷⁰Ga, and the energy splitting exhibits an open-down parabola, in agreement with the prediction of the parabolic rule.⁷

The members of the $\pi \tilde{p}_{3/2} \nu \tilde{f}_{5/2}$ multiplet in ⁷⁰As are fragmented into different states. Nevertheless, in the 1_1^+ , 3_4^+ , and 4_2^+ states the $\pi \tilde{p}_{3/2} \nu \tilde{f}_{5/2}$ components are dominant (Table III).

On the basis of lifetime measurements and other considerations Hübner⁴⁴ came to the conclusion that the 1^+_1 state of ⁷²As has either the $\pi p_{3/2}^{-1} \nu f_{5/2}^{-1}$ or the $\pi p_{3/2}^{-1} \nu p_{1/2}$ configuration. The IBFFM calculations give $\pi \tilde{p}_{3/2} \nu \tilde{f}_{5/2}$ as the dominant configuration for this state. The parabolic rule predicts

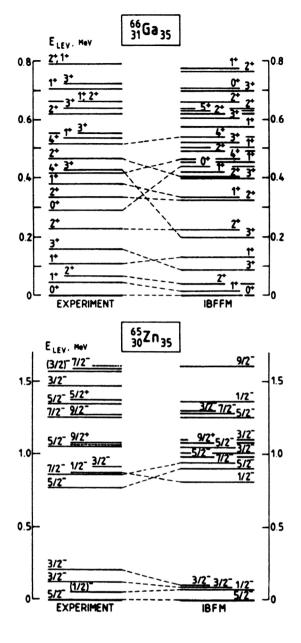


FIG. 9. Experimental energy spectra of ⁶⁵Zn (Ref. 31) and ⁶⁶Ga (Ref. 4) and the corresponding theoretical IBFM and IBFFM results.⁴

an open-down parabola for the energy splitting of the $\pi \tilde{p}_{3/2} \nu \tilde{f}_{5/2}$ multiplet, with a minimum energy for the 1⁺ state. The 4⁺ member of this multiplet may be the 4₁ state. The 2⁺ and 3⁺ members are fragmented into different states (e.g., 2_1^+ , 2_2^+ , etc.; see Table IX in Ref. 9).

The $\pi \tilde{p}_{3/2} \nu \tilde{p}_{1/2}$ doublet. In the case of doublets, for example, in $\pi \tilde{p}_{3/2} \nu \tilde{p}_{1/2}$, the energy splitting does not depend on the occupation probability, so that the shape of the splitting of the given multiplet is very similar for all investigated nuclei (see Fig. 7). The 1⁺ and 2⁺ members of the $\pi \tilde{p}_{3/2} \nu \tilde{p}_{1/2}$ doublet are clearly seen in ⁷⁰Ga (Ref. 7) and ⁷⁴As (Ref. 11). In both cases $E(1^+) < E(2^+)$, in agreement with the predictions of the parabolic rule and IBFFM calculations (for ⁷⁴As; see the 1⁺ and 2⁺ states in Fig. 17).

Some members of the $\pi \tilde{f}_{5/2} \nu \tilde{f}_{5/2}$ multiplet were identified in ⁷⁰As and ⁷²As.

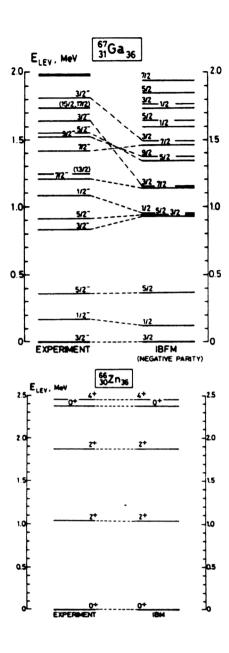


FIG. 10. Experimental energy spectra of ⁶⁶Zn (Ref. 30) and ⁶⁷Ga (Ref. 32) and the corresponding theoretical IBM and IBFM results. ^{5,6} The dashed lines show the assignment of the theoretical levels to the experimental ones. If the assignment is made on the basis of the energies only, it is marked by a dot-dash line.

According to the parabolic rule, the lowest-lying 0^+ states in 70 As are expected to be relatively pure, and they belong to the $\pi \tilde{f}_{5/2} \nu \tilde{f}_{5/2}$ and/or $\pi \tilde{p}_{1/2} \nu \tilde{p}_{1/2}$ quasiparticle multiplets. Many low-lying 1^+ levels are expected, with a strong configuration mixing among them. The same is true for the 2^+ and 3^+ states. The lowest-lying 4^+_1 state belongs to the $\pi \tilde{f}_{5/2} \nu \tilde{f}_{5/2}$ multiplet, and it is probably well separated from the 4^+ members of the $\pi \tilde{p}_{3/2} \nu \tilde{f}_{5/2}$ and $\pi \tilde{f}_{5/2} \nu \tilde{p}_{3/2}$ multiplets.

The IBFFM calculations (Fig. 12 and Table III) show that the 0_1^+ 345-keV and 4_1^+ ground states of 70 As have $\pi \tilde{f}_{5/2} \nu \tilde{f}_{5/2}$ as dominant configurations. The $\pi \tilde{f}_{5/2} \nu \tilde{f}_{5/2}$ components are fragmented in the 1_1^+ , 1_2^+ , 2_1^+ , 2_2^+ , 3_1^+ , 3_2^+ states with other components. The 0_1^+ state decays by a strong M1

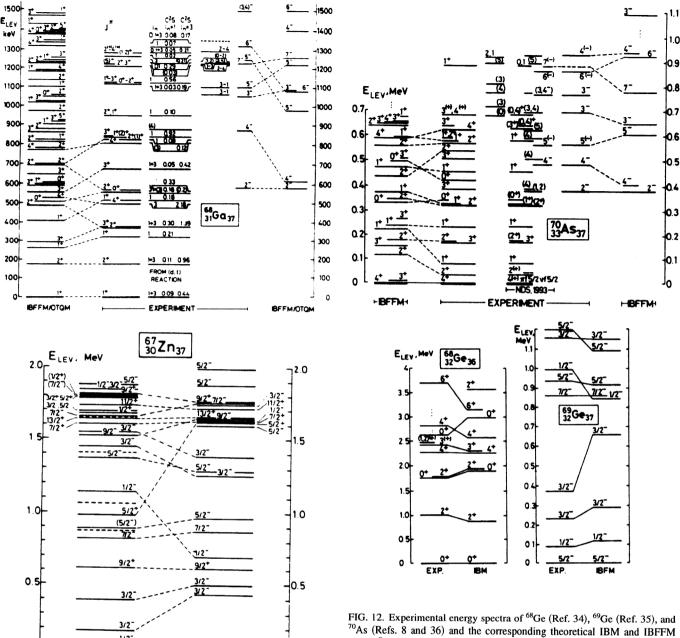


FIG. 11. Experimental energy spectra of 67 Zn (Ref. 32) and 68 Ga (Refs. 5 and 6) and the corresponding theoretical IBFM and IBFFM results. $^{5.6}$ The (d,t) reaction data are taken from Daehnick $et\ al.$ See also the caption to Fig. 10.

BFM

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transition to the 1_1^+ state, as expected. The low-lying 5^+ member of the $\pi \tilde{f}_{5/2} \nu \tilde{f}_{5/2}$ multiplet is missing in the experimental spectrum.

According to Bertschat *et al.*, ⁴⁵ the best agreement between the predicted and measured g factors of the 3_1^+ state at 214 keV in ⁷²As is obtained if one assumes the $\pi f_{5/2} \nu f_{5/2}^{-1}$ and $\pi f_{5/2} \nu p_{1/2}$ configurations. Our IBFFM calculations indicate a mixed configuration for this state, with the $\pi \tilde{f}_{5/2} \nu \tilde{f}_{5/2}$ configuration as the strongest one. The parabolic rule predicts an open-up parabola for the splitting of the

⁷As (Refs. 8 and 36) and the corresponding theoretical IBM and IBFFM results.⁸

 $\pi \tilde{f}_{5/2} \nu \tilde{f}_{5/2}$ multiplet, with a minimum energy for the 3⁺ (or 4⁺) member.

Some members of the $\pi \tilde{f}_{5/2} \nu \tilde{g}_{9/2}$ multiplet were observed in almost all the investigated odd-odd Ga and As nuclei.

On the basis of the parabolic rule we may expect that the 2_1^- , 3_1^- , 6_1^- , and 7_1^- states in $^{70}\mathrm{As}$ are relatively pure and probably belong to the $\pi \tilde{f}_{5/2} \nu \tilde{g}_{9/2}$ multiplet (see Fig. 6). In the low-lying 4_1^- and 5_1^- states there is probably a stronger configuration mixing. As Fig. 12 shows, the low-lying negative-parity states of $^{70}\mathrm{As}$ have been reasonably well reproduced by the IBFFM calculations. The 2_1^- , 3_1^- , 6_1^- , and 7_1^- levels are rather pure and belong mainly to the $\pi \tilde{f}_{5/2} \nu \tilde{g}_{9/2}$ multiplet (Table III). The 699-keV 3_1^- level decays by an M1 transition to the 2_1^- state, in accordance with expectations for the neighboring members of the same multiplet. The 4_1^- , 4_2^- , and 5_1^- states have mixed $\pi \tilde{p}_{1/2} \nu \tilde{g}_{9/2}$ and

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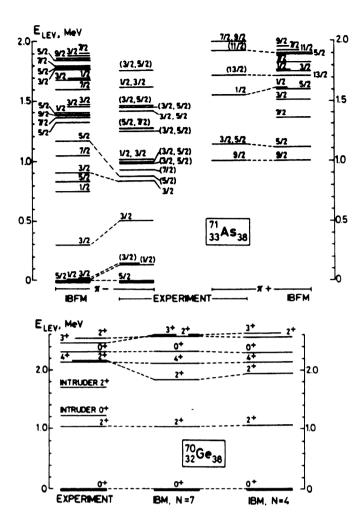


FIG. 13. Experimental energy spectra of 70 Ge (Ref. 36) and 71 As (Ref. 37) and the corresponding theoretical IBM and IBFM results. $^{9.13}$ The symbols π^+ and π^- denote positive- and negative-parity states, respectively.

 $\pi \tilde{f}_{5/2} \nu \tilde{g}_{9/2}$ multiplets as the main configurations.

According to the parabolic rule, the energy splitting of the $\pi \tilde{f}_{5/2} \nu \tilde{g}_{9/2}$ multiplet shows an open-down parabola with a minimum energy for the 2^- state in ⁷²As. On the basis of a measurement of the magnetic dipole moment by Hogervorst $et~al.^{46}$ one can conclude that the 2^- ground state of ⁷²As has predominantly a $\pi \tilde{f}_{5/2} \nu \tilde{g}_{9/2}$ configuration. The IBFFM calculations confirm these results: in the 2^-_1 ground, 3^-_1 , and 7^-_1 states the dominant configuration is $\pi \tilde{f}_{5/2} \nu \tilde{g}_{9/2}$ (Table IX in Ref. 9). In a recent publication Döring $et~al.^{47}$ assigned to the 562.8-keV level values $7^{(-)}$ for the spin and parity [instead of the former assignment (6)]. This state may be the 7^- member of the $\pi \tilde{f}_{5/2} \nu \tilde{g}_{9/2}$ multiplet (Fig. 14).

The 2^- and 7^- members of the $\pi \tilde{f}_{5/2} \nu \tilde{g}_{9/2}$ multiplet could be identified also in ⁶⁸Ga (Fig. 11 and Ref. 6) and ⁷⁰Ga (Ref. 7); and the 2^- member in ⁷⁴As (Fig. 16 and Ref. 11).

The 4⁻ and 5⁻ members of the $\pi \tilde{p}_{1/2} \nu \tilde{g}_{9/2}$ doublet were seen in ⁶⁸Ga (Refs. 5 and 6), ⁷⁰As (Ref. 8), and ⁷²As (Ref. 9), although they are mixed with other components. According to the parabolic rule, $E(4^-) < E(5^-)$ (see Figs. 6 and 7), in agreement with the experimental data and IBFFM calcula-

tions (Fig. 11 for ⁶⁸Ga, Fig. 12 and Table III for ⁷⁰As, and Fig. 14 and Table IX in Ref. 9 for ⁷²As).

In the low-lying 5⁻ and 6⁻ states, components of the $\pi \tilde{p}_{3/2} \nu \tilde{g}_{9/2}$ multiplet were seen in ⁶⁸Ga (Fig. 11 and Refs. 5 and 6) and ⁷⁰Ga (Ref. 7).

The *d*-boson composition of the IBFFM wave functions of some low-lying states in 66 Ga is given in Table IV. The 0_1^+ ground state of 66 Ga is basically of one-*d*-boson type. The total contribution of the sizable one-*d*-boson components is 75%. We note that this is an effect of the boson–fermion interaction, similar to the j-1 anomaly, which was studied previously in the cluster-vibration model for odd–even nuclei.²⁴ and appears also in the IBFM for odd–even nuclei.²⁴

The low-lying triplet of ⁷³Ge positive-parity levels, $9/2^+$, $5/2^+$, and $7/2^+$ (see Fig. 16), is associated also with the J = j - 1, j-2 anomaly due to lowering of the $|\tilde{g}_{9/2}, 12; 5/2\rangle$ and $|\tilde{g}_{9/2}, 12; 7/2\rangle$ one-*d*-boson multiplet states. The $5/2^+$ lowering is produced by the dynamical quadrupole interaction (and admixture of the $d_{5/2}$ configuration from the shell above), while the $7/2^+$ lowering is due to the exchange interaction.

The *electromagnetic moments* and the corresponding IB(FF)M results are given in Table V. Here we present calculated moments only for those states which have at least one experimentally measured moment (either electric quadrupole or magnetic dipole).

As Table V shows, the signs of the moments were correctly reproduced in 37 cases of the total of 39. (The exceptions are the electric quadrupole moments of the ⁷⁰As 4₁⁺ and ⁷⁰Ge 2₁⁺ states.) In 11 cases the signs of the moments were not determined experimentally. The calculations gave definite predictions for the signs, and allowed us to predict more than 200 moments which have so far not been measured (see Refs. 4, 9, and 11).

The calculated and available experimental B(E2), B(M1) reduced transition probabilities and γ -branching ratios for ⁶⁸Ga transitions are given in Table VI. Similar data have been obtained also for ⁶⁶Ga (Ref. 4), ⁷²As (Ref. 9), and ⁷⁴As (Ref. 11) transitions. As Table VI shows, there is reasonable agreement between the theory and experiment; at least all the leading branches are correctly reproduced. The γ -transition probabilities depend critically even on weak components of the wave function; we note a disagreement for the $3_1^+ \rightarrow 1_1^+$ transition.

In Table VII we present the experimental and calculated spectroscopic factors for the (d,t) transfer reaction populating the low-lying levels in 68 Ga. 6 Four levels (the 1_1^+ , 2_1^+ , $2_2^+ + 3_1^+$, and 4_1^+ states) are excited with large spectroscopic factors. According to the calculation, the low-lying 1_1^+ , 2_1^+ , 3_1^+ , and 4_1^+ states have the largest spectroscopic factors, in qualitative agreement with experiment. We have also calculated spectroscopic factors for the 75 As $(p,d)^{74}$ As and 73 Ge $(^{3}$ He, $d)^{74}$ As reactions. 11

Summarizing the results, we have described the energy spectra and electromagnetic properties of even-even core isotopes, odd-A neighbors, and odd-odd nuclei in a consistent way. For example, in the 64 Zn, 65 Zn, 65 Ga, and 66 Ga quartet more than 400 nuclear data (energy levels, moments, reduced transition probabilities, γ -branching ratios, etc.) have been calculated, using \lesssim 25 (more or less freely fitted)

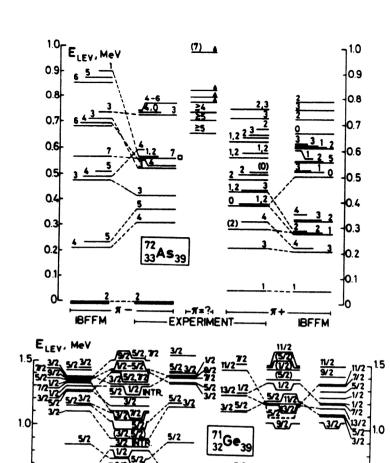


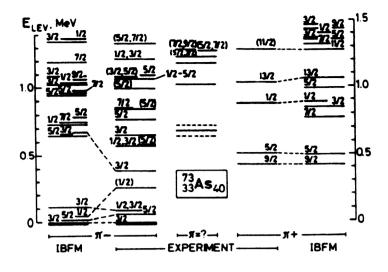
FIG. 14. Experimental energy spectra of 71 Ge (Ref. 37) and 72 As (Refs. 9 and 13) and the corresponding theoretical IBFM and IBFFM results. $^{9.13}$ The 72 As levels, marked with \triangle and \square , were taken from the (α, xnp) work of Mariscotti *et al.* 38 and Döring *et al.*, 47 respectively.

TABLE IV. The d-boson composition of the IBFFM wave functions of some low-lying states in 66 Ga.

$J_k^{\mathbf{x}}$	n_d								
	0	1	2	3	4				
0,	0.065	0.750	0.072	0.110	0.003				
I ₁ ⁺	0.344	0.466	0.121	0.065	0.004				
21	0.502	0.282	0.172	0.038	0.006				
12	0.339	0.460	0.132	0.064	0.005				
31	0.504	0.303	0.146	0.043	0.004				
22	0.381	0.396	0.163	0.054	0.006				
02	0.473	0.313	0.168	0.040	0.006				

For each state J_k^π , the numbers in the table give the value of $\sum_{j_\pi i_\nu j_{\pi\nu}} f \xi^2 [(j_\pi j_\nu) j_{\pi\nu}, n_d l; J]$ for each of the possible values of the d-boson number: $n_d = 0, 1, 2, \ldots, N$.

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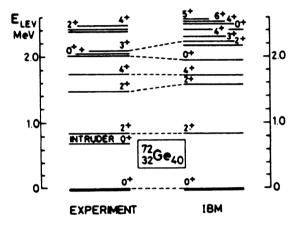


FIG. 15. Experimental energy spectra of ⁷²Ge (Ref. 39) and ⁷³As (Refs. 10 and 40) and the corresponding theoretical IBM and IBFM results. ^{11,13}

parameters. As the parameters of the calculations mostly show only rather small variations for the neighboring quartets (see Table I), the method allows a consistent description of a larger group of nuclei.

For odd-odd As nuclei a renormalization of the parameters was needed when we added one nucleon to the neighboring single-odd nuclei. Such a renormalization seems to be a consequence of the softness of the nuclei in this region. The appearance of an additional nucleon may change the dynamical deformation appreciably.

4.4. Comparison with earlier theoretical calculations

Theoretical interpretation of the structure of the ⁶⁶Ga and ⁶⁸Ga nuclei was completely missing before our work. ⁴⁻⁶

Using the number-conserving BCS quasiproton—quasineutron model, Ten Brink *et al.*⁵¹ calculated the energy spectra of ^{70,72,74,76}As and the electromagnetic properties of ^{70,72}As nuclei. They used a Schiffer force for the effective proton—neutron residual interaction. It was assumed that the odd—odd As nuclei are spherical and that the lowest states have the lowest seniority (two). The phonon degrees of freedom were neglected.

Kimura *et al.*⁵² calculated the level spectrum of ⁷²As, using harmonic-oscillator wave functions and a proton-neutron residual interaction of the form

$$V_{nn}(|\mathbf{r}_n - \mathbf{r}_n|) = V_0[(1 - \alpha) + \alpha(\boldsymbol{\sigma}_n \boldsymbol{\sigma}_n)] \delta(|\mathbf{r}_n - \mathbf{r}_n|).$$

In our IB(FF) calculation the following facts were taken into account:

- a) the Zn and Ge core nuclei may have a small effective deformation (the parameters h_2 and h_3 may differ from zero);
- b) the boson degree of freedom is important and cannot be neglected;
- c) the tensor residual interaction may play an important role in the description of some of the lowest-spin states $(0^+,1^+)$.

The experimental energy spectrum of ⁷⁴As is compared with the results of the present IBFFM and with previous theoretical calculations in Fig. 17. There are about 48 experimentally observed states in ⁷⁴As below 800 keV. The present IBFFM calculations reproduce 46 levels, while the earlier calculations⁵¹ give only 18 levels. A comparison with the previous theoretical results for ⁷²As is given in Ref. 9. Many

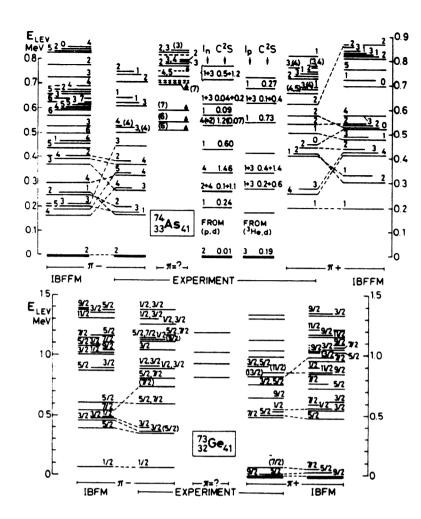


FIG. 16. Experimental energy spectra of 73 Ge (Ref. 40) and 74 As (Refs. 11 and 13) and the corresponding theoretical IBFM and IBFFM results. $^{11.13}$ The 74 As levels, marked with \triangle , were taken from García Bermúdez *et al.*; 41 the (p,d) and $(^{3}$ He,d) results are from Fournier *et al.* 42 and Rosner *et al.*, 43 respectively.

nuclear moments and reduced transition probabilities have been calculated for the first time in our work.^{9,11}

5. SUPERSYMMETRY IN 74 Se, 75 Se, 73 As, AND 74 As NUCLEI

On the basis of the vibrational-symmetry limit of the interacting boson-fermion model, Vervier *et al.*,⁵³ Čule and Paar,⁵⁴ and Van Isacker and Jolie⁵⁵ have developed formulas for the description of the level schemes of nuclei around ⁷⁶As.

The new, more complete level schemes of ⁷⁴As and ⁷³As obtained in our work¹¹ offered a new possibility of checking the validity of the supersymmetry scheme. According to this scheme, the energy spectra of four nuclei (in the present case ⁷⁴Se, ⁷⁵Se, ⁷³As, and ⁷⁴As) are interrelated and are described by the same Hamiltonian. The main advantage of this symmetry-based approximation is that the eigenvalue problem can be solved analytically.

According to Van Isacker and Jolie,⁵⁵ the eigenvalues of the Hamiltonian, the energies of the excited states, are given by the formula

$$E = A_{\pi\nu} \sum_{i} N_{i}(N_{i} + 7 - 2i) + A_{\pi} \sum_{i} N_{i\pi}(N_{i\pi} + 7 - 2i)$$

$$+ A_{\nu} \sum_{i} N_{i\nu}(N_{i\nu} + 7 - 2i) + B_{1} \sum_{i} n_{i} + B_{2} \sum_{i} n_{i}$$

$$\times (n_{i} + 6 - 2i) + C[v_{1}(v_{1} + 3) + v_{2}(v_{2} + 1)]$$

$$+ DL(L+1) + ES(S+1) + FJ(J+1), \qquad (11)$$

where $N_i, N_{i\pi}, N_{i\nu}, n_i, v_1, v_2, L, S, J$ are quantum numbers and $A_{\pi\nu}, A_{\pi}, A_{\nu}, B_1, B_2, C, D, E, F$ are parameters, which are not determined by the symmetry. The formula (11) was obtained for the U(5) limit of the $U_{\pi}(6/12) \otimes U_{\nu}(6/12)$ supersymmetry (SUSY) (proton particle, neutron hole case and $[N_i] \neq [N+1,1^5]$. (A somewhat different formula has been derived for $[N_i] = [N+1,1^5]$, but only a few levels belong to this group representation below 600 keV.)

First, we fitted the parameters of Eq. (11) to the levels of the even-even ⁷⁴Se, odd-A ⁷⁵Se, and ⁷³As nuclei by a least-squares method. Quantum numbers were assigned to the states of the supermultiplet on the basis of the energies, spins, parities, decay properties, available one-nucleon transfer-reaction spectroscopic factors, as well as the IBF(F)M and SUSY wave functions of the levels considered. The following parameters were obtained $A_{\pi} + A_{\pi\nu} = 55$, $A_{\nu} + A_{\pi\nu} = 26$, $B_{1} = 525$, $B_{2} = 0$, C = 4, D = -28, a high negative value for E, and F = 41 (all in keV). Then we used these

TABLE V. Experimental electric quadrupole (Q) and magnetic-dipole (μ) moments of Zn, Ga, Ge, and As nuclei compared with IB(FF)M theoretical results.

Nucleus	J*	E.	Q(eb)		μ(μ	N)		Main config.
		keV	Exp. [49]	Calc	Ref.	Exp. [49]	Calc	Ref.	comig.
64Zn ₃₄	21	992	-0.124(12)	-0.07		+0.84(18)	+0.94		One d-boson
			-0.143(21)			+0.92(20)			
65Zn ₃₅	5/2-	0	-0.023(2) ^a	-0.003		+0.7690(2)	a)+1.38		v j 5/2
				0.10	-+	0.78/20)	b)+0.84 a)-1 49		ν ρ̃ 3/2
	3/2	115		+0.10		-0.78(20)	b)-0.82		V p 3/2
	3/2-	207		-0.06		+0.73(25)	a)+0.52		v f 5/2@2+
	2				[4]		b)+0.43	[4]	
	9/2;	1066				-1 73(49)	a)-1.80		v] 9/2
							b⊢1.03		
66 31 Ga ₃₅	2;	66		-0 15		±1 011(18) ^h	a)+1 03		πφ̄ 3/2vj̄ 5/2
3							b)+0.72		
	7-	1464	±0.78(4) st	-0.40		+0.903(21)	a)-0.06		πj 5/2vg 9/2
							b)+0.70		
	9†	3043		-0 35		±4.23(90)	a)+5 12		
							b)+4.59		
66Zn 30 ^{Zn} 36	2†	1039		-0.064		+0 94(22)	+0.91		One d- boson
$^{67}_{30}$ Zn $_{37}$	5/2-	0	+0.150(15)*	+0.07		+0.8754790(84)	+1.35		vf 5/2
	1/2	93				+0.587(11)	+0 72		νρ̃ 1/2
	3/2	185		-0.04	[6]	+0.50(6)	+0.13	[6]	vƒ 5/2⊗2+
	9/2	604	= ±0.3[32]	-0 20		-1.097(9)	-1 02		vē 9/2
67 31 Ga ₃₆	3/2;	0	±0.195 st	+0.07		+1.8507(3)	+2.08		πρ̄ 3/2
68Ga ₃₇	I;	0	±0.0277(14)k,st	+0.008	2	±0.01175(5)k	-0.0134		πρ̄3/2v j̄ 5/2
	7-	1230	±0.72(2) ^{rt}	-0.514		+0.707(14)	+0.67		ng 5/2vg9/2
69 32 Ge 37	5/2	0	±0.024(5)"	+0.04		±0 735(7)	0.73		vj̃ 5/2
70 33 As ₃₇	41	0	+0.094(24)	-0.023	[8]	+2.1061(2)	+2.1	[8]	nj 5/2vj 5/2
	7-	888				+0.75(5) [50]	+0.77		πj 5/2∨g9/2
⁷⁰ Ge ₃₈	21	1039	+0.03(6) or +0.09(6)	-0.197		+0.936(52)	+0.914		One d-boson
71 32 Ge ₃₉	1/2	0				+0.547(5)	+0.445		νρ̃1/2
	5/2	175				+1.018(10)	+0.987		vj̃ 5∕2
	9/2	199	±0.34(5)	-0.230	[9]	-1.0413(7)	-1.221	[9]	v <u>₹</u> 9/2
71 As ₃₈	5/2	0	-0.021(6)	-0.156		(+)1.6735(18)	+1.010		rý 5/2
	9/2	1001		-0.365		+5.15(9)	+5.985		π <u>ē</u> 9∕2
72 33 As 39	27	0	-0.082(24)	-0.132		-2.1566(3)	-1.809		# 5/2∨89/
	3+	214		+0.229)	+1.580(18)	+1.525		RJ 5/2vJ 5/2
72 32 Ge ₄₀	21	834	-0.13(6)	-0.209)	+0.798(66)	+0.888		One d- boson
73Ge ₄₁	9/2	0	-0.173(26) ^a	-0.165	i	-0.8794677(2	-1 204		vg9/2
	5/2	13	-0 4(3) or +0.22 [58]	+0.230)	-0.0941(25)	-1 091		
73 33 ^{As} 40	5/2	67		-0.112	2 [11]	+1.63(10)	+0 964	[11]	πj 5/2
	9/2	428		-0.392	2	+5.234(14)	+5.955	3	πḡ9/2
74 As ₄₁	2-	0		+0.11	ı	-1.597(3)	-1.393	3	ng 5/2vg9/
	4†	259		+0.41	7	+3.238(40)	+3.79	3	πρ3/2vf 5/

a) $g_s^r = 0.9 g_s^r$ (free); b) $g_s^r = 0.5 g_s^r$ (free). h) Does not include Knight-shift correction. k) $\mu / Q < 0$ (signs of μ and Q are different). st) "Steinheimer" or other polarization correction included. u) No polarization correction included.

TABLE VI. Calculated E2 and M1 transitions between the low-lying levels in 68 Ga and comparison with available data. The assignment of calculated to experimental levels is made according to Fig. 11.

$J_i^{\mathfrak{X}} \to J_f^{\mathfrak{X}}$	10 ² B(E	$(2) (e^2b^2)$	10 B(M	11) (μ _N)		, ₁
IBFFM	IBFFM	Exp. [34]	IBFFM	Exp [34]	IBFFM	Exp. [34]
$2_1^+ \rightarrow 1_1^+$	0.13		1 822	≥ 0 014	100	100
$1_2^+ \rightarrow 2_1^+$	0.03		0.130		5.5	5.5
→ 1 ⁺ ₁	0.04	*	0 221		100	100
$2^{+}_{2} \rightarrow 1^{+}_{2}$	0.25		0.024		0.1	
→ 2 ⁺ ₁	0.002		0 044		119	3.0
$\rightarrow l_1^*$	0.00006		0.056		100	100
$3_1^{\star} \rightarrow 2_2^{\star}$	0.01		0.008		10-7	
→ 1 ⁺ ₂	0.0004				10-7	
$\rightarrow 2_1^+$	0.44		1.593	≥ 0.005	100	100
$\rightarrow l_1^+$	0.03	≥ 0.05			0,1	45
$4_1^+ \rightarrow 3_1^+$	0.40		1.389	≥ 0.038	100	100
$\rightarrow 2^{+}_{2}$	0.01				0.0008	
→ 2 ⁺ ₁	0.03	> 0 01			0.3	4.6
$1_3^{\bullet} \rightarrow 3_1^{\bullet}$	0.09				0.006	
$\rightarrow 2_2^+$	0.48		0 050		26	4.2
→ l ₂ *	0 23		0 036		5.0	1.0
→ 2 ⁺ ₁	0.03		0.134		100	100
$\rightarrow 1_1^+$	0.002		0.004		10.5	20
$0_1^* \rightarrow 1_3^*$			0.217		0.08	
$\rightarrow 2^{+}_{2}$	0.02				0.001	
→ 1 ⁺ ₂			0.925		60	15
$\rightarrow 2_1^+$	1.01				2.8	
$\rightarrow 1_1^+$			0.116		100	100
$2_3^+ \rightarrow 0_1^+$	0.00007				3×10^{-12}	
→ 1;	0.08		0.009		0.01	
→ 4 ⁺ ₁	0.004				5 × 10 ⁻⁶	
$\rightarrow 3^{+}_{1}$	0.23		0.271		21	0.9
$\rightarrow 2_2^+$	0.04		0.444		35	5.0
→ 1 ⁺ ₂	0.005		0.050		8.3	5.8
→ 2 ⁺ ₁	1.34		0.0003		9.8	13.4
→ I ⁺	0.20		0 044		100	100
$3_2^{\star} \rightarrow 2_3^{\star}$	0.12		0.377		0.7	
→ 1;	0.25		out when the out of		0.003	
→ 4 ⁺ ₁	0.22		0.00002		0.004	
→ 3 ⁺	0.19		0 0003		0.1	9.3
$\rightarrow 2_2^+$	0.05		0.065		2.5	
→ 1 ⁺ ₂	0.09				0.05	
→ 2 ⁺ ₁	0.19		0.561		100	100
→ 1 ⁺ ₁	0.04				0.6	2.8
$4_2^{\star} \rightarrow 3_2^{\star}$	1.06		0.167		2.1	
→ 2 ⁺ ₃	0.01				0.003	
$\rightarrow 4_1^{\bullet}$	0.06		0.583	≥ 0.0014	100	100
→ 3 ⁺ ₁	0.58		0.257	≥ 0.0004	121	90
→ 2 ⁺ ₂	0.02				0.1	
→ 2 ⁺ ₁	0.60				24	20

TABLE VI. (Continued.)

$J_i^{\mathfrak{X}} \to J_f^{\mathfrak{X}}$	102 B(E	$(2) (e^2b^2)$	10 B(N	il) (μ <mark>2</mark>)	1,		
IBFFM	IBFFM	Exp [34]	IBFFM	Ехр. [34]	IBFFM	Exp. [34]	
$2_4^+ \rightarrow 4_2^+$	0.32				4 × 10 ⁻⁸		
→ 3 ⁺ ₂	0.03		0.009		0.02		
→ 2 ⁺ ₃	0.26		0.062		0.6	2.4	
→ 0 ⁺	0.001				0.00005		
→ 1;	0.12		0.034		0.6		
→ 4 ⁺	0.21				0.03		
→ 3 ⁺ ₁	0.66		0.029		1.9	8.6	
→ 2 ⁺ ₂	0.002		0.013		0.6	4.6	
→ 1 ⁺ ₂	0.006		0.032		2.2	10.5	
→ 2 ⁺	0.0003		0.026		3.8	4.7	
→ 1 ⁺ 1	0.93		0.291		100	100	
$4\overline{1} \rightarrow 2\overline{1}$	2.69	≥ 0.21			100	100	
$3\frac{1}{1} \rightarrow 4\frac{1}{1}$	1.01		1 212		1.1		
→ 2 ₁	2.10		5.934		100	100	
$5\frac{1}{2} \rightarrow 3\frac{1}{1}$	0.004				10-6		
→ 4 ₁	0.19		0 579		100	100	
$7\frac{1}{1} \rightarrow 5\frac{1}{2}$	0.21				100	100	
$5\frac{1}{1} \rightarrow 7\frac{1}{1}$	4.27				10⊸		
→ 5 ₂	0.18		0.304		1.2		
→ 3 ₁	1.36				0 03		
→ 4 ₁	0.32		1.533		100	100	
6 ₁ → 5 ₁	0.04		0.072		0.06	≤ 17	
→ 7 ₁	0.02		0.076		0.1		
→ 5 ₂	0.47		5.019		100	100	
$\rightarrow 4^{-}_{1}$	0.96				2.2	8	

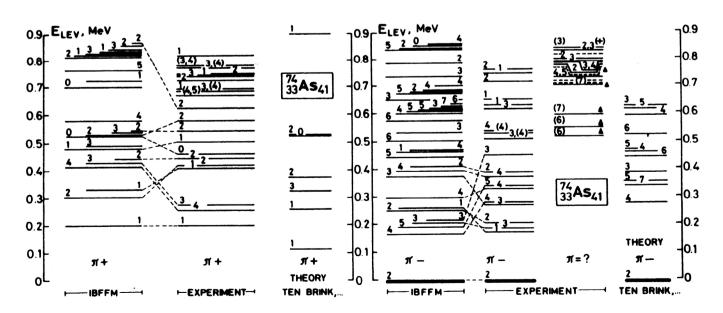


FIG. 17. Level scheme of 74 As (Ref. 11), compared with the present IBFFM and previous calculations by Ten Brink *et al.*⁵¹ The levels marked with triangles were observed in Ref. 41 from the $(\alpha, np\gamma)$ reaction.

363

TABLE VII. Spectroscopic factors for the $^{69}_{31}\mathrm{Ga}_{38}(d,t)^{68}_{31}\mathrm{Ga}_{37}$ transfer reaction populating the low-lying levels in $^{68}\mathrm{Ga}$.

				C2S			
J*		IBFN	И			Exp [33]	
IBFFM	P _{1/2}	P 3/2	f _{5/2}	89/2	$p_{1/2} + p_{3/2}$	f _{\$22}	89/2
11	0.012	0.003	3.333		0.09	0.44	
21	0.023	0.002	3.230		0.11	0.96	
12	0.308	0.015	0.170		0.21		
2*	0.344	0.430	0.063		0.30	1.39	
3 ⁺		0.114	2.772) 5.55)	
4+			3 227			2.18	
13	0.055	0.283	0 032		0.18		
0†		0.403			0.18	0.2	
2*	0.004	0.210	0.007		0.33		
3 ⁺ ₂		1.050	0.278		0.05	0.42	
4 2			0.438			0.12	
24	0.005	0.0004	0.178		0.09		
4-				0.094			
3-				0.020			
5-2				0.476			
5-				0.003			
6 ₁				0.360			

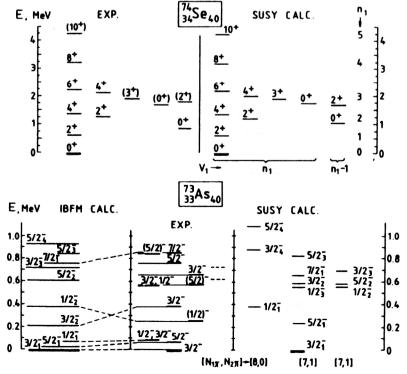


FIG. 18. Upper part: experimental energy levels of ⁷⁴Se (Ref. 56) in comparison with the SUSY calculations. Lower part: experimental energy levels of ⁷³As (Refs. 10 and 40) in comparison with the IBFM and SUSY calculations.

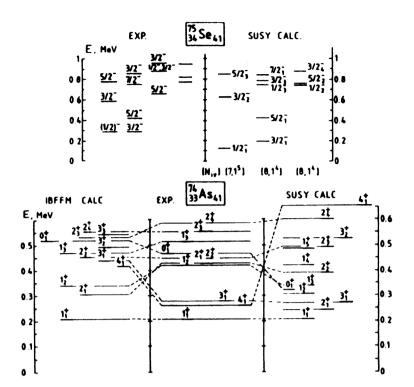


FIG. 19. Upper part: experimental energy levels of ⁷⁵Se (Ref. 57) in comparison with the SUSY calculations. Lower part: experimental energy levels of ⁷⁴As (Ref. 11) in comparison with the IBFFM¹¹ and SUSY¹² calculations. If the assignment of calculated levels to experimental ones is made on the basis of the energies only, it is marked with a dot-dash line.

parameters to generate the level spectrum of 74 As. (The value of $A_{\pi\nu}$ was taken as zero, after testing its role in the generated 74 As spectrum.)

The experimental and calculated energy spectra are compared in Figs. 18 and 19.

The supersymmetry calculations describe 44 levels of four different nuclei reasonably well with only seven fitted parameters. However, a problem appears with additional 1⁺ and 3⁺ states in the low-energy SUSY spectrum, and the assignment of some SUSY states to experimental ones cannot be made unambiguously.

The supersymmetry calculations and results are described in detail in Ref. 12.

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