

*Two neutron separation
energies and phase transition in
the Interacting Boson Model.*

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**“Mapping the triangle”, Jackson Lake Lodge,
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Motivation

- Renewed interest in the study of transitional regions and phase transitions.
- The traditional IBM analysis of phase transitions can be improved.
- Better understanding of the evidences of phase transitions.
- Possible connection between excited states and ground state properties.
- Connection with shape and configuration mixing.

Generic Hamiltonian

$$\begin{aligned}\hat{H} = & \varepsilon_s \hat{n}_s + \varepsilon_d \hat{n}_d + \kappa_0 \hat{P}^\dagger \hat{P} + \kappa_1 \hat{L} \cdot \hat{L} \\ & + \kappa_2 \hat{Q} \cdot \hat{Q} + \kappa_3 \hat{T}_3 \cdot \hat{T}_3 + \kappa_4 \hat{T}_4 \cdot \hat{T}_4,\end{aligned}$$

where \hat{n}_s and \hat{n}_d are the s and d boson number operators, respectively, and

$$\begin{aligned}\hat{P}^\dagger &= \frac{1}{2} d^\dagger \cdot d^\dagger - \frac{1}{2} s^\dagger \cdot s^\dagger, \\ \hat{L} &= \sqrt{10} (d^\dagger \times \tilde{d})^{(1)}, \\ \hat{Q} &= s^\dagger \tilde{d} + d^\dagger \tilde{s} + \chi (d^\dagger \times \tilde{d})^{(2)}, \\ \hat{T}_3 &= (d^\dagger \times \tilde{d})^{(3)} \\ \hat{T}_4 &= (d^\dagger \times \tilde{d})^{(4)}.\end{aligned}$$

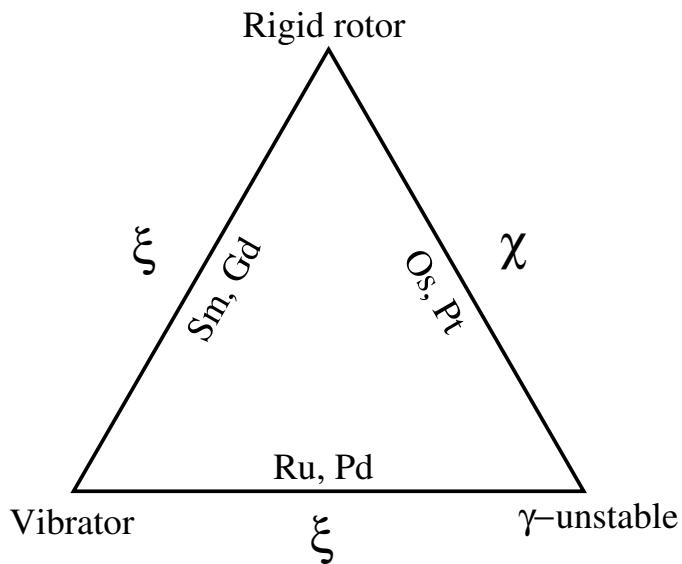
Transitional regions in the IBM

$$\hat{H} = \kappa \left(N \frac{1 - \xi}{\xi} \hat{n}_d - \hat{Q} \cdot \hat{Q} \right) + \kappa' \hat{L} \cdot \hat{L}$$

$\hat{n}_d = d$ boson number operator

$$\hat{L} = \sqrt{10} (d^\dagger \times \tilde{d})^{(1)},$$

$$\hat{Q} = s^\dagger \tilde{d} + d^\dagger \tilde{s} + \chi (d^\dagger \times \tilde{d})^{(2)}$$



Intrinsic state in the IBM

- The trial wave function

$$|c\rangle = \frac{1}{\sqrt{N!}} (\Gamma_c^\dagger)^N |0\rangle,$$

where

$$\Gamma_c^\dagger = \frac{1}{\sqrt{1 + \beta^2}} \left(s^\dagger + \beta \cos \gamma d_0^\dagger + \frac{1}{\sqrt{2}} \beta \sin \gamma (d_2^\dagger + d_{-2}^\dagger) \right).$$

- The energy surface

$$\begin{aligned} \langle c | H | c \rangle &= \frac{N}{5(1 + \beta^2)} \left(5\varepsilon_s + 25\kappa_2 + \beta^2 (5\varepsilon_d - 3\kappa_1 + 5\kappa_2 + 5\chi^2\kappa_2 \right. \\ &\quad \left. - 7\kappa_3 + 9\kappa_4) \right) \\ &+ \frac{N(N-1)}{140(1 + \beta^2)^2} \left(35\kappa_0 + \beta^2 (-70\kappa_0 + 560\kappa_2) \right. \\ &\quad \left. - 80\sqrt{14}\beta^3\chi \cos(3\gamma)\kappa_2 + \beta^4 (35\kappa_0 + 40\chi^2\kappa_2 + 72\kappa_4) \right). \end{aligned}$$

- The symmetry limits

$U(5)$ limit $\rightarrow \beta = 0$.

$SU(3)$ limit $\rightarrow \beta = \sqrt{2}, \gamma = 0, \pi/3$.

$O(6)$ limit $\rightarrow \beta = 1, \gamma$ unstable nucleus.

S_{2n} calculations using IBM

- Definition:

$$S_{2n}(N) = BE(N) - BE(N - 1).$$

- “Global” part of the IBM Hamiltonian: It is related with the Casimir operators of the U(6) group.

$$\hat{H}^{gl} = -E_0 - \mathcal{A} - \frac{\mathcal{B}}{2}\hat{N}(\hat{N} - 1).$$

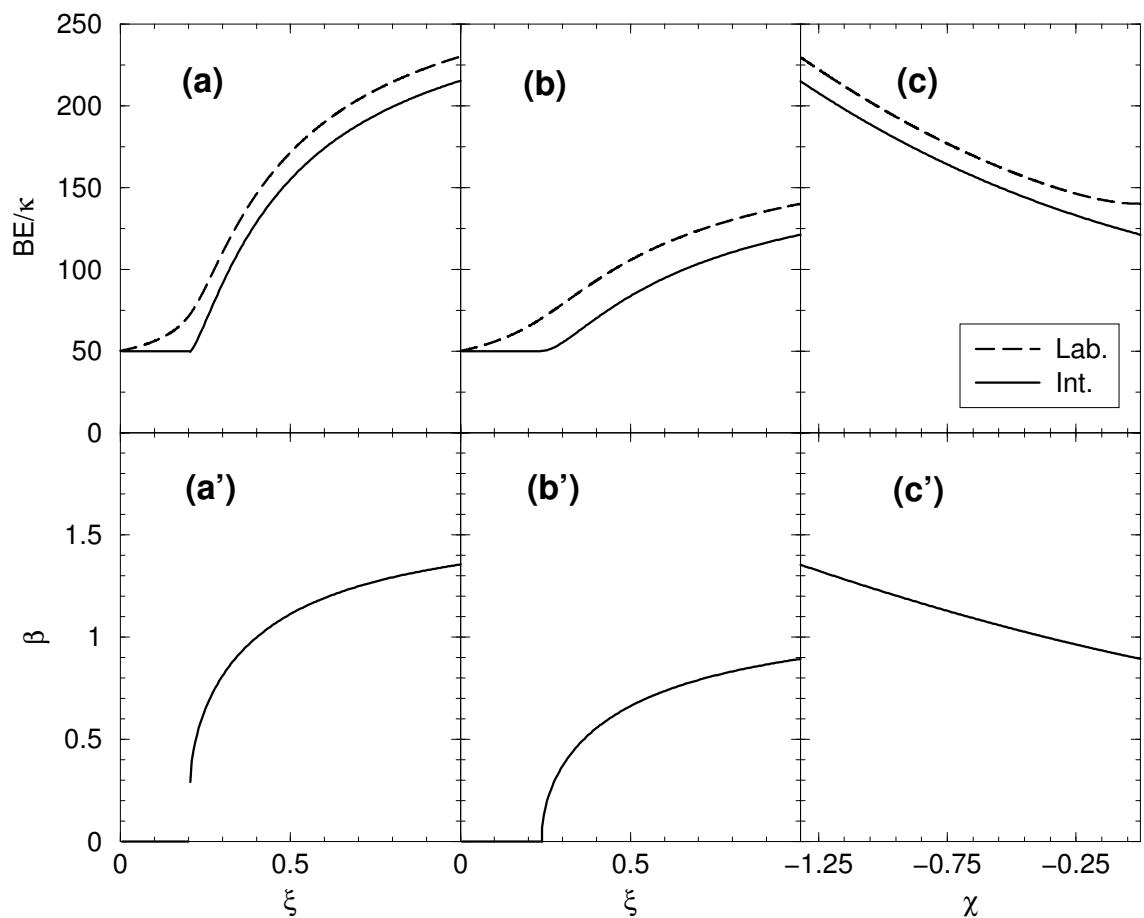
$$BE^{gl}(N) = E_0 + \mathcal{A}N + \frac{\mathcal{B}}{2}N(N - 1).$$

$$S_{2n}^{gl}(N) = (\mathcal{A} - \mathcal{B}/2) + \mathcal{B}N.$$

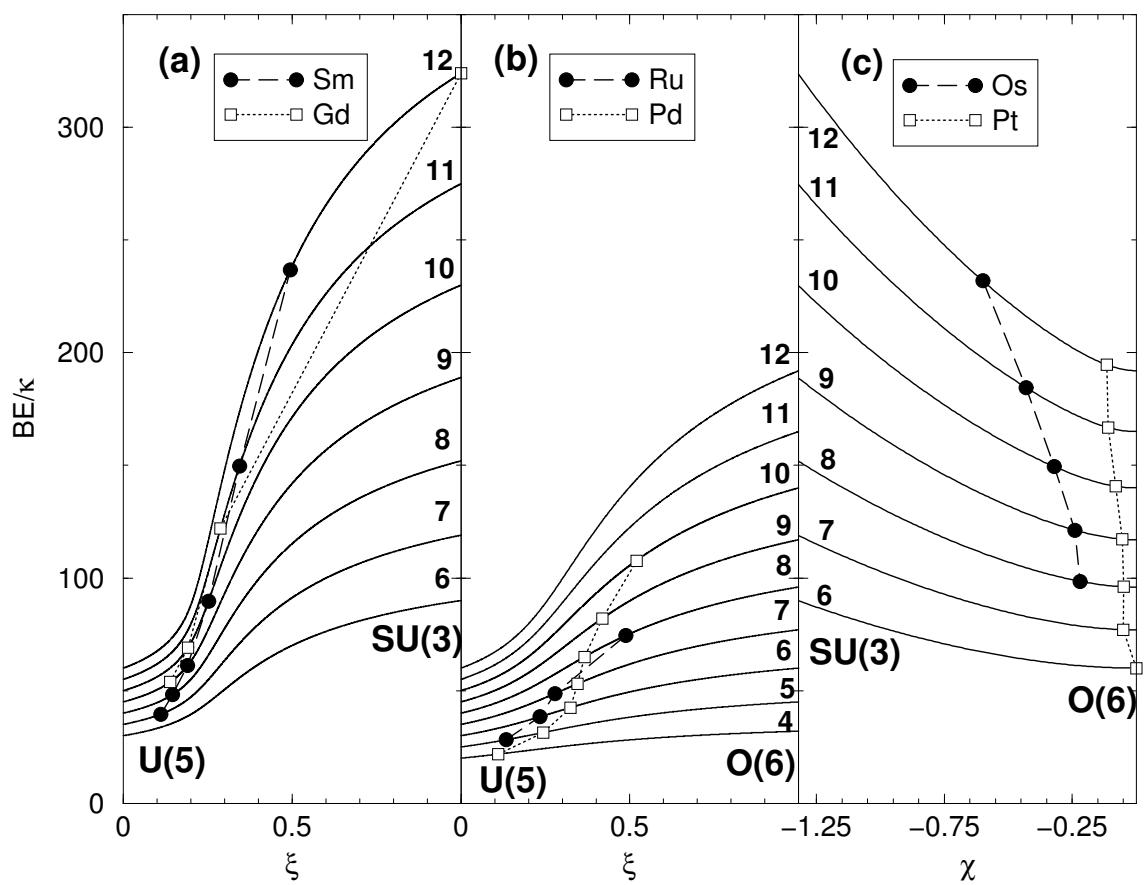
$$S_{2n} = S_{2n}^{gl} + S_{2n}^{lo}.$$

- \mathcal{A} and \mathcal{B} are kept as constants for a chain of isotopes at a given major shell.

Standard diagram “Binding energy vs. control parameter”



New diagram “Binding energy vs. N-control parameter”

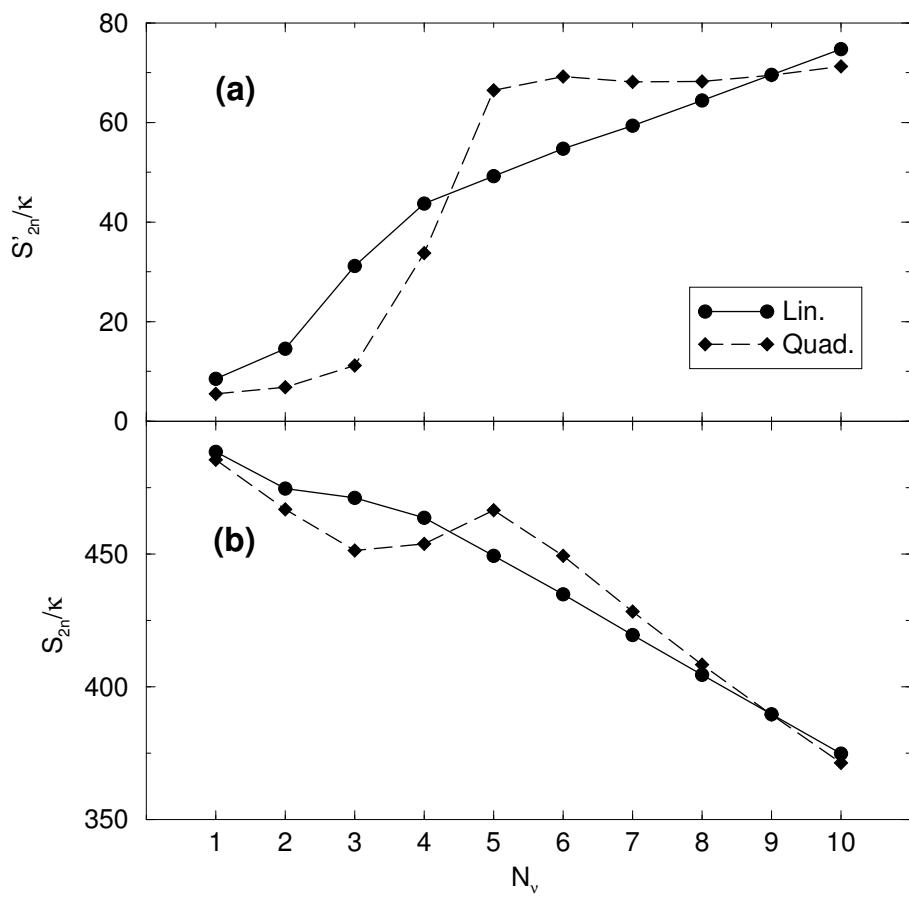


Simulation of a phase transition I

- Relation between the control parameter and the number of bosons.

$$\begin{aligned}\xi_{lin} &= 0.099N_\nu + 0.01, \\ \xi_{qua} &= 0.0099N_\nu^2 + 0.01.\end{aligned}$$

- Linear part for S_{2n} : $S_{2n}^{lin}/\kappa = 200 - 20N_\nu$.
- $U(5) - SU(3)$.



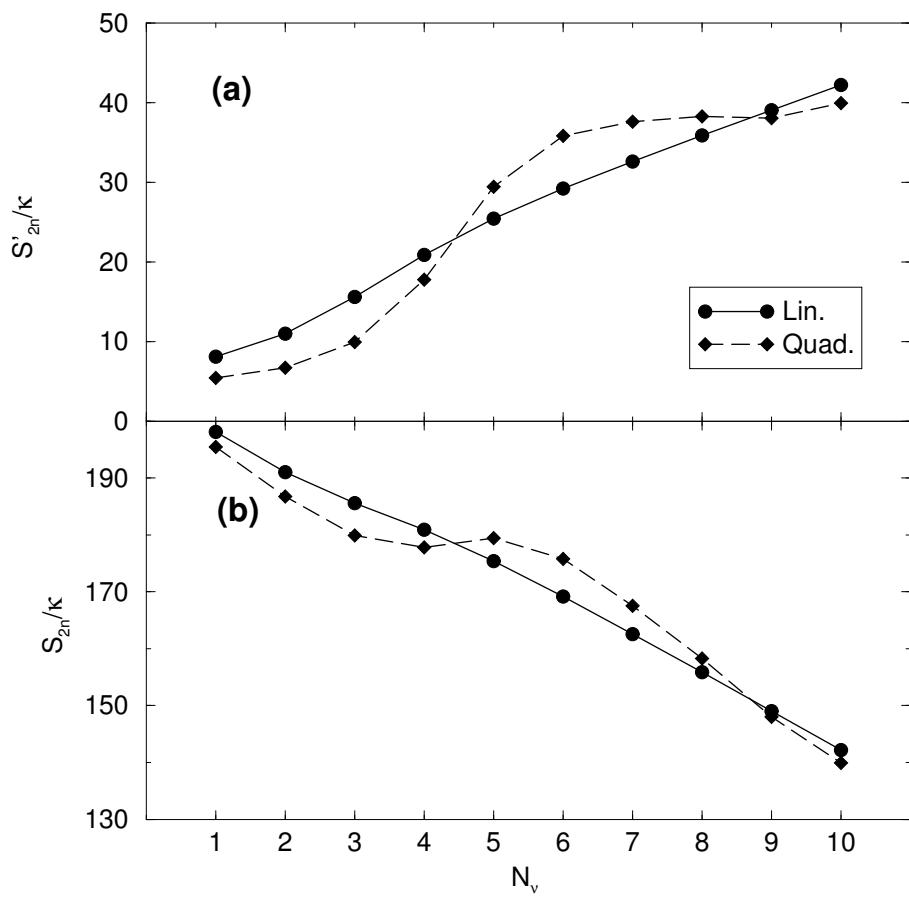
Simulation of a phase transition II

- Relation between the control parameter and the number of bosons.

$$\xi_{lin} = 0.099N_\nu + 0.01,$$

$$\xi_{qua} = 0.0099N_\nu^2 + 0.01.$$

- Linear part for S_{2n} : $S_{2n}^{lin}/\kappa = 200 - 10N_\nu$.
- $U(5) - O(6)$

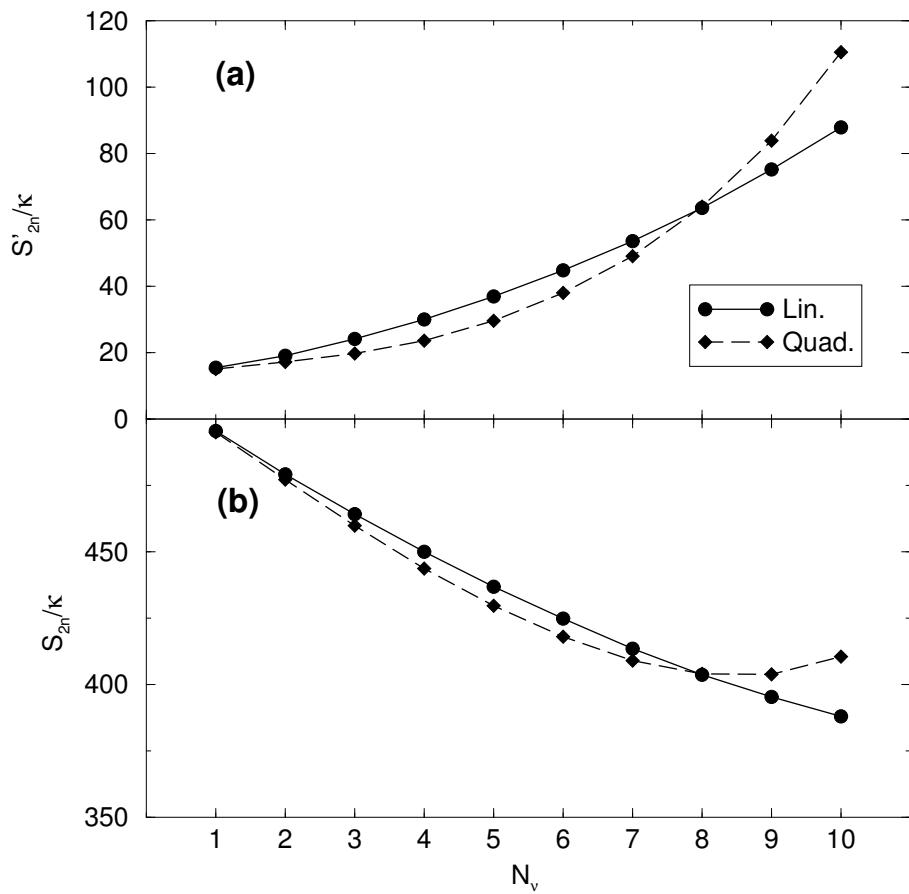


Simulation of a phase transition III

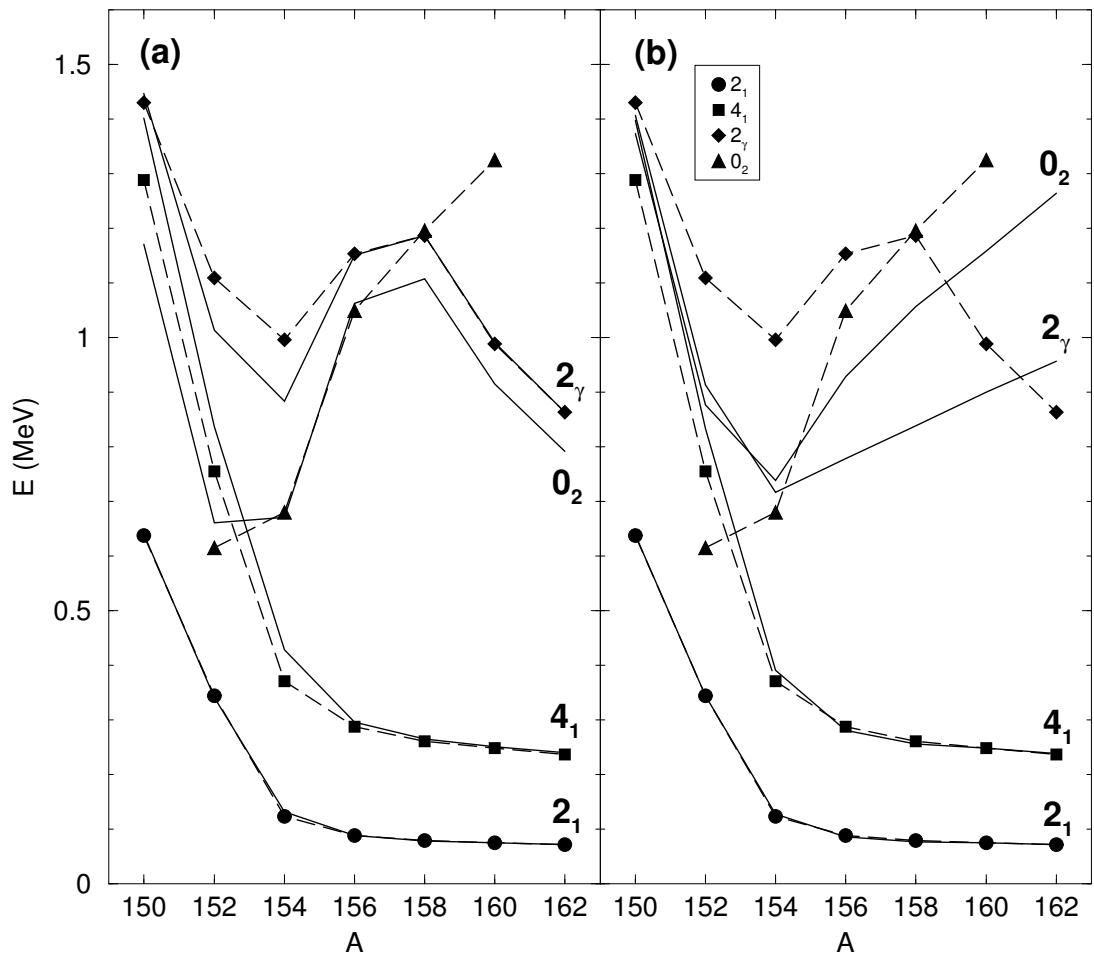
- Relation between the control parameter and the number of bosons.

$$\chi_{lin} = -\frac{\sqrt{7}}{20}N_\nu, \quad \chi_{qua} = -\frac{\sqrt{7}}{200}N_\nu^2.$$

- Linear part for S_{2n} : $S_{2n}^{lin}/\kappa = 200 - 20N_\nu$.
- $SU(3) - O(6)$



Gd isotopes: spectra



Gd isotopes: parameters of the Hamiltonian

Theo.-a

A	150	152	154	156	158	160	162
N_ν	2	3	4	5	6	7	8
κ	15.4	15.4	15.4	15.4	14.8	11.3	9.1
ξ	0.139	0.192	0.287	1	1	1	1
κ'	9.0	9.0	9.0	9.0	7.7	8.3	8.6

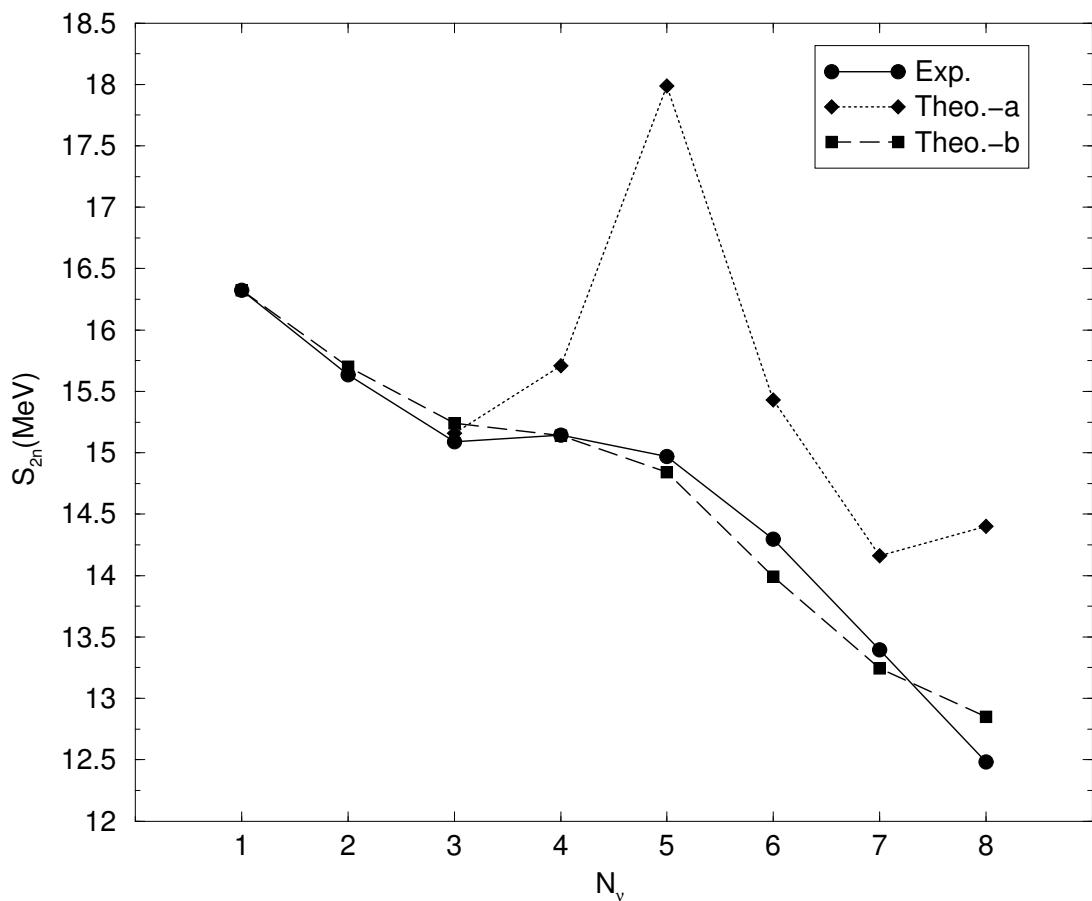
κ, κ' in keV and ξ dimensionless, $\chi = -\sqrt{7}/2$.

Theo-b

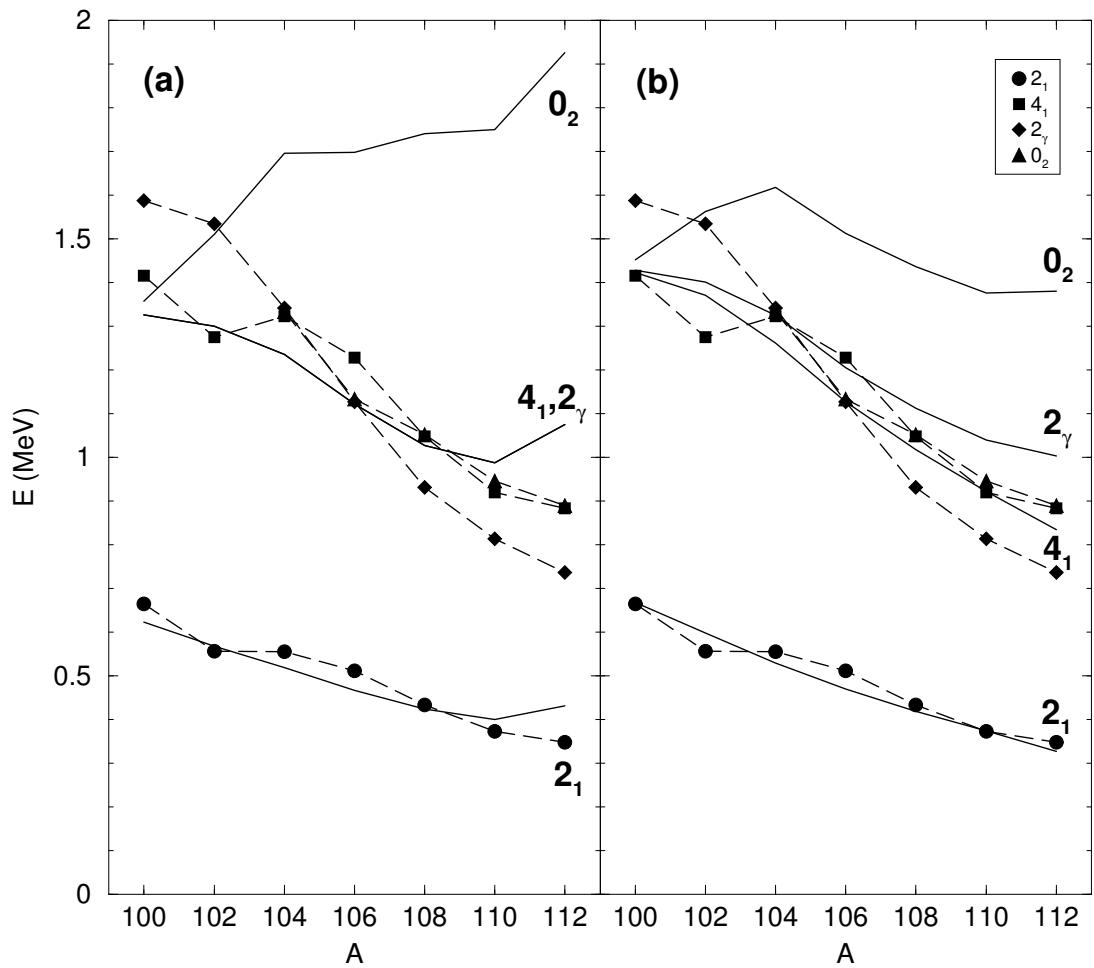
A	146	148	150	152	154	156	158	160	162
N_ν	0	1	2	3	4	5	6	7	8
ξ	0.60	0.137	0.166	0.236	0.373	0.535	0.625	0.658	0.724

$\kappa = 19.2$ keV, $\kappa' = 0$ and ξ dimensionless, $\chi = -0.6$.

Gd isotopes: S_{2n}



Pd isotopes: spectra



Pd isotopes: parameters of the Hamiltonian

Theo.-a

A	100	102	104	106	108	110	112
N_ν	2	3	4	5	6	7	8
κ	20.0	42.0	52.0	49.0	47.0	52.0	70.0
ξ	0.110	0.244	0.324	0.345	0.366	0.419	0.52

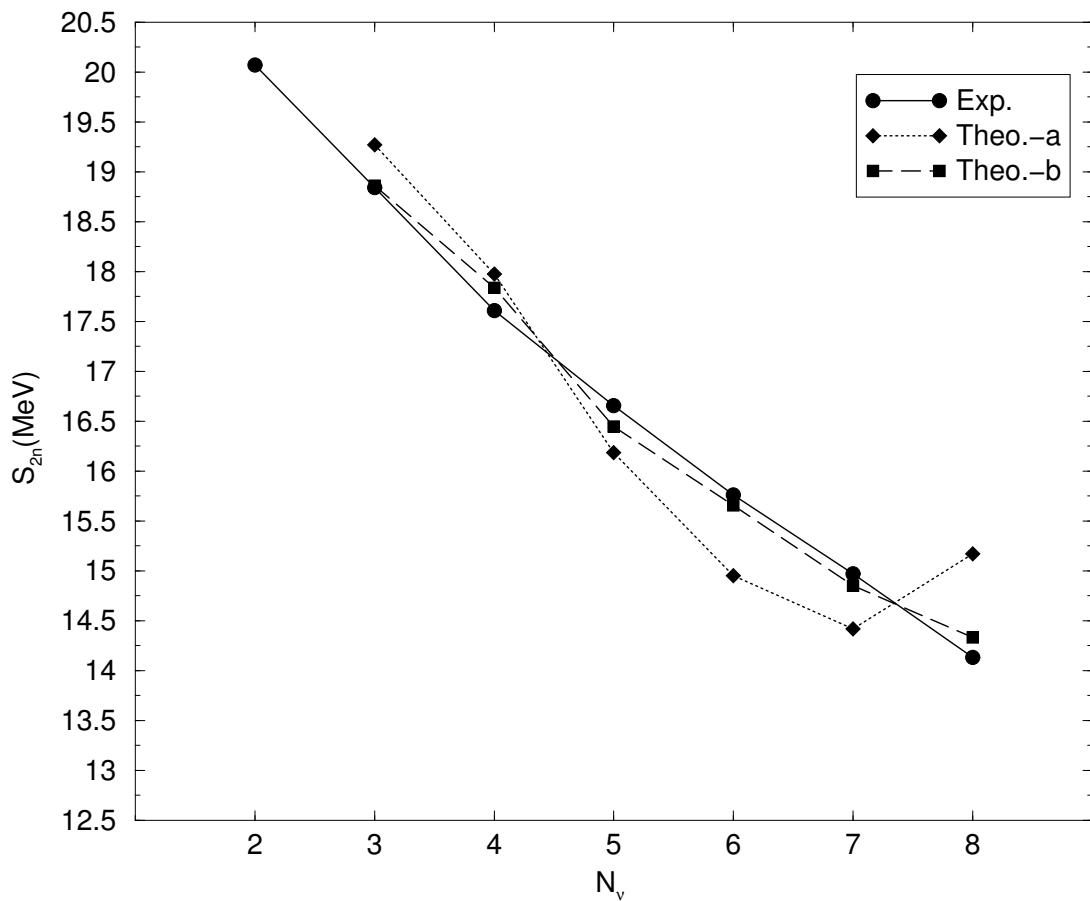
κ in keV, $\kappa' = 0$ and ξ dimensionless, $\chi = 0$.

Theo-b

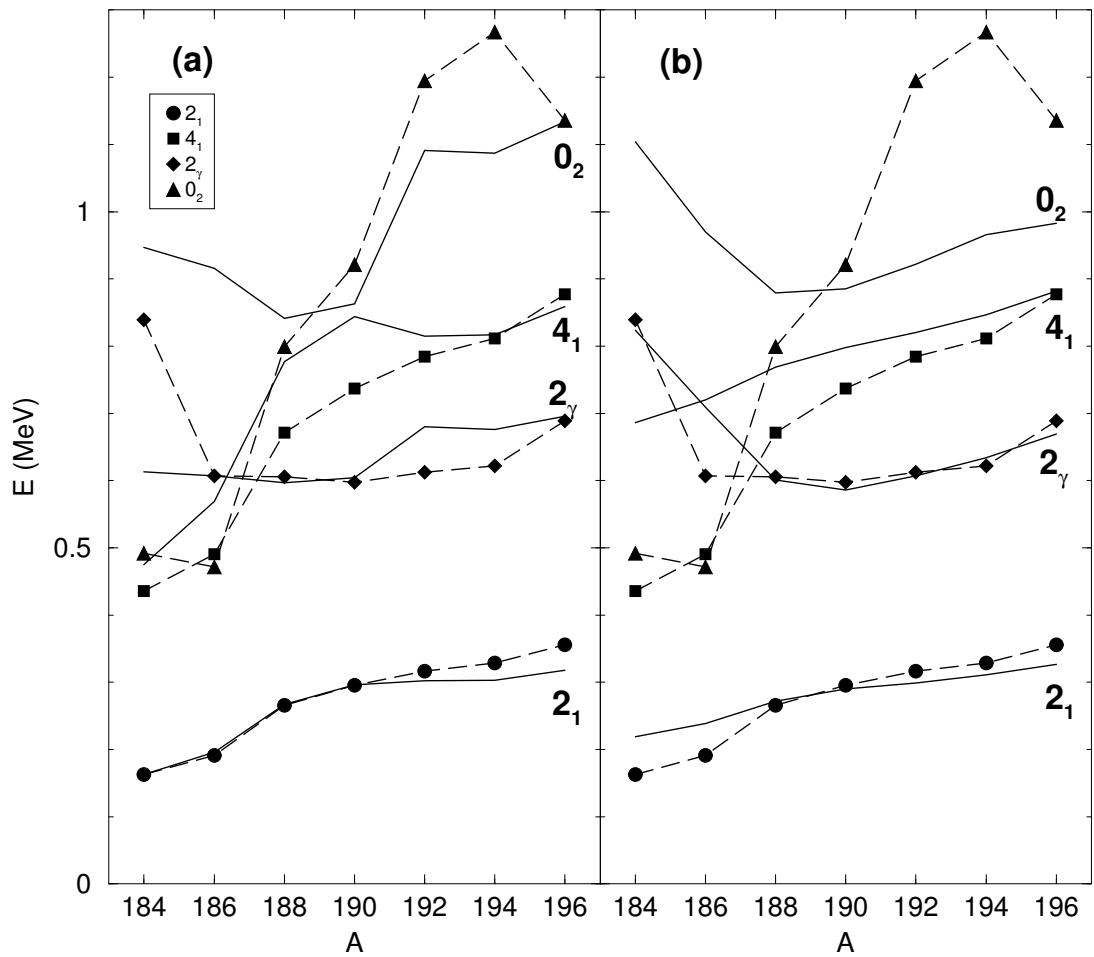
A	100	102	104	106	108	110	112	114
N_ν	2	3	4	5	6	7	8	9
κ	22.0	44.0	50.0	44.0	40.0	37.0	37.0	33.0
ξ	0.112	0.239	0.300	0.306	0.314	0.322	0.346	0.342

κ in keV, $\kappa' = 0$ and ξ dimensionless, $\chi = -0.3$.

Pd isotopes: S_{2n}



Pt isotopes: spectra



Pt isotopes: parameters of the Hamiltonian

Theo.-a

A	184	186	188	190	192	194	196
N_ν	10	9	8	7	6	5	4
κ	43.0	44.0	44.0	47.0	60.0	60.0	63.0
χ	-0.115	-0.110	-0.080	-0.055	-0.049	-0.050	0
κ'	4.2	0.8	17.6	19.0	11.0	11.0	11.0

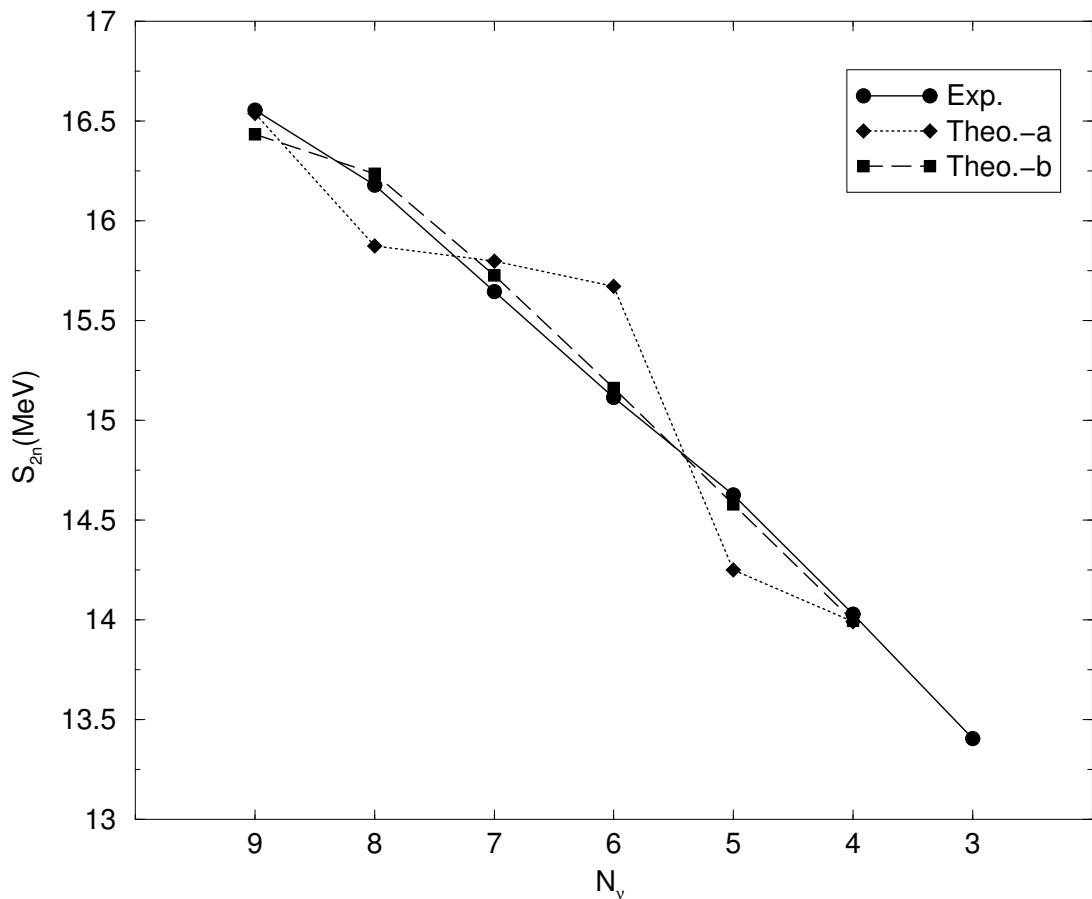
κ and κ' in keV and ξ dimensionless, $\varepsilon_d = 0$.

Theo-b

A	184	186	188	190	192	194	196
N_ν	10	9	8	7	6	5	4
κ	33.5	33.5	33.5	33.5	33.5	33.5	33.5
χ	-0.25	-0.20	-0.10	0	0	0	0
κ'	15.2	15.2	15.2	15.2	15.2	15.2	15.2

κ and κ' in keV and ξ dimensionless, $\varepsilon_d = 25.7$ keV.

Pt isotopes: S_{2n}



Conclusions

- A new approach, that includes the connection between the control parameter and N , has been presented.
- The crossing of a phase transition point produces a deviation in the S_{2n} values from the overall linear background.
- The study of long chains of isotopes should include the analysis of ground state properties, *i.e.* binding energies.
- This method can be used for calculating S_{2n} values in extended areas.