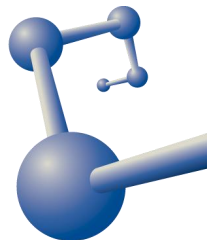


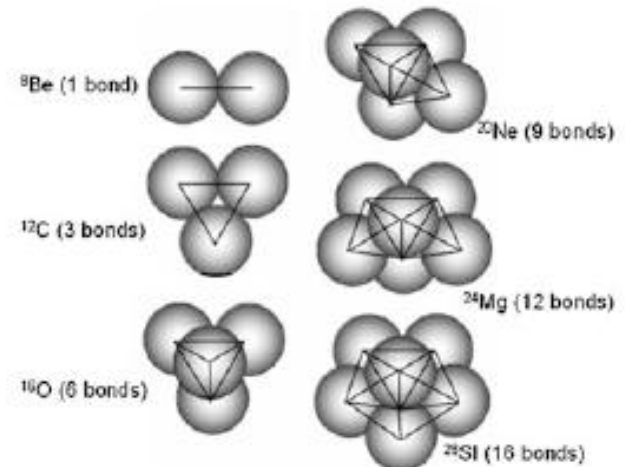
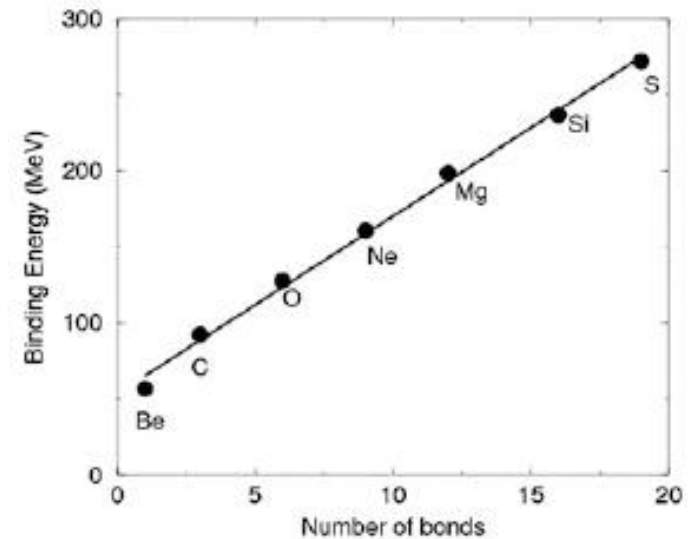
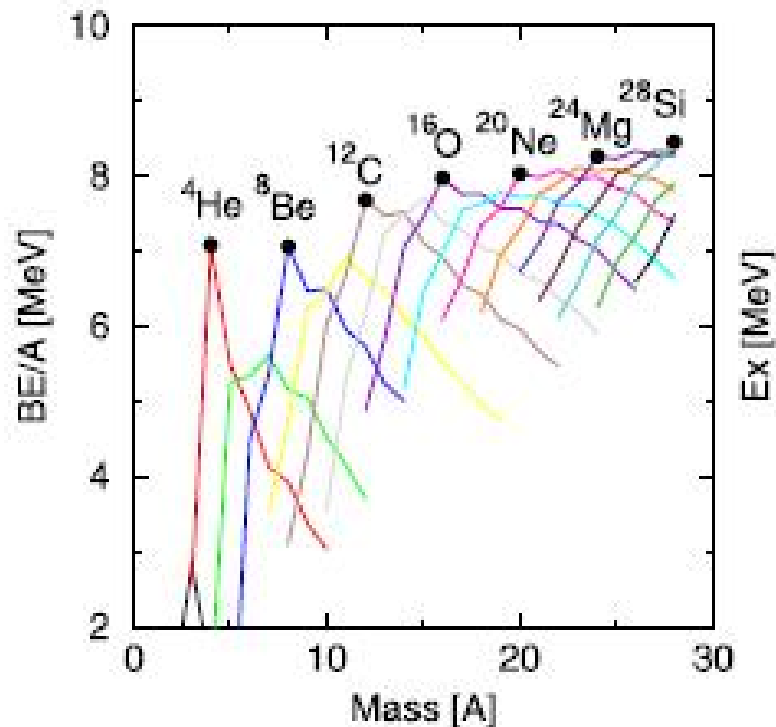
Algebraic Treatment of Alpha-Cluster Nuclei

- Introduction
- Algebraic Cluster Model (ACM)
- Point group symmetry
- Applications: ^{12}C
- Cluster Shell Model (CSM)
- Double point group symmetry
- Applications: ^{13}C
- Summary and conclusions

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Alpha-Cluster Nuclei



3 (Left panel) Binding energy per nucleon of the

Recent Theoretical Work

- Alpha-cluster model (Wheeler 1937, Brink 1966, Robson 1978)
- AMD (Kanada-Enyo, PTP, 2007)
- FMD model (Chernykh et al, PRL, 2007)
- BEC-like cluster model (Funaki et al, PRC, 2009)
- Ab initio no-core shell model (Roth et al, PRL, 2011)
- Lattice EFT (Epelbaum et al, PRL, 2011, 2012)
- No-core symplectic model (Dreyfuss et al, PLB, 2013)
- Algebraic Cluster Model (2000, 2002, 2014, 2017, 2023)
- and many others
- Recent reviews: Freer & Fynbo, PPNP 78, 1 (2014), Freer, Horiuchi, Kanada-En'yo, Lee & Meissner, RMP 90, 035004 (2018)

Few-Body Systems

- Integro-differential methods
- Algebraic Cluster Model (ACM)
- For k dof introduce a SGA of $U(k+1)$
- Two-body systems: $U(4)$ vibron model (diatomic molecules, mesons)
- Three-body systems: $U(7)$ model (baryons, nuclear clusters, molecules)
- Four-body systems: $U(10)$ model
- n -body systems: $U(3n-2)$ model

ACM for n-Body Systems

$3n-3$ relative degrees of freedom: Jacobi coordinates

$$\vec{\rho}_k = \frac{1}{\sqrt{k(k+1)}} \left(\sum_{i=1}^k \vec{r}_i - k\vec{r}_{k+1} \right), \quad k = 1, 2, \dots, n-1$$

Introduce $n-1$ dipole bosons and a scalar boson

$$b_{k,m}^\dagger, s^\dagger, \quad k = 1, 2, \dots, n-1$$

Number conserving one- and two-body Hamiltonian

$$\begin{aligned} H = & \epsilon_0 s^\dagger \tilde{s} - \epsilon_1 \sum_k b_k^\dagger \cdot \tilde{b}_k + u_0 s^\dagger s^\dagger \tilde{s} \tilde{s} \\ & - u_1 \sum_k s^\dagger b_k^\dagger \cdot \tilde{b}_k \tilde{s} + v_0 \sum_k [b_k^\dagger \cdot b_k^\dagger \tilde{s} \tilde{s} + \text{h.c.}] \\ & + \sum_L \sum_{ijkl} v_{ijkl}^{(L)} [b_i^\dagger \times b_j^\dagger]^{(L)} \cdot [\tilde{b}_k \times \tilde{b}_l]^{(L)}, \end{aligned}$$

Mixing between
oscillator shells

Wave Functions

- Labeled by $[N], \alpha, L_t^P \rangle$
- Total number of bosons: N
- Angular momentum: L
- Parity: P
- Identical clusters: Permutation symmetry: $\dagger$



Roelof Bijker, ICN-UNAM

Three-Body Clusters

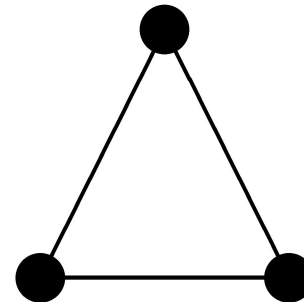
Permutation symmetry

$$\begin{aligned} H = & \xi_1 (R^2 s^\dagger s^\dagger - b_1^\dagger \cdot b_1^\dagger - b_2^\dagger \cdot b_2^\dagger) (\text{h.c.}) \\ & + \xi_2 [(b_1^\dagger \cdot b_1^\dagger - b_2^\dagger \cdot b_2^\dagger) (\text{h.c.}) + 4(b_1^\dagger \cdot b_2^\dagger) (\text{h.c.})] \\ & + \kappa \vec{L} \cdot \vec{L} \end{aligned}$$

Classical limit: potential energy surface

Equilibrium shape: two perpendicular
Jacobi coordinates of equal length

Equilateral triangle:
Oblate top



Rotation-Vibration Spectrum

$$E = E_{\text{vib}} + E_{\text{rot}}$$

$$E_{\text{vib}} \approx \omega_1 \left(\nu_1 + \frac{1}{2} \right) + \omega_2 (\nu_2 + 1)$$

$$\omega_1 = 4NR^2\xi_1$$

$$\omega_2 = \frac{4NR^2}{1 + R^2}\xi_2$$

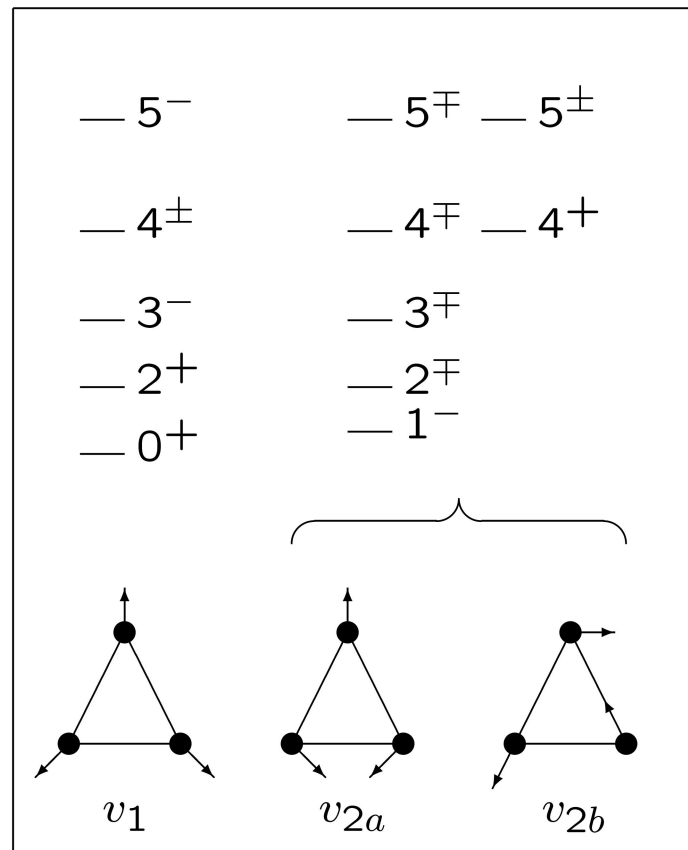
$$E_{\text{rot}} = \kappa L(L + 1)$$

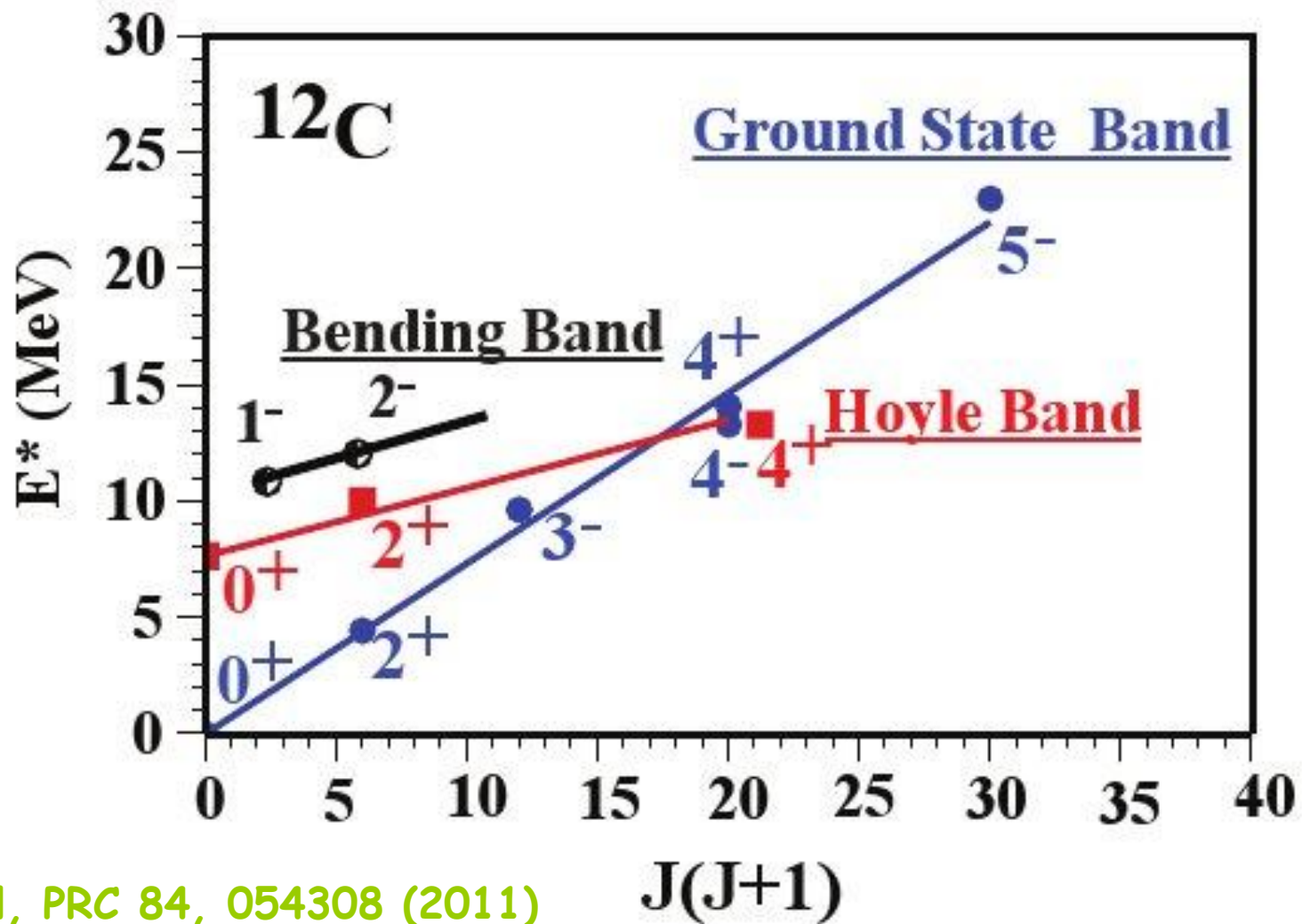
Rotational sequence

$$L^P = 0^+, 2^+, 3^-, 4^\pm, 5^-, \dots$$

Fingerprint of triangular symmetry (D_{3h})

Bijker & Iachello,
AP 298, 334 (2002)



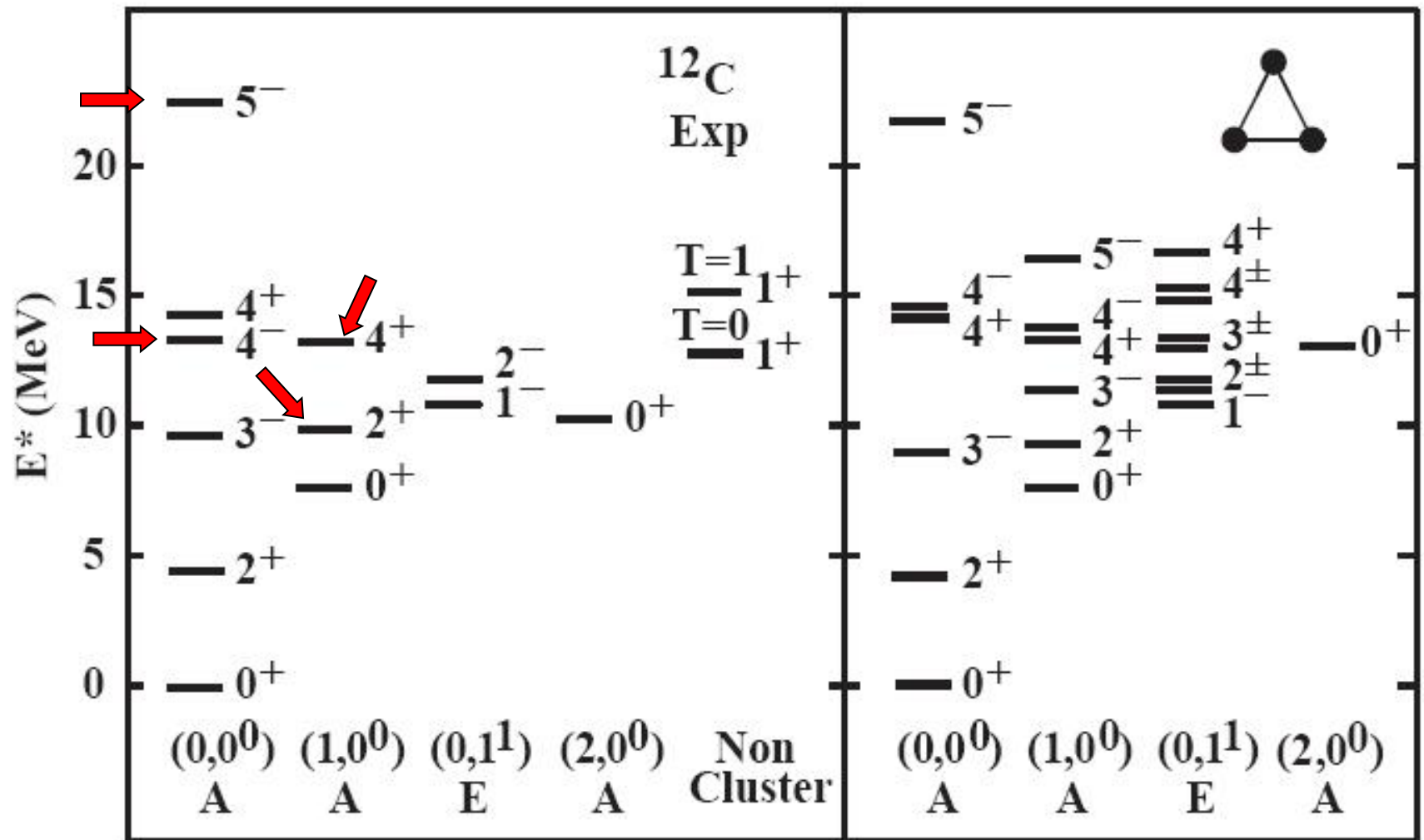


Itoh et al, PRC 84, 054308 (2011)

Freer et al, PRC 86, 034320 (2012)

Zimmerman et al, PRL 110, 152502 (2013)

Marín-Lámbarri et al,
PRL 113, 012502 (2014)



Marín-Lámbbarri, Bijker et al, PRL 113, 012502 (2014)

Electric Transitions

$$\rho(\vec{r}) = \frac{Ze}{k} \left(\frac{\alpha}{\pi}\right)^{3/2} \sum_{i=1}^k e^{-\alpha(\vec{r}-\vec{r}_i)^2}$$

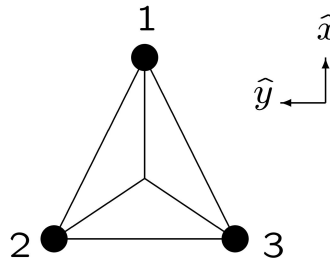
$$\vec{r}_1 = \left(\beta, \frac{\pi}{2}, 0\right)$$

$$\vec{r}_2 = \left(\beta, \frac{\pi}{2}, \frac{2\pi}{3}\right)$$

$$\vec{r}_3 = \left(\beta, \frac{\pi}{2}, \frac{4\pi}{3}\right)$$

$$\alpha = 0.56 \text{ 1/fm}^2$$

$$\beta = 1.82 \text{ fm}$$



$$\mathcal{F}_L(q) = c_L j_L(q\beta) e^{-q^2/4\alpha}$$

$$\langle r^2 \rangle^{1/2} = \sqrt{\beta^2 + 3/2\alpha}$$

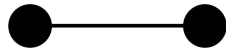
$$B(EL; 0^+ \rightarrow L^P) = \frac{(Ze)^2 c_L^2 \beta^{2L}}{4\pi}$$

$$c_L^2 = \frac{2L+1}{3} \left[1 + 2P_L(-\frac{1}{2})\right]$$

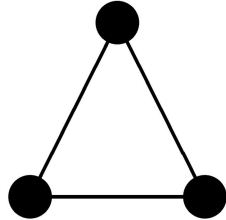
$$Q_{2+} = +\frac{2}{7}Ze\beta^2$$

		ACM	Exp.	
^{12}C	$B(E2; 2_1^+ \rightarrow 0_1^+)$	7.8	7.63 ± 0.19	$e^2\text{fm}^4$
	$B(E3; 3_1^- \rightarrow 0_1^+)$	65	104 ± 14	$e^2\text{fm}^6$
	$B(E4; 4_1^+ \rightarrow 0_1^+)$	48		$e^2\text{fm}^8$
	$Q_{2_1^+}$	5.7	5.97 ± 0.30	$e\text{fm}^2$
	$\langle r^2 \rangle^{1/2}$	2.448	2.468 ± 0.012	fm

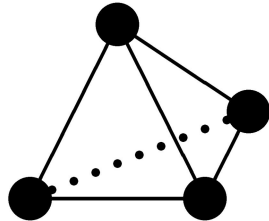
Bijker & Iachello,
AP 298, 334 (2002)



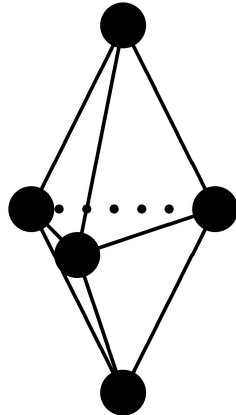
Z_2 Dumbbell ^8Be
NPA 973, 1 (2018)



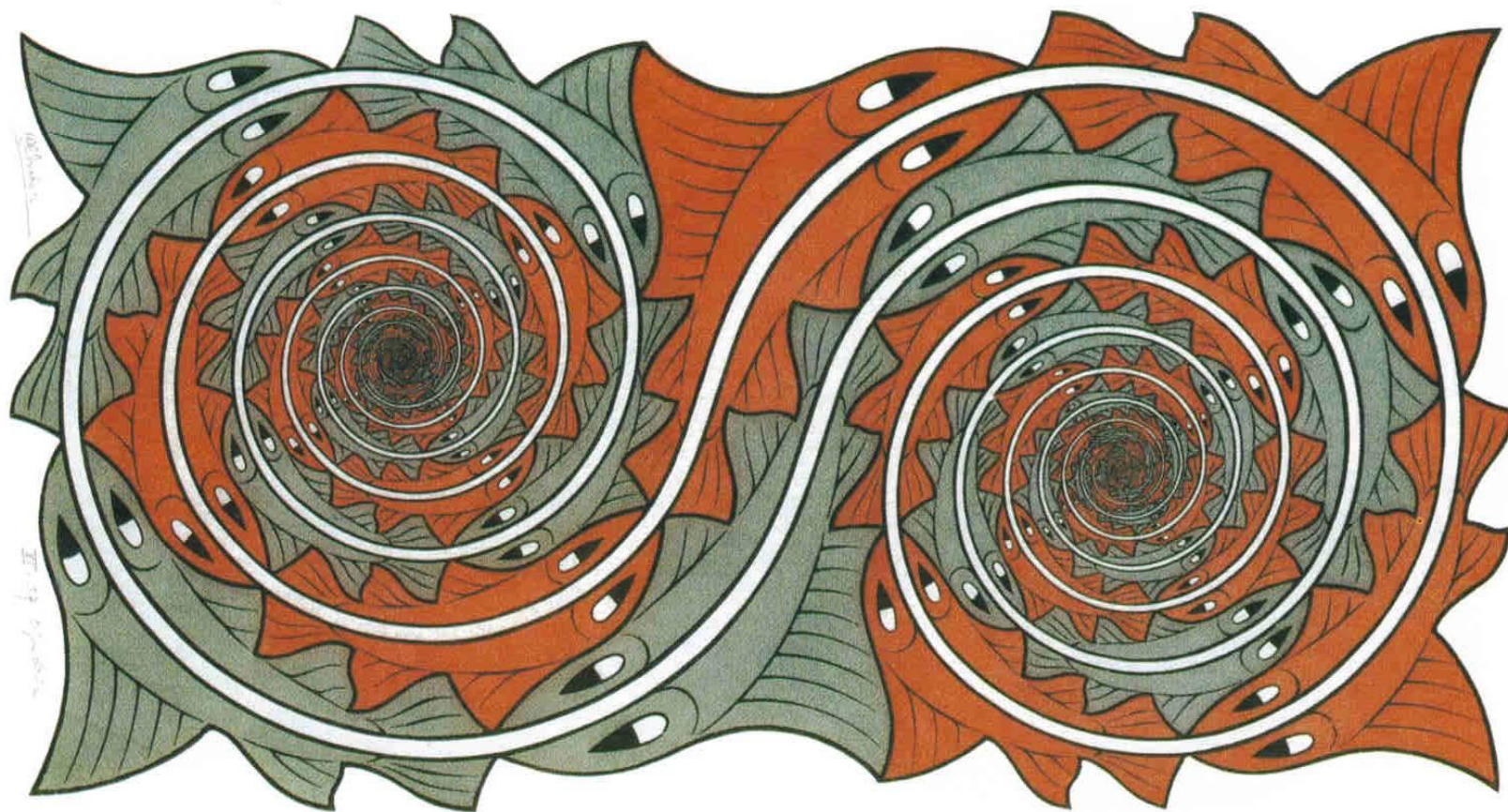
D_{3h} Triangle ^{12}C
PRC 61, 067305 (2000)
Ann Phys 298, 334 (2002)
PRL 113, 012502 (2014)



T_d Tetrahedron ^{16}O
PRL 112, 152501 (2014)
NPA 957, 154 (2017)

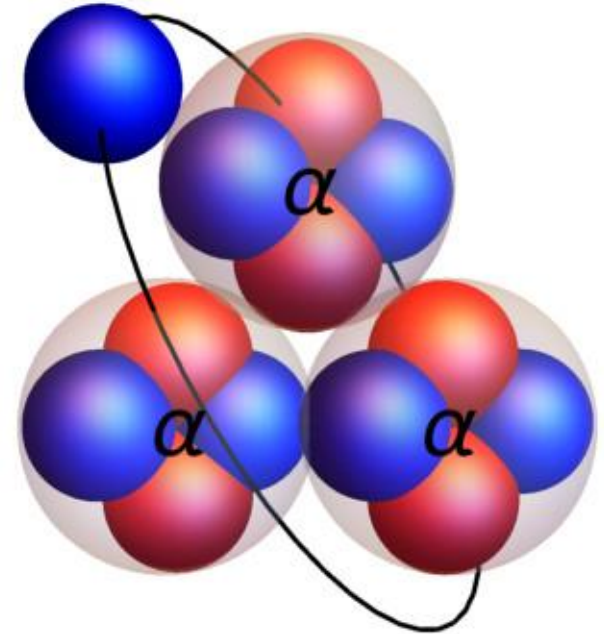


D_{3h} Bi-pyramid ^{20}Ne
NPA 1006, 122077 (2021)



Odd Cluster Nuclei

- What are the signatures of α -clustering in odd-mass nuclei?
- **Cluster Shell Model (CSM)**
- Splitting of sp levels in cluster potentials (analogous to Nilsson model)
- Double point groups



Cluster Shell Model

Cluster density

$$\begin{aligned}\rho(\vec{r}) &= \frac{Ze}{k} \left(\frac{\alpha}{\pi}\right)^{3/2} \sum_{i=1}^k \exp \left[-\alpha (\vec{r} - \vec{r}_i)^2 \right] \\ &= \frac{Ze}{k} \left(\frac{\alpha}{\pi}\right)^{3/2} e^{-\alpha(r^2 + \beta^2)} 4\pi \sum_{\lambda\mu} i_{\lambda}(2\alpha\beta r) Y_{\lambda\mu}(\theta, \phi) \sum_{i=1}^k Y_{\lambda\mu}^*(\theta_i, \phi_i)\end{aligned}$$

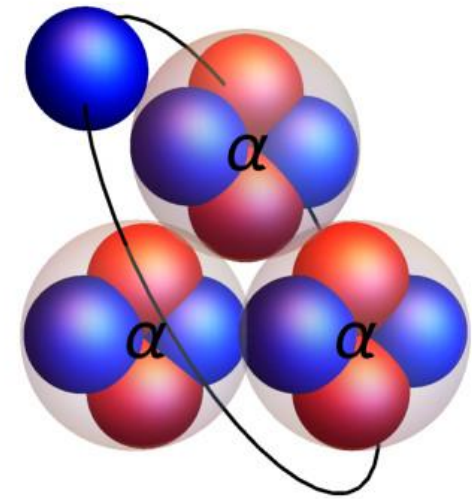
$$\vec{r}_i = (\beta, \theta_i, \phi_i)$$

Cluster potential

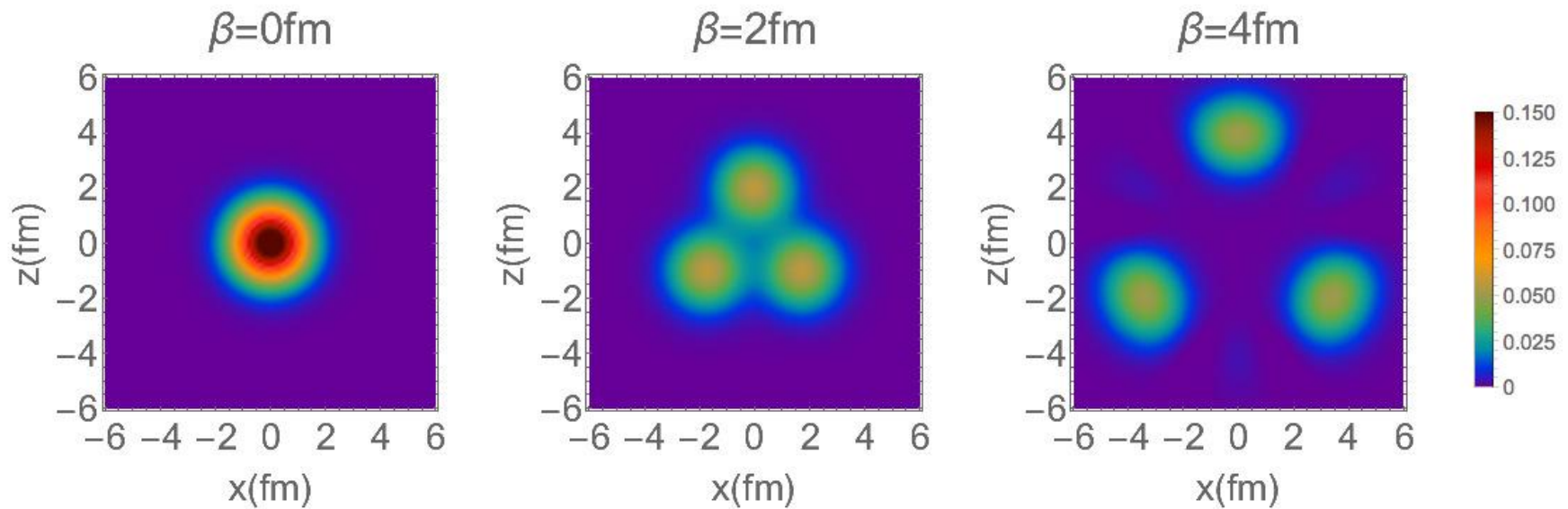
$$H = \sum_i \frac{\vec{p}_i^2}{2m} + V(\vec{r}) + V_{\text{so}}(\vec{r}) + V_{\text{C}}(\vec{r})$$

Della Rocca, Bijker & Iachello
NPA 966, 158 (2017)

Adrian Horacio Santana Valdés
M.Sc. Thesis, UNAM (2018)

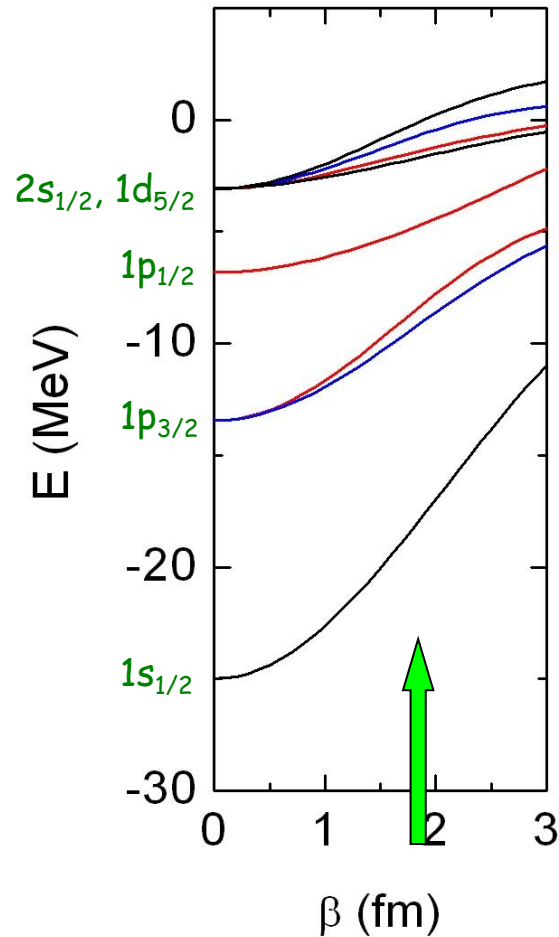


Densities of 3- α cluster



Della Rocca, Bijker & Iachello, NPA 966, 158 (2017)

Triangular Symmetry



Symmetry: D'_{3h}

Ω	$E_{1/2}$	$E_{5/2}$	$E_{3/2}$
Deg	2	2	2
$s_{1/2}$	1		
$p_{1/2}$		1	
$p_{3/2}$		1	1
$d_{3/2}$	1		1
$d_{5/2}$	1	1	1

Bijker & Iachello,
PRL 122, 162501 (2019)

Vibrations

^{12}C	$^{12}\text{C} \otimes \text{E}_{5/2}$	$^{12}\text{C} \otimes \text{E}_{1/2}$	$^{12}\text{C} \otimes \text{E}_{3/2}$
Bend $\checkmark \text{E}'$	$\text{---} \text{E}_{3/2}$ $\text{---} \text{E}_{1/2}$	$\text{---} \text{E}_{3/2}$ $\text{---} \text{E}_{5/2}$	$\text{---} \text{E}_{5/2}$ $\text{---} \text{E}_{1/2}$
Hoyle $\checkmark \text{A}'_1$	$\checkmark \text{E}_{5/2}$	$\text{---} \text{E}_{1/2}$	$\text{---} \text{E}_{3/2}$
Gsb $\checkmark \text{A}'_1$	$\checkmark \text{E}_{5/2}$	$\checkmark \text{E}_{1/2}$	$\text{---} \text{E}_{3/2}$

Rotations

gsb

$\text{gsb} \otimes E_{5/2}$

$\text{gsb} \otimes E_{1/2}$

$\text{gsb} \otimes E_{3/2}$

$D_{3h} : A'_1$	$D'_{3h} : E_{5/2}$	$D'_{3h} : E_{1/2}$	$D'_{3h} : E_{3/2}$
— 4^+ — 4^-	— $9/2^-$ — $9/2^+$ — $9/2^+$	— $9/2^+$ — $9/2^-$ — $9/2^-$	— $9/2^\pm$ — $9/2^\pm$
— 3^-	— $7/2^-$ — $7/2^+$ — $7/2^+$	— $7/2^+$ — $7/2^-$ — $7/2^-$	— $7/2^\pm$
— 2^+	— $5/2^-$ — $5/2^+$	— $5/2^+$ — $5/2^-$	— $5/2^\pm$
— 0^+	— $3/2^-$	— $3/2^+$	— $3/2^\pm$
— $1/2^-$	— $1/2^-$	— $1/2^+$	
$K^P = 0^+ \quad 3^-$	$K^P = 1/2^- \quad 5/2^+ \quad 7/2^+$	$K^P = 1/2^+ \quad 5/2^- \quad 7/2^-$	$K^P = 3/2^\pm \quad 9/2^\pm$

^{12}C

^{13}C

Rotational Energy

$$H_{\text{rot}} = \frac{L_1^2 + L_2^2}{2\mathcal{I}} + \frac{L_3^2}{2\mathcal{I}_3} = \sum_{i=1}^2 \frac{(J_i - j_i)^2}{2\mathcal{I}} + \frac{(J_3 - j_3)^2}{2\mathcal{I}_3}$$

Wave function

$$|\Omega, \mu; J^P K M\rangle = \frac{1}{\sqrt{2}} \left(1 + \hat{P} e^{i\pi J_2} \hat{p} e^{-i\pi j_2} \right) |J^P K M\rangle |\Omega, \mu\rangle$$

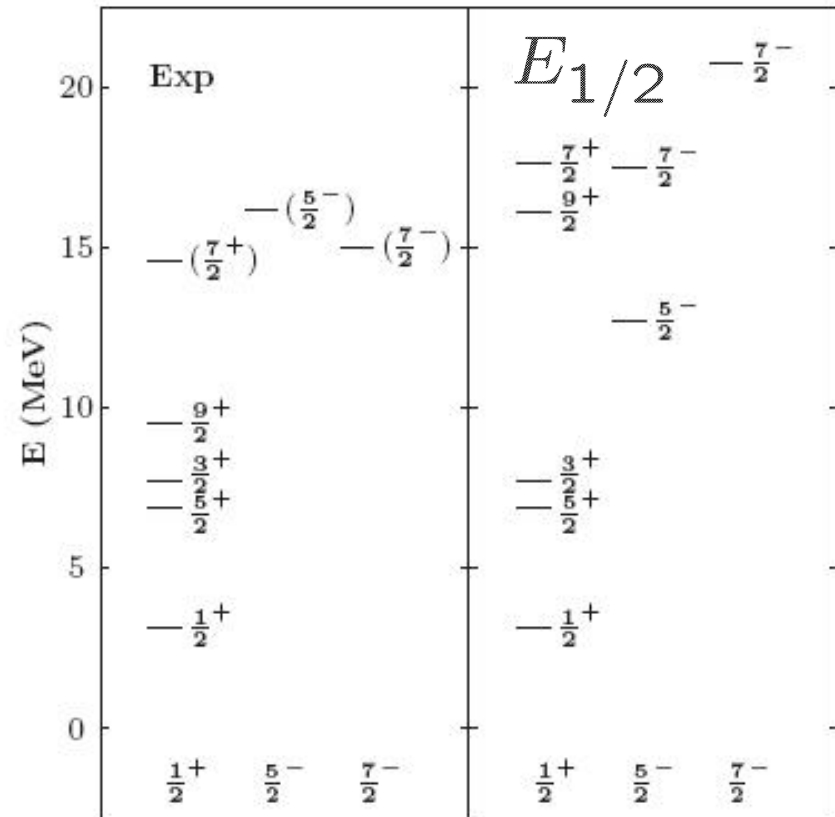
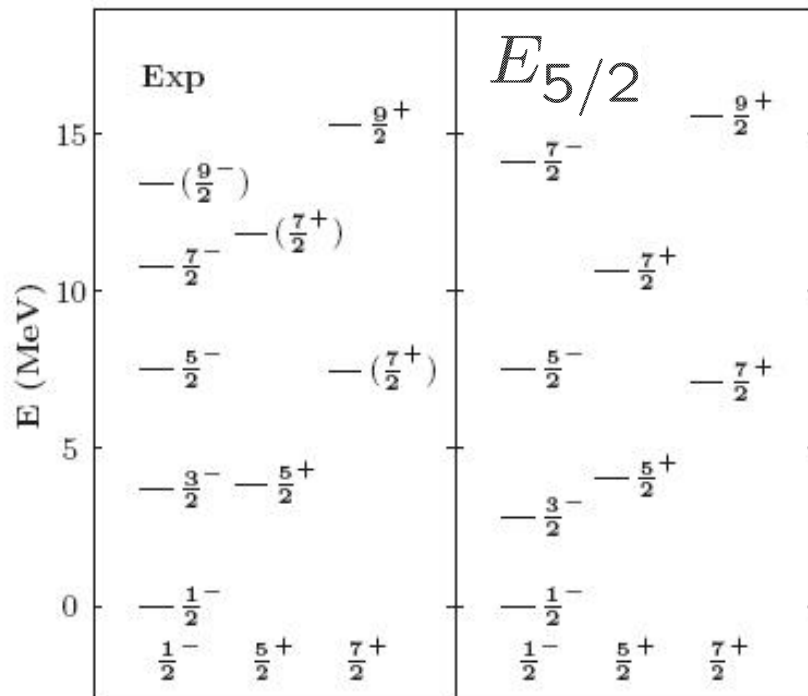
Energies

$$E_{\Omega}(J) \approx \frac{1}{2\mathcal{I}} \left[J(J+1) - 2K^2 + \delta_{K,1/2} a_{\Omega} (-1)^{J+1/2} \left(J + \frac{1}{2} \right) \right]$$

a_{Ω} Decoupling parameter

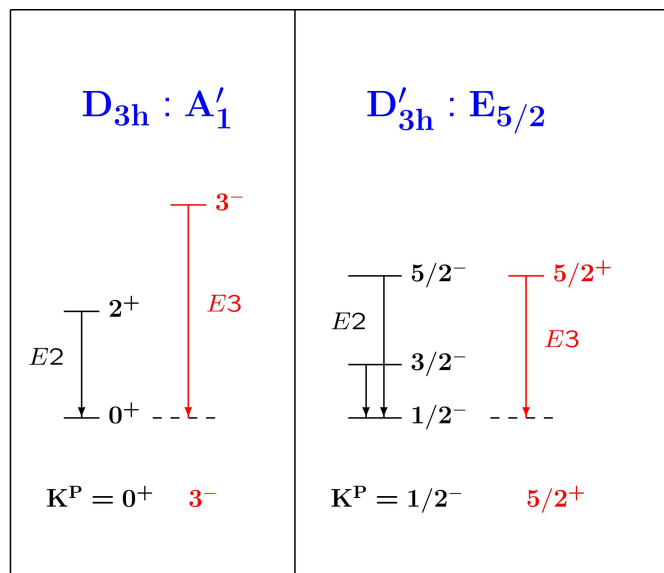
Generalized formula for triangular symmetry

Bandas Rotacionales en ^{13}C



Electric Transitions

$$\frac{B(E3; 3_1^- \rightarrow 0_1^+)}{B(E2; 2_1^+ \rightarrow 0_1^+)} = \frac{5}{2}\beta^2$$



Dominated by collective part

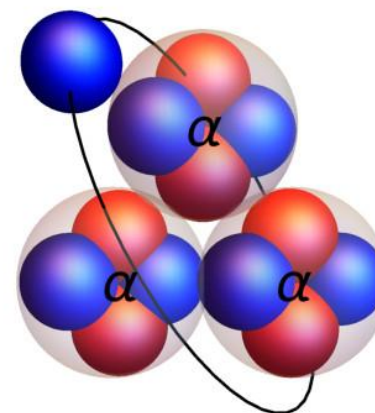
$$\begin{aligned}
 B(E2; 3/2_1^- \rightarrow 1/2_1^-) &= B(E2; 5/2_1^- \rightarrow 1/2_1^-) \\
 &= B(E2; 2_1^+ \rightarrow 0_1^+) \\
 B(E3; 5/2_1^+ \rightarrow 1/2_1^-) &= B(E3; 3_1^- \rightarrow 0_1^+) \\
 Q_{5/2_1^-} &= \frac{10}{7} Q_{3/2_1^-} = Q_{2_1^+}
 \end{aligned}$$

	$B(EL)$	Th	Exp	
^{12}C	$B(E2; 2_1^+ \rightarrow 0_1^+)$	7.8	7.63 ± 0.19	$e^2\text{fm}^4$
	$B(E3; 3_1^- \rightarrow 0_1^+)$	65.0	104 ± 14	$e^2\text{fm}^6$
	$Q_{2_1^+}$	5.7	5.97 ± 0.30	efm^2
^{13}C	$B(E2; 3/2_1^- \rightarrow 1/2_1^-)$	7.8	6.4 ± 1.5	$e^2\text{fm}^4$
	$B(E2; 5/2_1^- \rightarrow 1/2_1^-)$	7.8	5.6 ± 0.4	$e^2\text{fm}^4$
	$B(E3; 5/2_1^+ \rightarrow 1/2_1^-)$	65.0	100 ± 40	$e^2\text{fm}^6$
	$Q_{5/2_1^-}$	5.7		efm^2
	$Q_{3/2_1^-}$	4.0		efm^2

Form Factors

Charge distribution

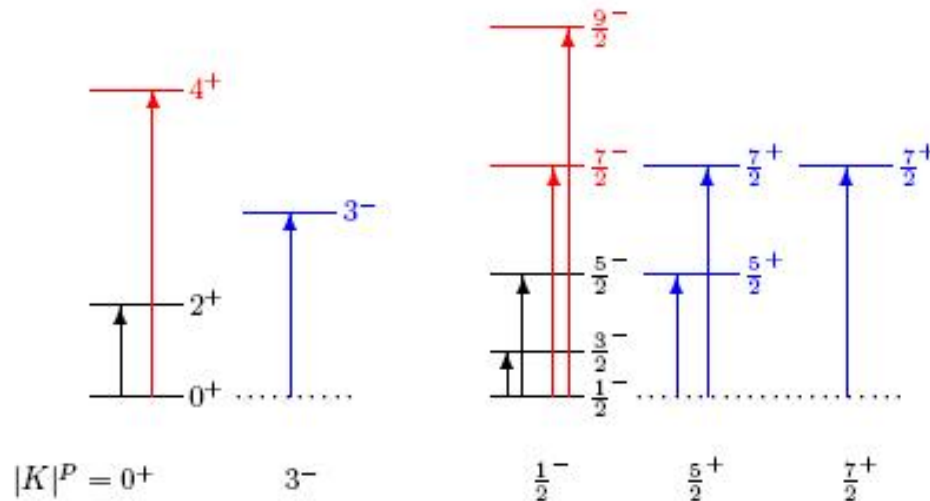
$$\begin{aligned}\rho_{\text{ch}}(\vec{r}) &= \rho_{\text{ch}}^{\text{c}}(\vec{r}) + \rho_{\text{ch}}^{\text{sp}}(\vec{r}) \\ \rho_{\text{ch}}^{\text{c}}(\vec{r}) &= \frac{(Ze)_{\text{c}}}{3} \left(\frac{\alpha}{\pi}\right)^{3/2} \sum_{i=1}^3 e^{-\alpha(\vec{r}-\vec{r}_i)^2} \\ \rho_{\text{ch}}^{\text{sp}}(\vec{r}) &= \tilde{e} \delta(\vec{r} - \vec{r}_{\text{sp}})\end{aligned}$$



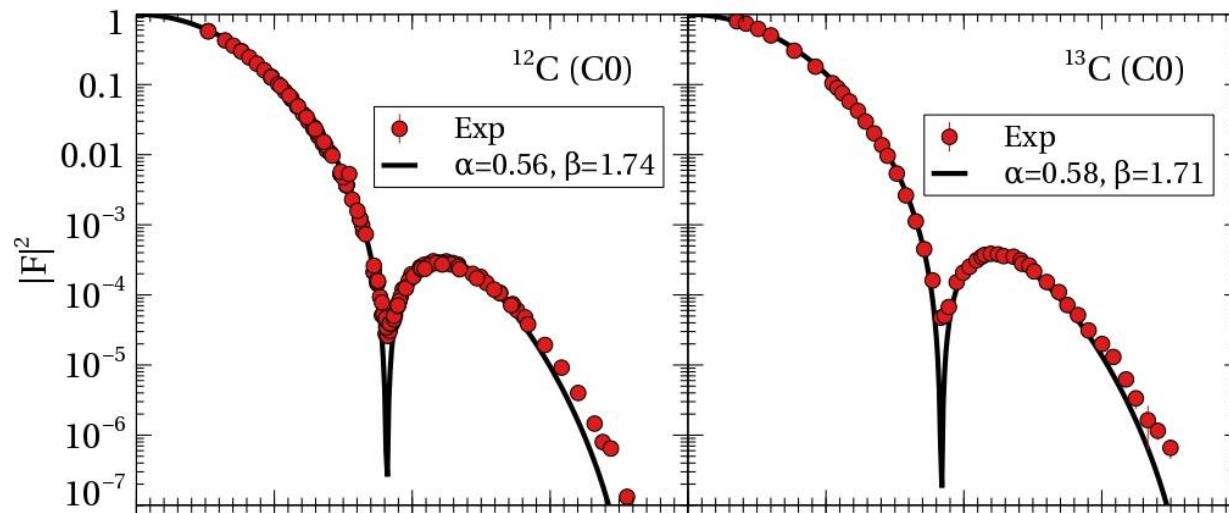
Form factors

$$\left\langle \psi_f \left| \int \rho_{\text{ch}}(\vec{r}) e^{i\vec{q}\cdot\vec{r}} d^3r \right| \psi_i \right\rangle$$

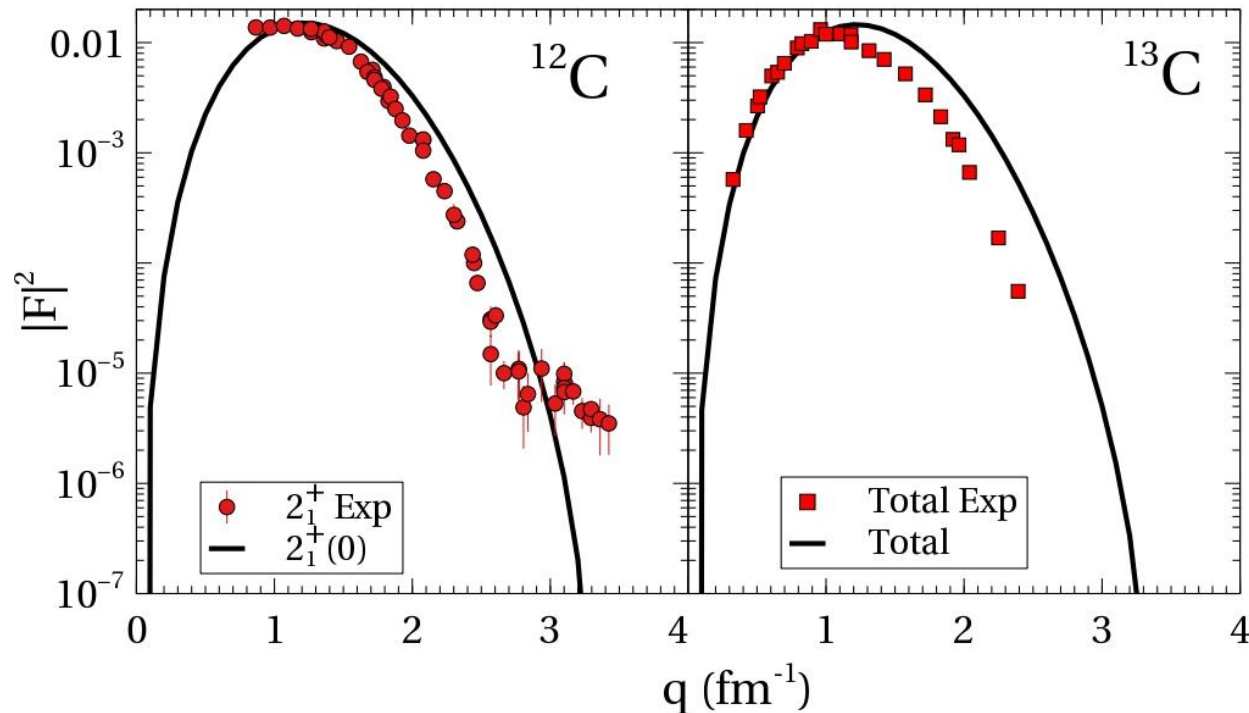
Correspondence ^{12}C y ^{13}C



$$\begin{aligned} & \left| F_{C\lambda}(^{12}\text{C}, q; 0^+, 0 \rightarrow L^P, K) \right|^2 \\ &= \sum_{I, K'} \left| F_{C\lambda}(^{13}\text{C}, q; \frac{1}{2}^-, \frac{1}{2} \rightarrow I^{P'}, K') \right|_C^2 \end{aligned}$$

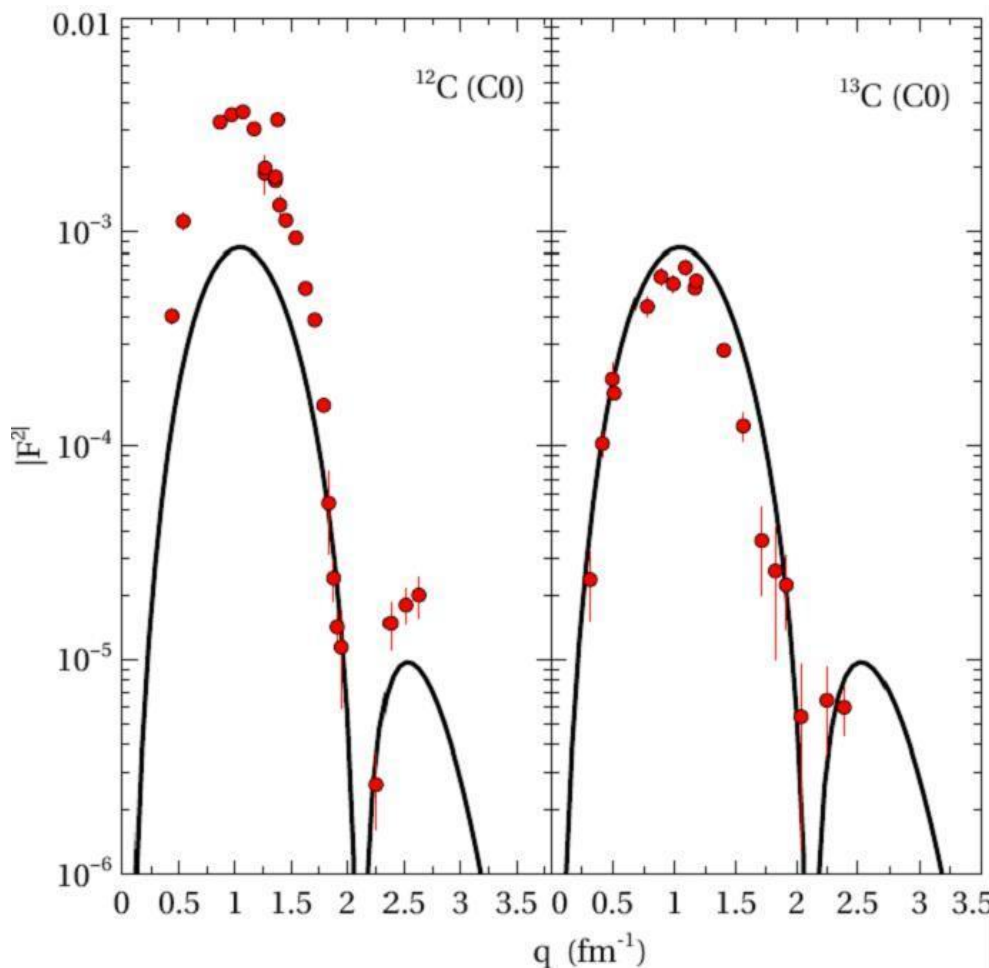


Factor de forma
elástico C0



Factor de forma
inelástico C2

Analogue of Hoyle state in ^{13}C



$$^{12}\text{C} : F(0_1^+ \rightarrow 0_2^+)$$

$$0_2^+ @ 7.65 \text{ MeV}$$

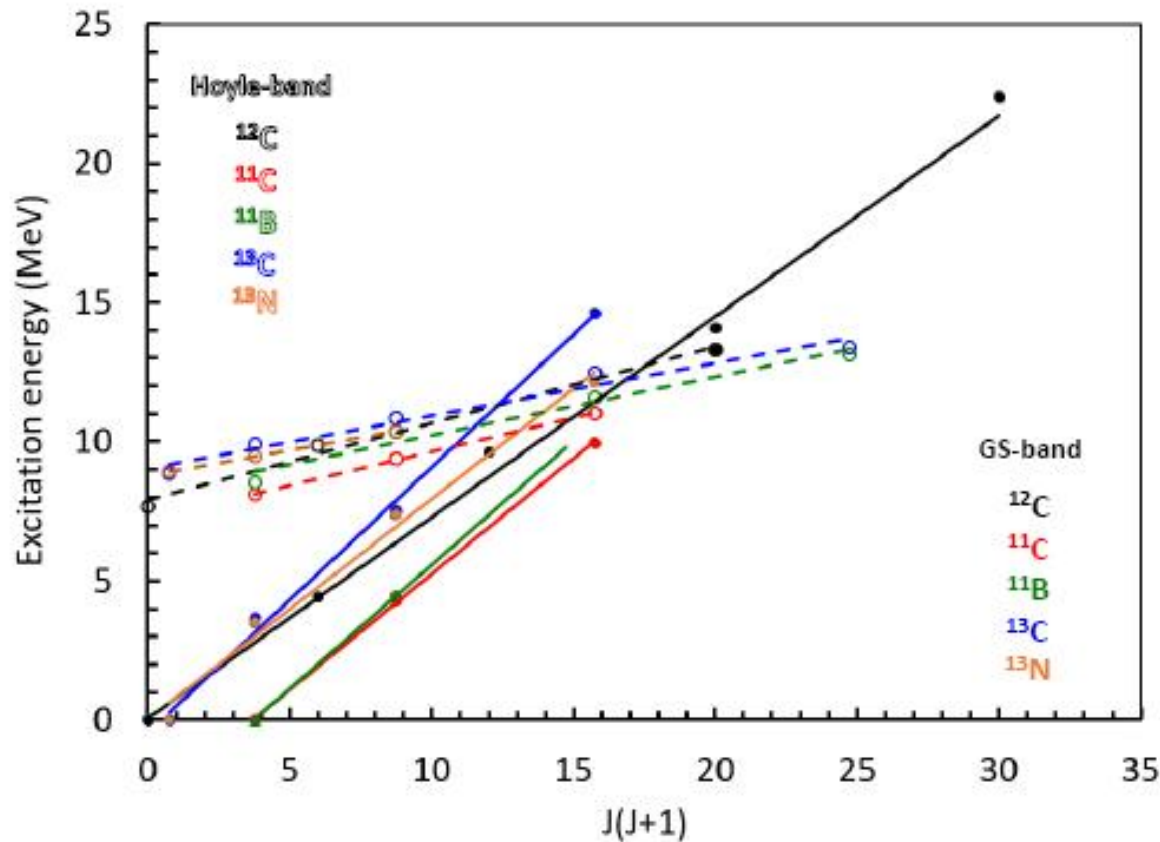
$$^{13}\text{C} : F(1/2_1^- \rightarrow 1/2_2^-)$$

$$1/2_2^- @ 8.86 \text{ MeV}$$

Santana & Bijker,
Phys. Lett. B 843,
138026 (2023)

Energy Systematics

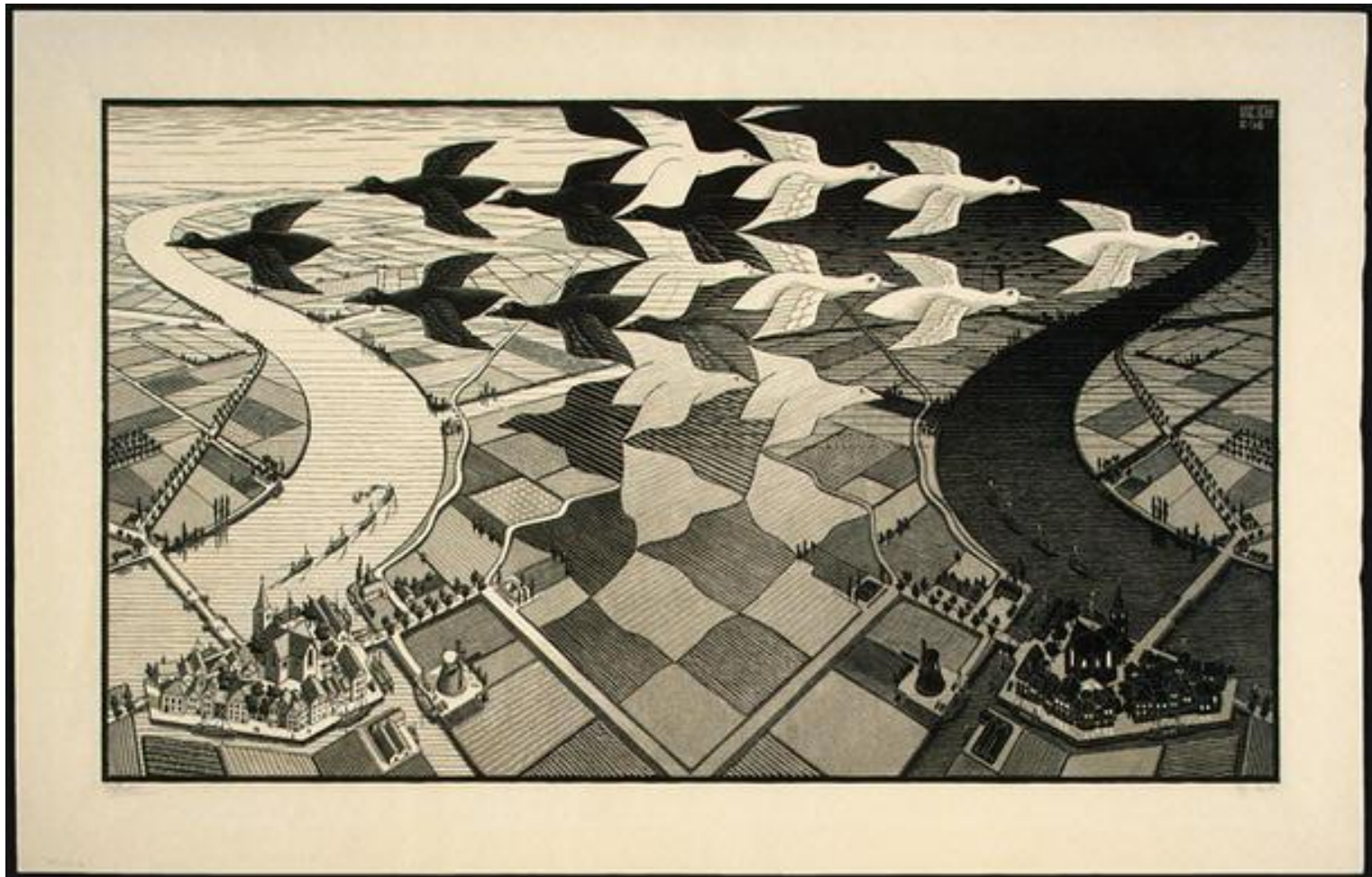
GS bands versus Hoyle-bands for 3α -like nuclei



Courtesy
Ivano Lombardo, 2019

Coincides in the
identification of the
analogue of the Hoyle
state in ^{13}C

$4n + 1$	Nuclei	References
$n = 2$	${}^9\text{Be}, {}^9\text{B}$	NPA 973, 1 (2018)
$n = 3$	${}^{13}\text{C}$	PRL 122, 162501 (2019)
		EPJ-ST 229, 2353 (2020)
		PLB 843, 138026 (2023)
	${}^{13}\text{N}$	B.Sc. Thesis, Villavicencio (2025)
$n = 5$	${}^{21}\text{Ne}, {}^{21}\text{Na}$	NPA 1010, 122193 (2021)
$4n + 2$	Nuclei	References
$n = 2$	${}^{10}\text{Be}$	JPCS 2619, 012006 (2023)
	${}^{10}\text{B}, {}^{10}\text{C}$	in progress, Omar Díaz



Summary and Conclusions I

- Algebraic Cluster Model
- SGA of $U(3n-2)$ for n -body systems
- Discrete and continuous symmetries
- Rotational bands fingerprints of point group symmetries
- Hoyle band: linear, bent or triangular?
- Applications in molecular, nuclear, hadronic physics

Summary and Conclusions II

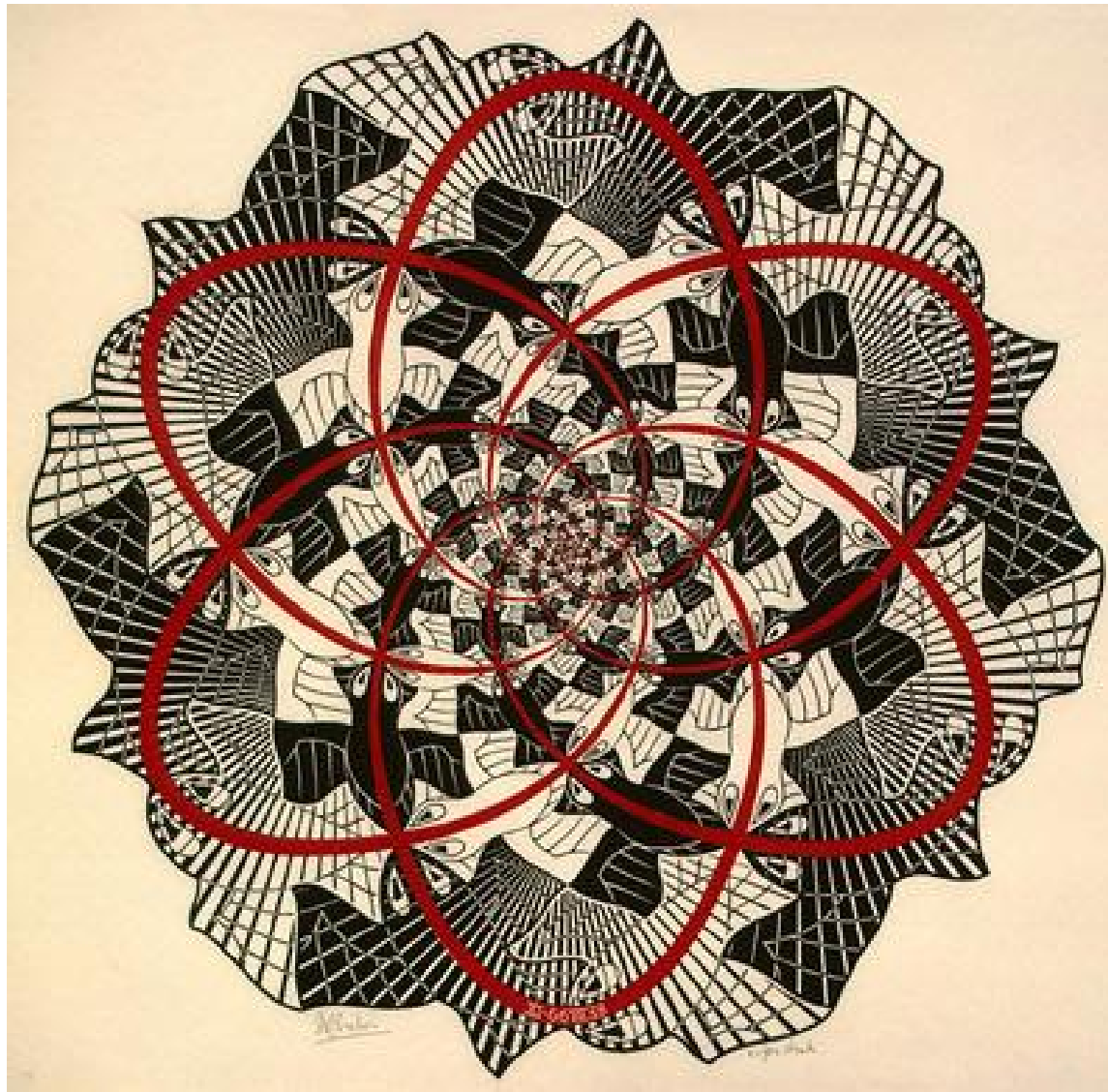
- Cluster Shell Model: ^{13}C
- Symmetries
- Rotational bands: fingerprints of a triangular configuration of three alpha particles plus a neutron
- Large electric transitions
- Form factor to identify the analogue of Hoyle state in ^{13}C
- Benchmark for microscopic studies of nuclear clustering



Bijker & Iachello,
PRL 122, 162501 (2019)

Who's Who?

- UNAM, Mexico
 - Adrian Santana
 - Omar Díaz
- UABJ, Mexico
 - Emiliano Villavicencio
- Yale
 - Valeria Della Rocca
 - Francesco Iachello



Roelof Bijker, ICN-UNAM

Pauli Principle

- ACM: effective α - α interaction of the Morse type
- CSM similar to Nilsson model
- Nucleon moves in the deformed field generated by the cluster of α particles

