

# The high spin level scheme of $^{156}\text{Dy}$ and neighbouring nuclei.

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Phys. Rev. C110, 034313 (2024))

# Outline

- ▶ The CNS, CNSB (and CNS(B)) models.
- ▶ Interpretation of the high-spin  $(\pi, \alpha) = (+, 0)$  bands in  $^{156}\text{Dy}$ .
- ▶ Coexistence at low spin for the  $(\pi, \alpha) = (+, 0)$  bands in  $^{156}\text{Dy}$  and neighbouring nuclei.
- ▶ The  $\pi(h_{11/2})$  crossing in  $^{156}\text{Dy}$  and neighbouring nuclei
- ▶ Summary and conclusions.

## Cranked Nilsson-Strutinsky (CNS) -Bogoliubov (CNSB) models.

Modified oscillator or Nilsson potential:

$$V_{MO} = V(\varepsilon_2, \gamma, \varepsilon_4) - \kappa \hbar \omega_0 \left( 2\vec{I}_t \cdot \vec{s} + \mu (I_t^2 - \langle I_t^2 \rangle_N) \right).$$

The cranking Hamiltonian is diagonalized (CNS):

$$H_{CNS}^\omega = H_{MO} - \omega j_x; \quad H_{CNS}^\omega \chi_\nu^\omega = e_\nu^\omega \chi_\nu^\omega; \quad e_\nu = \langle \chi_\nu^\omega | H_{CNS} | \chi_\nu^\omega \rangle$$

CNSB:  $H_{CNSB}^\omega = H_{MO} - \omega j_x - \Delta(P^\dagger + P) - \lambda \hat{N}$

Total quantities:

$$E_{tot} = E_{rld} + E_{shell} (+E_{pair}); \quad E_{shell} = \sum_{occ} e_\nu - \left\langle \sum_{occ} e_\nu \right\rangle \quad I = \sum_{occ} \langle j_x \rangle$$

Rotating liquid drop energy:

$$E_{rld}(Z, N, I, \varepsilon_i) = E_{ld}(Z, N, \varepsilon_i) + \frac{\hbar^2 I(I+1)}{2\mathcal{J}_{rig.}(Z, N, \varepsilon_i)}$$

$E_{ld}$ : Lublin-Strasbourg drop (LSD model) [or FRLDM]

$\mathcal{J}_{rig.}$ : diffuse surface:  $r_0 = 1.16$  fm,  $a = 0.6$  fm

T. Bengtsson and I.R., NPA 436, 14 (1986); B.G. Carlsson and I.R., PRC 74, 011302 (2006); B.G. Carlsson *et al.* PRC 78, 034316 (2008), Hai-Liang Ma, B.G. Carlsson, I.R. and H. Ryde, PRC 90, 014316 (2014).

## Cranked Nilsson-Strutinsky (CNS) -Bogoliubov (CNSB) models.

Removal of virtual crossings - diabatic configurations.

CNS: Mesh in deformation space:  $(\varepsilon_2, \gamma, \varepsilon_4)$  .

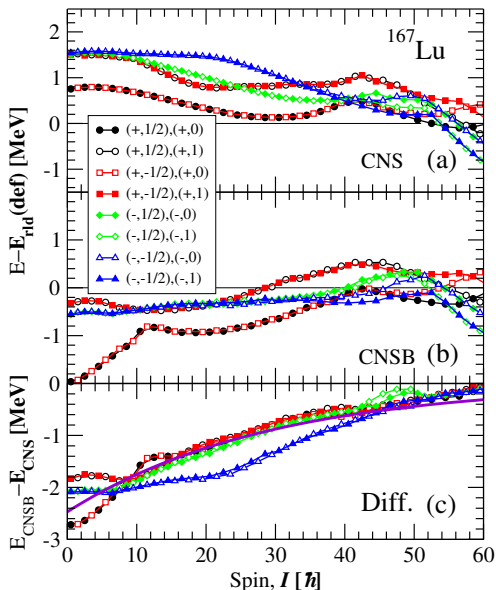
Unique features to label orbitals and thus to fix conf's, e.g.:

- ▶ Diagonalization in rotating harmonic oscillator basis,  $N_{rot}$ -shells treated as pure
- ▶ Labelling of high- $j$  shells and low- $j$  shells, respectively

CNSB: Mesh in def. and pairing space:  $(\varepsilon_2, \gamma, \varepsilon_4, \lambda_p, \lambda_n, \Delta_p, \Delta_n)$

- ▶ Minimization in  $(\varepsilon_2, \gamma, \varepsilon_4, \lambda_p, \lambda_n, \Delta_p, \Delta_n)$  space
- ▶ Particle number projection.
- ▶ Removal of virtual interactions for fixed parameters  
⇒ Possible to draw smooth PES's in the full  $(\varepsilon_2, \gamma)$ -plane.  
Conf's:  $(\pi, \alpha)_p(\pi, \alpha)_n$

## Comparison CNS) ↔ CNSB



Analyze rotational bands:  
 $\mathcal{J}^{(1)}$ ,  $\mathcal{J}^{(2)}$  or  $I$  vs.  $\omega$

More illustrative:  
 $E - E_{rld}(def)$

Comparison between CNS  
 and CNSB for  $^{167}\text{Lu}$ .

Conclusion:

Add average pairing to

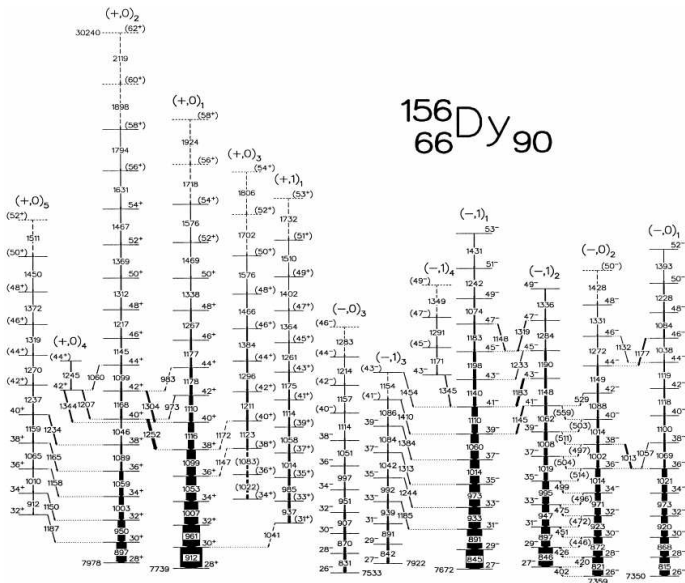
CNS:

$$\langle E_{pair} \rangle = -2.47 \exp(-I/29)$$

CNS(B)

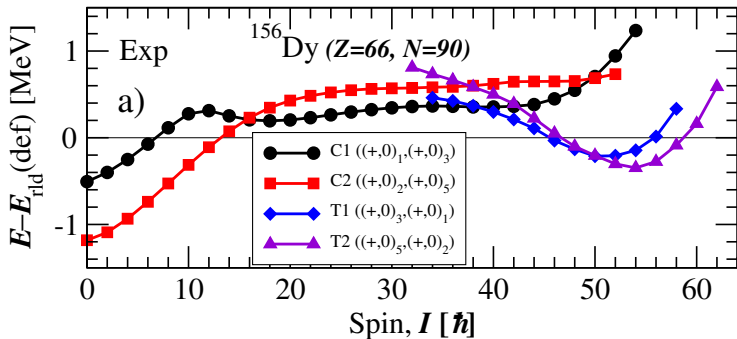
# High spin level scheme; $^{156}\text{Dy}$

$^{156}_{66}\text{Dy}_{90}$



F.G. Kondev, M.A. Riley, R.V.F. Janssens, J. Simpson, A.V. Afanasjev, I.R. et al., PLB 437, 35 (1998).

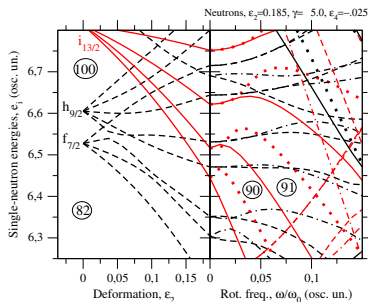
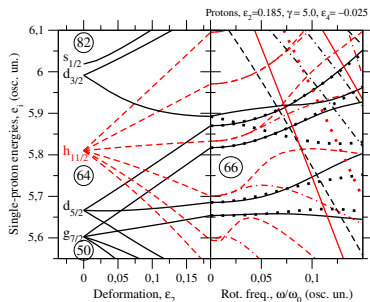
$E - E_{\text{rid}}(\text{def}); (+, 0)$  bands,  $^{156}\text{Dy}$ .



Collective bands: C1, C2

Terminating bands: T1, T2

## Nilsson diagrams; single particle routhians.



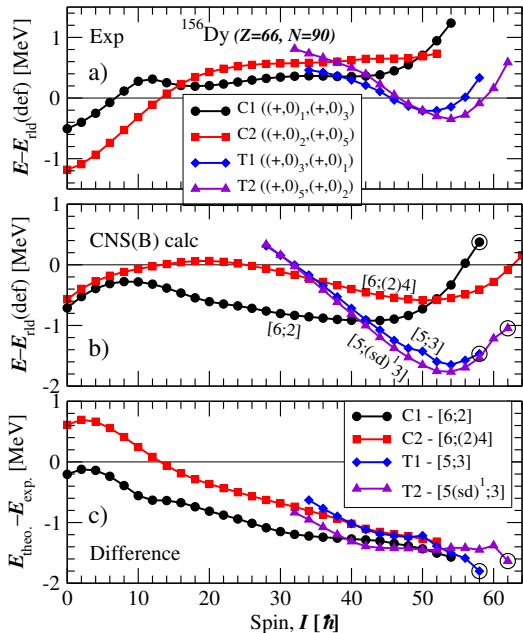
Configurations,  $^{156}\text{Dy}$ : [# of  $h_{11/2}$  protons; # of  $i_{13/2}$  neutrons]

Low-energy configurations at low spin: [6;2], [6;(2)4]

at high spin [5;3]

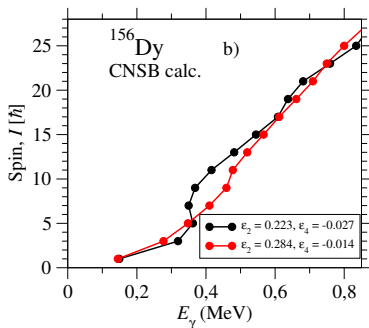
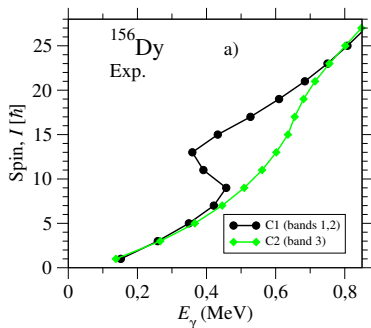


Comparison, exp  $\leftrightarrow$  calc; (+, 0) bands,  $^{156}\text{Dy}$ .



Convincing agreement between exp. and calc. Note that  $I = 58$  is the highest spin you can build in  $[5;3]$  conf.  $+4\hbar$  with  $(dg)^{-1}sd^1$  excitation.

## $i_{13/2}$ alignment in $^{156}\text{Dy}$ (backbend or upbend)



Calculations at fixed deformations:

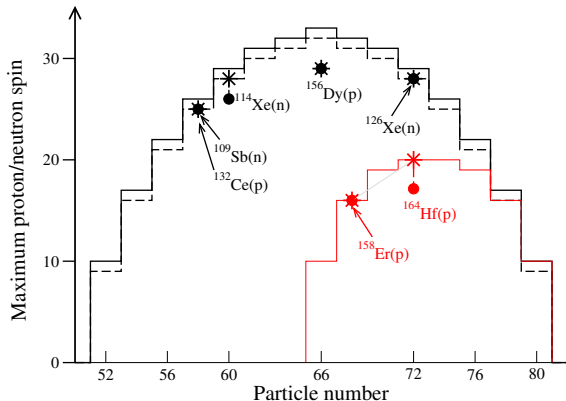
Minima of  $[6;2]$  and  $[6;(2)4]$  conf's in CNS calculations.

Alignment in the bottom of a high- $j$  shell:

F.S. Stephens, R.M. Diamond and S.G. Nilsson, Phys. Lett. 44B, 429.

S.G. Nilsson and I.R., Shapes and shells in Nuclear Structure, Cambridge Univ. Press, 1995, 2005.

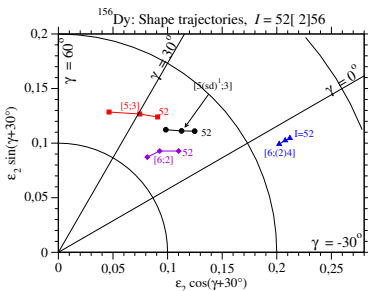
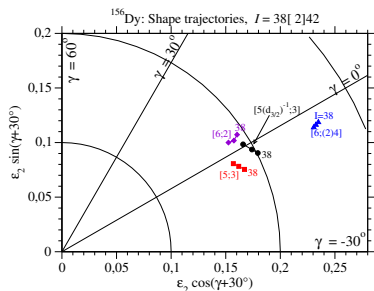
## Maximum spin in shells.



Maximum spin in a half-full shell. Thus  $Z = 66$  for  $^{156}\text{Dy}$  is in the middle of the  $Z/N = 50-82$  shell.

Max. value is  $33\hbar$ . In the highest spin band in  $^{156}\text{Dy}$ ,  $I_p = 29$  has been observed, i.e. only  $4\hbar$  below the maximum value where all 16 valence particles rotate in the same direction around the core.

# Shape coexistence, $^{156}\text{Dy}$ ; $I = 38, 40, 42, I = 52, 54, 56$



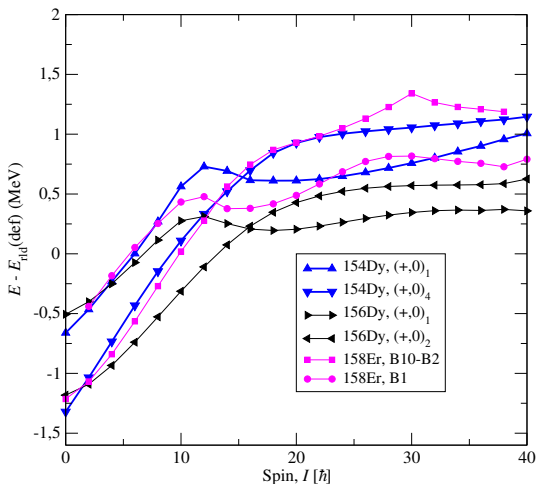
Valence space

configurations:  $\varepsilon_2 \approx 0.20$  for  $I \approx 40$

coming close to termination for  $I \approx 55$ .

$(h_{11/})^{-2}$  configuration: Stays at  $\varepsilon_2 > 0.20$  for spin values up to  $I \approx 60$ .

# Similar bands in $^{154,156}\text{Dy}$ , $^{158}\text{Er}$ .



Coexistence between:

$$\pi(h_{11/2})^6 \nu(i_{13/2})^2$$

and

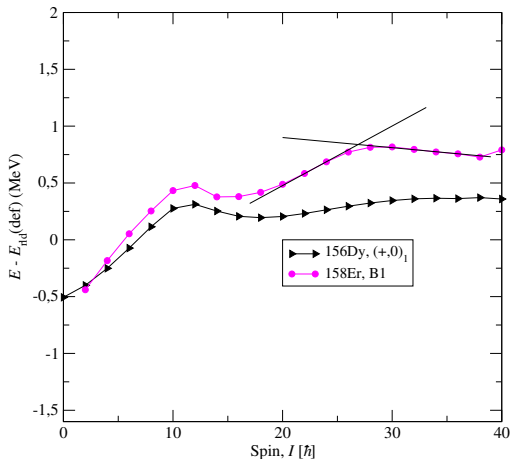
$$\pi(h_{11/2})^6 \nu(h_{11/2})^{-2}(i_{13/2})^4$$

configurations.

$^{158}\text{Er}$ : J. Simpson, M.A. Riley et al.,  
 PRC 107, 054305 (2023)

$^{154}\text{Dy}$ : W.C. Ma et al., PRC 65,  
 034312 (2002).

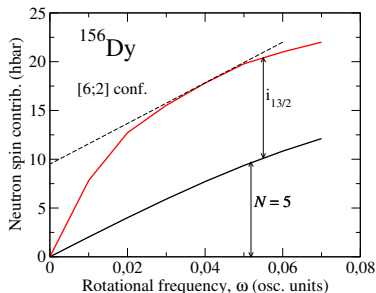
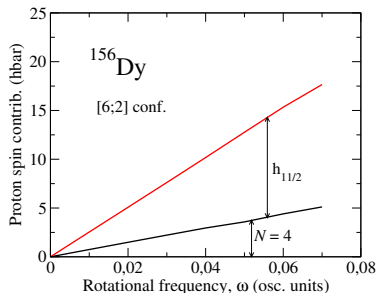
$^{156}\text{Dy}$ ,  $^{158}\text{Er}$ :  $\pi(h_{11/2})$  bandcrossing around  $I = 30$ ?



"Accepted" explanation:  
Alignment of a pair of  $h_{11/2}$  protons but interaction strength so large in  $^{156}\text{Dy}$  that the alignment is smeared out over an extended spin range.

Oscillating behaviour of the yrast-yrare interaction:  
R. Bengtsson, I. Hamamoto and B.R. Mottelson, PL 73B, 259 (1978)

## Building of spin in the [6;2] configuration of $^{156}\text{Dy}$ , CNS



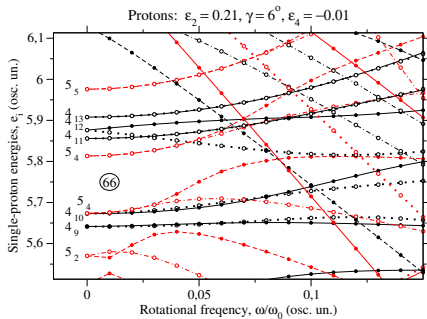
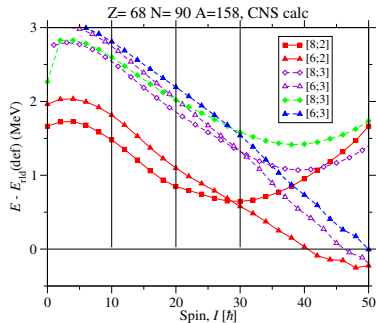
Large spin contribution from  $i_{13/2}$  already at small frequencies.

Linear increase with  $\omega$  from other  $j$ -shells.

$$N = 4: \pi(d_{5/2}, g_{7/2}); N = 5: \nu(f_{7/2}, h_{9/2});$$

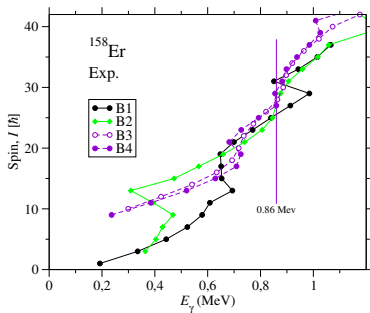
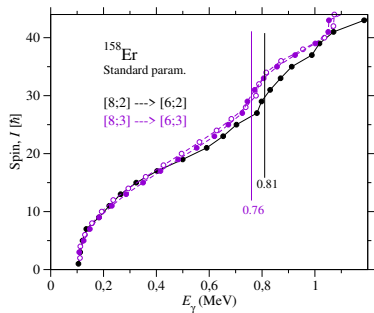
No special contribution from  $h_{11/2}$  protons!

# Crossing between $\pi(h_{11/2})^8$ and $\pi(h_{11/2})^8$ bands in $^{158}\text{Er}$ , CNS.





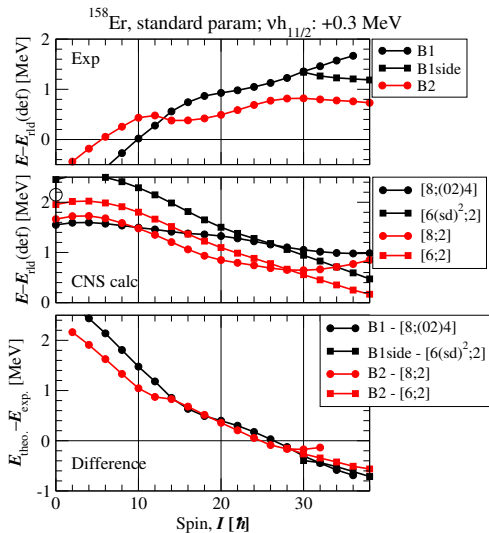
# The $\pi(h_{11/2})^8 \rightarrow \pi(h_{11/2})^6$ transition in $^{158}\text{Er}$



Calculation, crossing at  
 $E_\gamma \approx 0.78$  MeV  
 $\hbar\omega = 0.39$  MeV

Exp., crossing at  
 B2-B4:  $E_\gamma \approx 0.86$  MeV  
 B1: Crossing looks different  
 $\nu(h_{11/2})^{-2} \rightarrow \pi(h_{11/2})^6$

# Compare exp $\leftrightarrow$ calc., (+, 0) in $^{158}\text{Er}$ .



Differences well collected showing a pairing energy decreasing with spin.

Previous publications; # of protons in  $h_{11/2}$ .

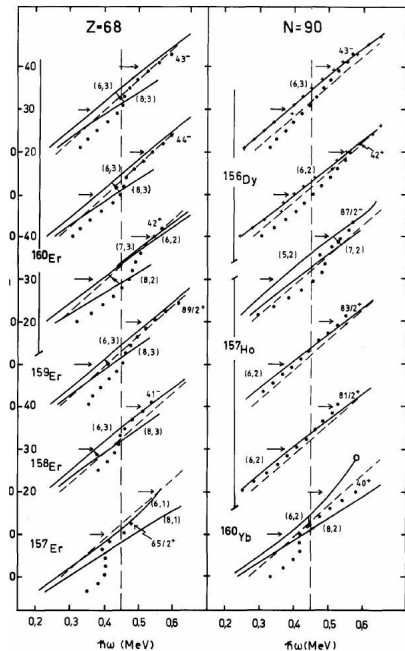
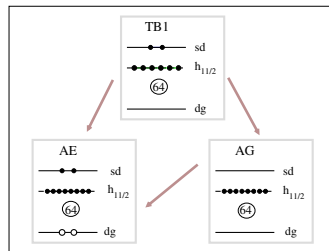
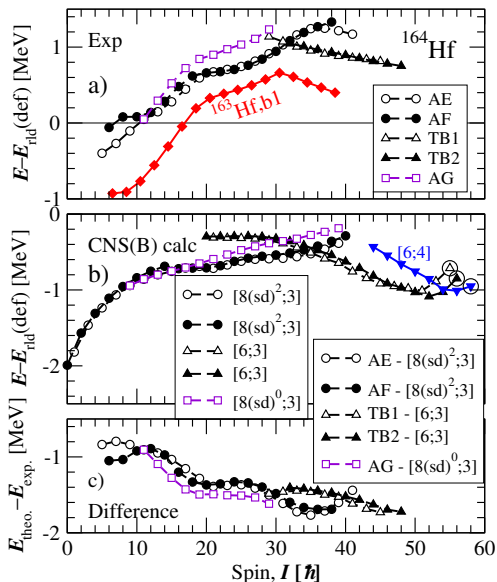


Figure from T. Bengtsson and I.R., Phys. Lett. **163B**, 31 (1985).

F. Grümmer, K.W. Schmid and A. Faessler, Nucl. Phys. A326, 1 (1979) :

It is shown that the oscillations of the yrast-yrare coupling can be explained as being due to the mixing of states with different particle numbers inside the valence shell.

$\pi(h_{11/2})^8 \rightarrow \pi(h_{11/2})^6$  transition in  $^{164}\text{Hf}$



Transition from AE configuration to TB1 configuration.

## Summary

- ▶ Detailed understanding of collective and terminating bands in  $^{156}\text{Dy}$  up to the highest spin,  $I \approx 60$
- ▶ The bands built on the two lowest  $0^+$  states in  $^{154,156}\text{Dy}$  and  $^{158}\text{Er}$  are assigned as collective coexisting bands for  $I \approx 10 - 40$  suggesting coexistence also for lower spin values down to  $I = 0$ .
- ▶ Assignments at higher spin are often more reliable because pairing is negligible. These assignments can be helpful for the understanding of lower spin states.
- ▶ The band crossings at  $I \approx 30$  in rare earth region can be understood in an unpaired formalism as caused by a change of the number of  $h_{11/2}$  protons at the crossing.
- ▶ In the nuclei analyzed by us, except for the first  $i_{13/2}$  alignment, the band-crossings can be described in an unpaired formalism (or adding an average pairing).

Thank you!