

# Configuration assignments for the extensive level scheme of $^{167}\text{Lu}$ .

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# Outline

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- The Cranked Nilsson-Strutinsky (CNS) and Cranked Nilsson-Strutinsky Bogoliubov (CNSB) formalisms.
- Typical features of high-spin bands,  $E - E_{rld}$  diagrams.
- The extensive level scheme  $^{167}\text{Lu}$ .
- Interpretation of the observed bands in  $^{167}\text{Lu}$  - evolution with spin, band crossings observed.
- Conclusions.

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

Modified oscillator or Nilsson Hamiltonian:

$$H_{MO} = -\frac{\hbar^2}{2M}\Delta + V_{HO}(\varepsilon_2, \gamma) + V(\varepsilon_4) - \kappa\hbar\omega_0 \left( 2\vec{l}_t \cdot \vec{s} + \mu (l_t^2 - \langle l_t^2 \rangle_N) \right).$$

$$\text{Cranking Hamiltonian: } H^\omega = \underbrace{H_{MO} - \omega j_x}_{\text{CNS}} - \overbrace{\Delta(P^\dagger + P) - \lambda\hat{N}}^{\text{CNSB}}$$

The cranking Hamiltonian is diagonalized

⇒ Single (quasi)-particle energies,  $e_i(E_i)$ , spin projections  $(j_x)_i$  etc.

Total quantities (macroscopic-microscopic formalism):

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

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.

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Removal of virtual crossings - diabatic configurations.

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

- Unique features to label orbitals ( $N_{osc}$ -shell, high- $j$  - low- $j$ , (pseudo-spin partners)) and thus to fix conf's.
- Minimization in  $(\varepsilon_2, \gamma, \varepsilon_4)$ -space (each configuration at each spin value).

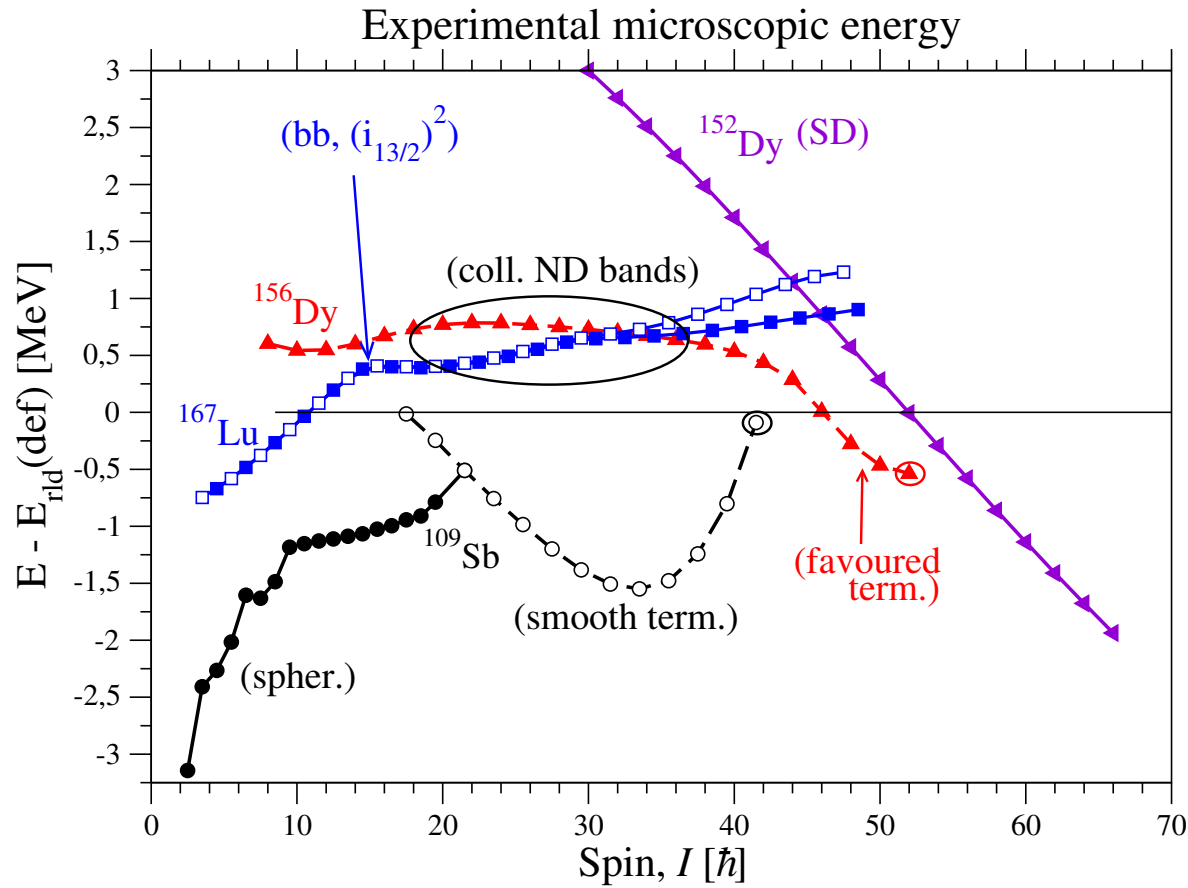
CNSB: Mesh in def. and pairing space:  $(\varepsilon_2, \gamma, \varepsilon_4, \lambda_p, \lambda_n, \Delta_p, \Delta_n)$

- 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:  $(\pi, \alpha)_p(\pi, \alpha)_n$

# Energy relative to rotating liquid drop energy.



Best way to get an understanding of nuclear level schemes,  $E - E_{rld}$  plots:

- Excitation energy.
- Moments of inertia  $\mathcal{J}^{(1)}$  (slope) and  $\mathcal{J}^{(2)}$  (curvature).
- Band crossings.
- Signature splitting.

# Comparison CNS - CNSB, $^{167}\text{Lu}$ .

No pairing and pairing included.

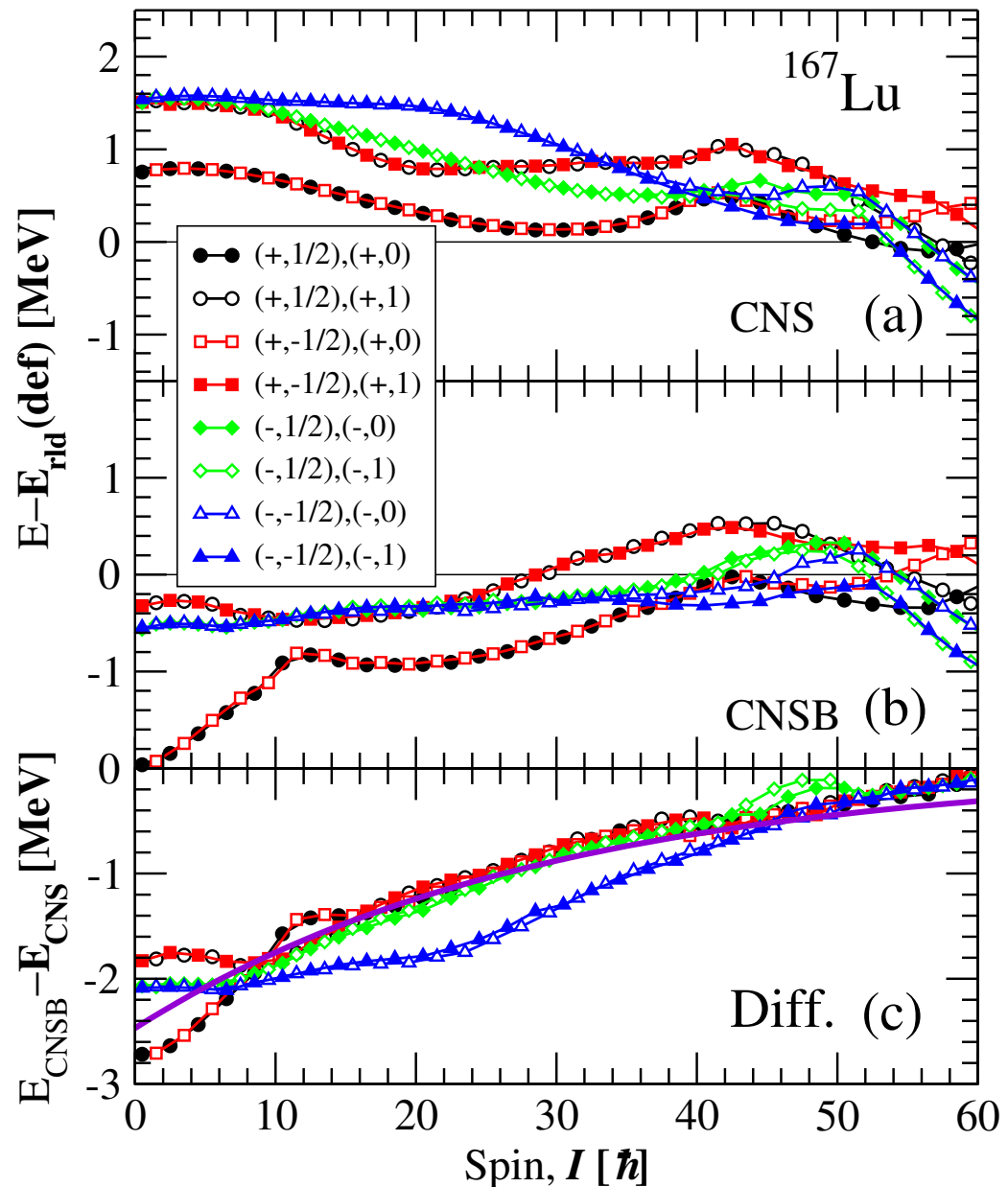
Difference shows a rather smooth trend approaching zero at high spin.

⇒ Experiment down to rather low spin is well described by CNS if average pairing is subtracted (moment of inertia renormalized):

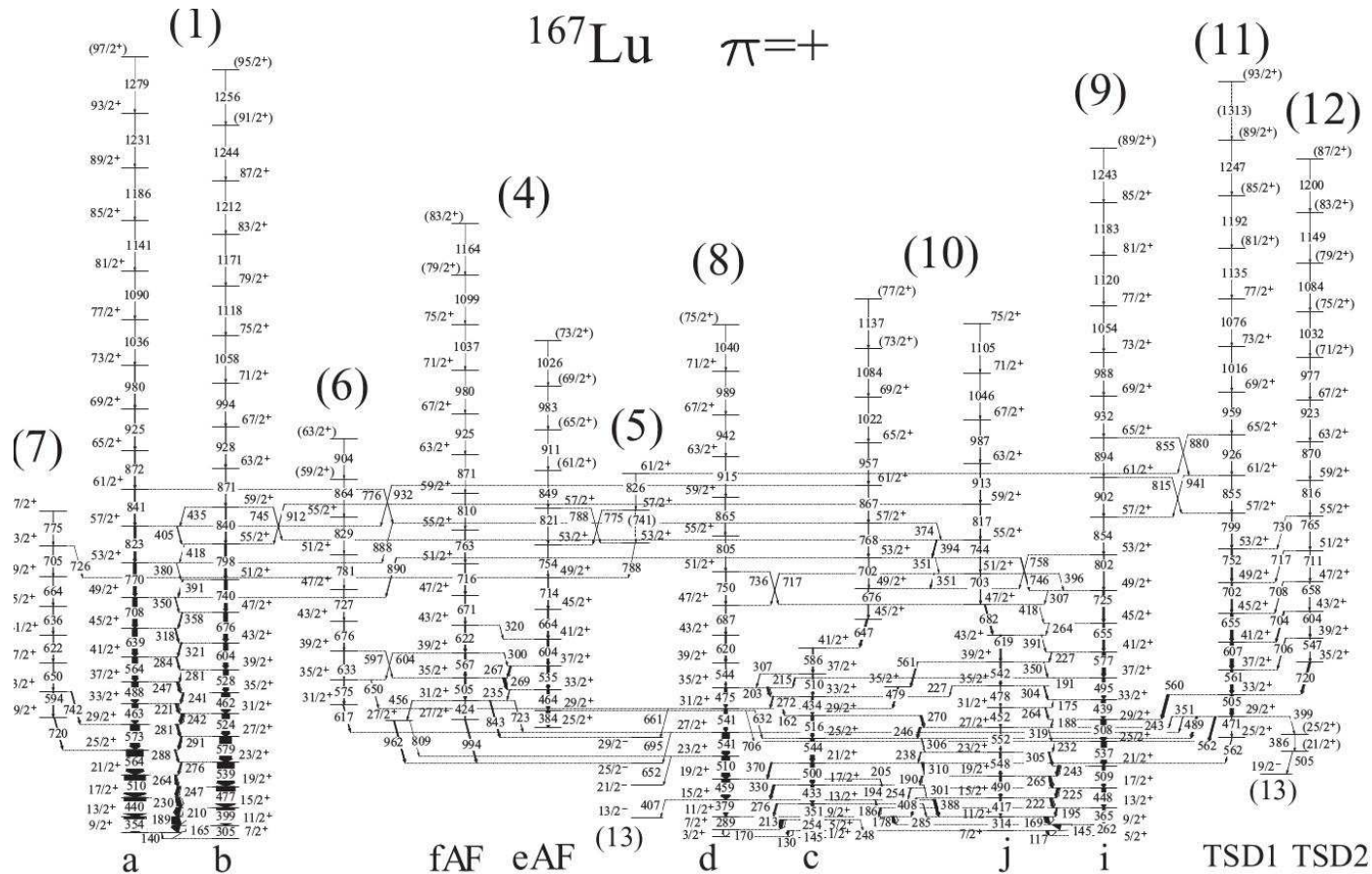
$$\langle E_{pair} \rangle = -2.47 \cdot \exp\left(\frac{I}{29}\right) \text{ MeV}$$

Somewhat stronger pairing for  $(\pi, \alpha)_p = (-, -1/2)$  conf's

No paired crossings at high spin



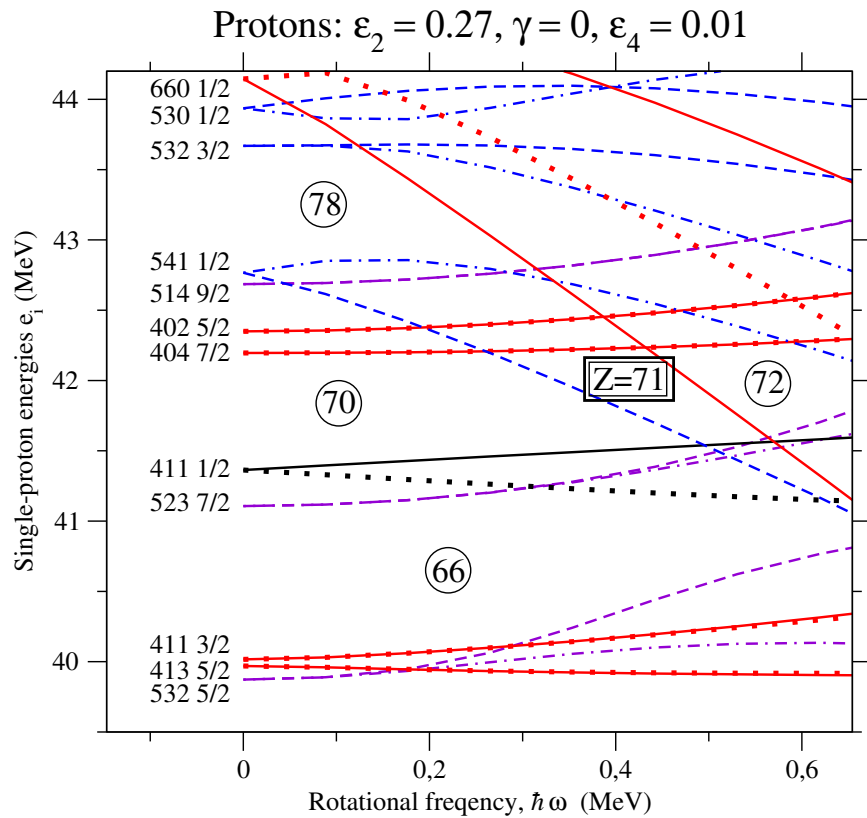
# Observed positive parity bands, $^{167}\text{Lu}$ .



Assignments mainly based on paired band crossings and number of aligned particles (number of quasiparticles).

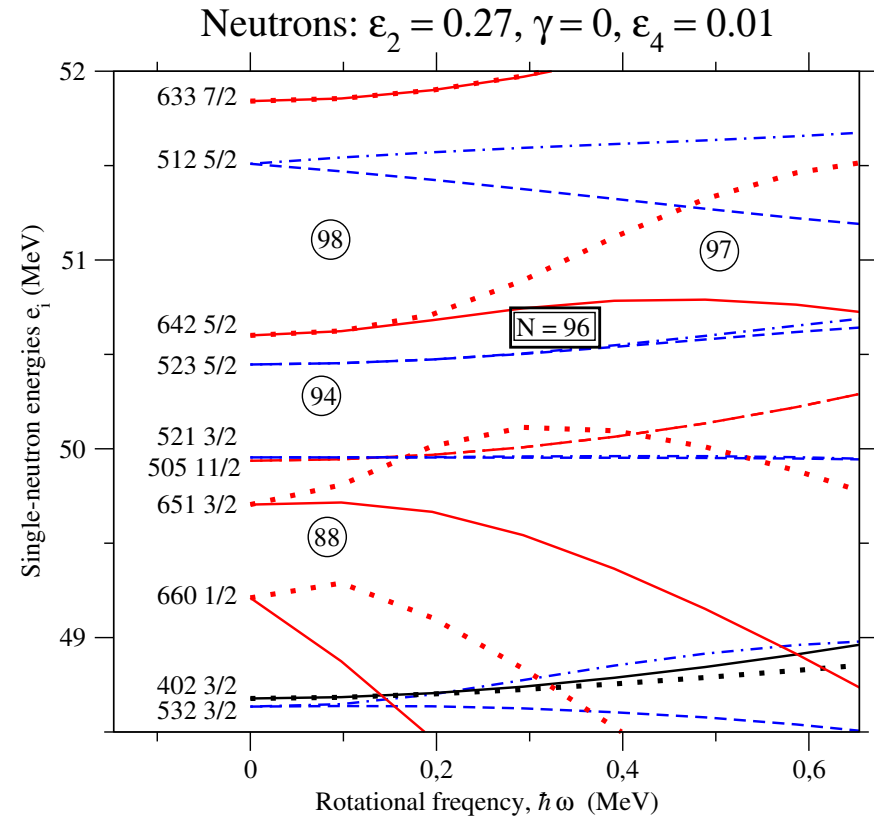
D.G. Roux, W.C. Ma *et al.*, Phys. Rev. C 92, 064313 (2015).

# Single-particle routhians at ground state.



Bands built on:

- 404 7/2, 402 5/2:  $dg$
- 411 1/2:  $sd$
- 523 7/2, 514 9/2:  $h_{11/2}$
- 541 1/2:  $fh$

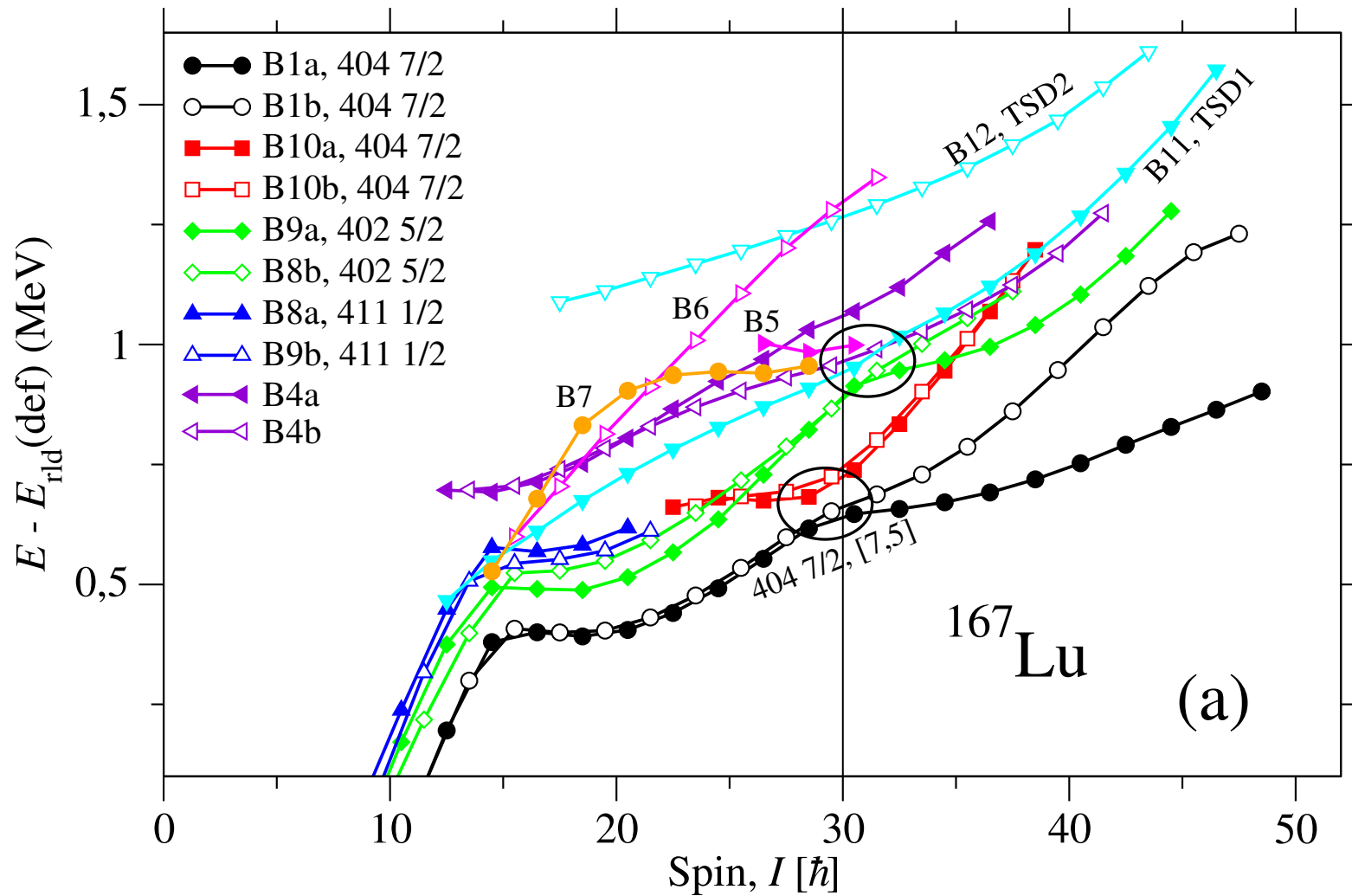


Bands of the type:

- $(fh)^{10} (i_{13/2})^4$
- $(fh)^9 (i_{13/2})^5$



$\pi = +$  bands relative to  $E_{rld}$ .

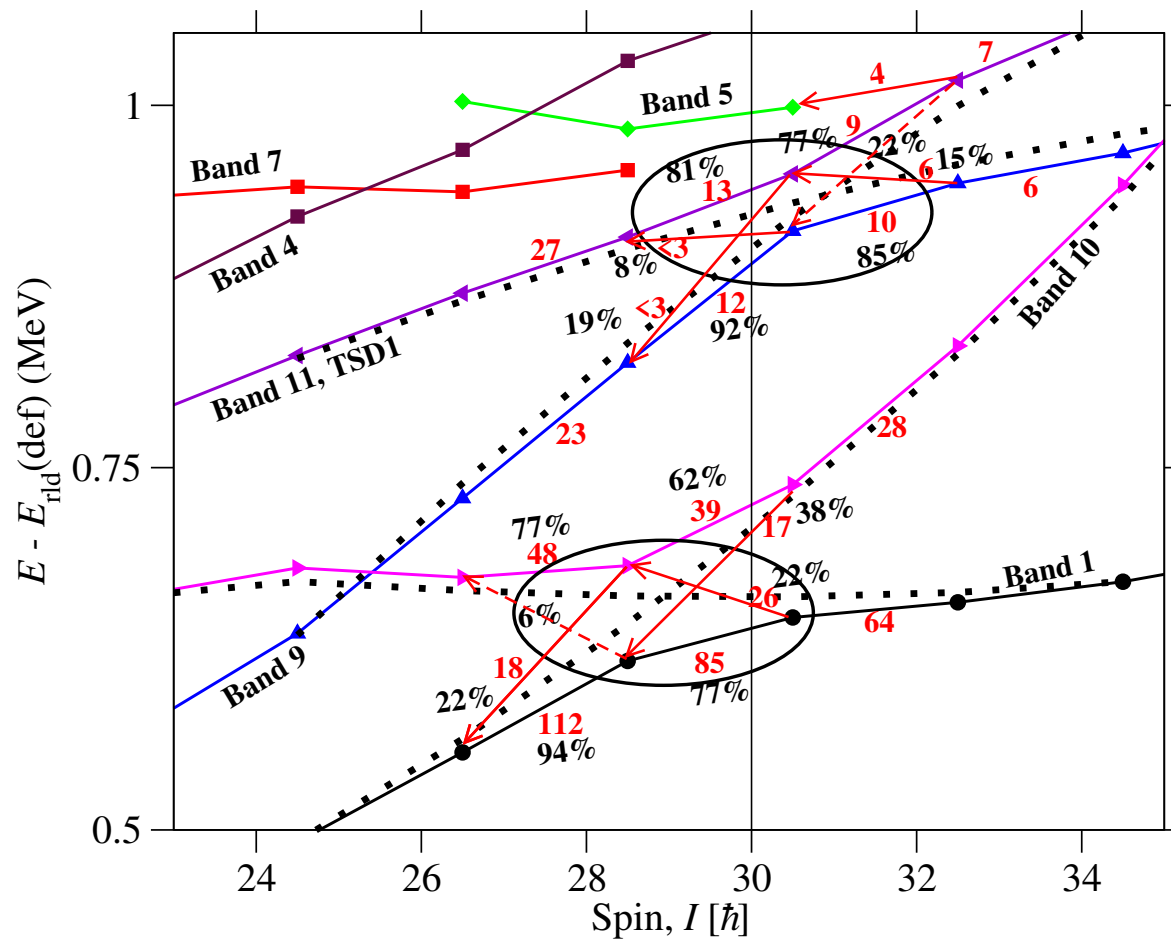


Low-spin bands: 404 7/2, 402 5/2, 411 1/2

Three q-p band: B4

TSD1 and TSD2 (wobbling)

# Band crossings, $\pi = +, \alpha = 1/2$ bands.

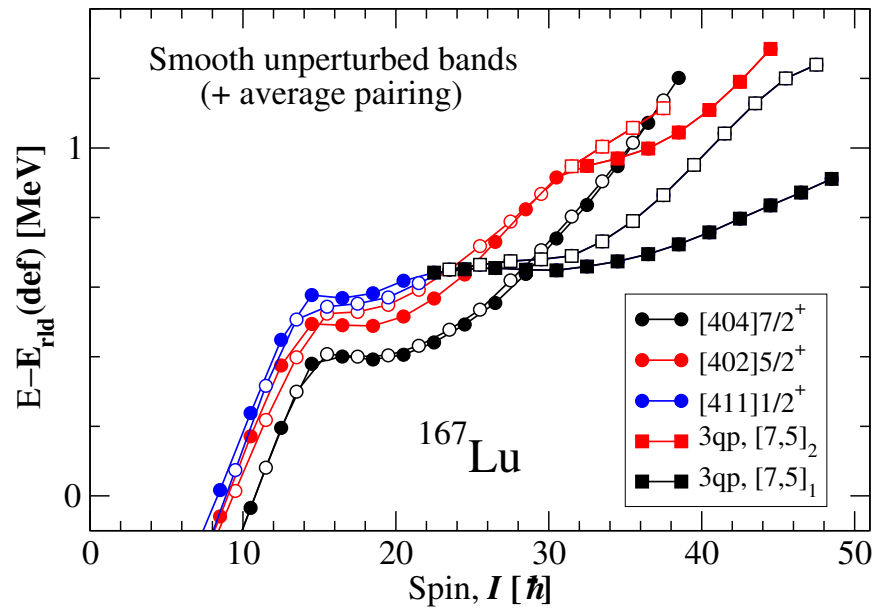


Two-band mixing calculations leading to smooth unperturbed bands:

Clear case for yrast crossing: Bands 1 and 10 (404 7/2 and 3qp band);  $V_{int} = 30$  keV;  $\pm 7$  keV

Questionable; Band 9 (402 5/2) and band 11 (TSD1);  $V_{int} = 24$  keV;  $\pm 10$  keV

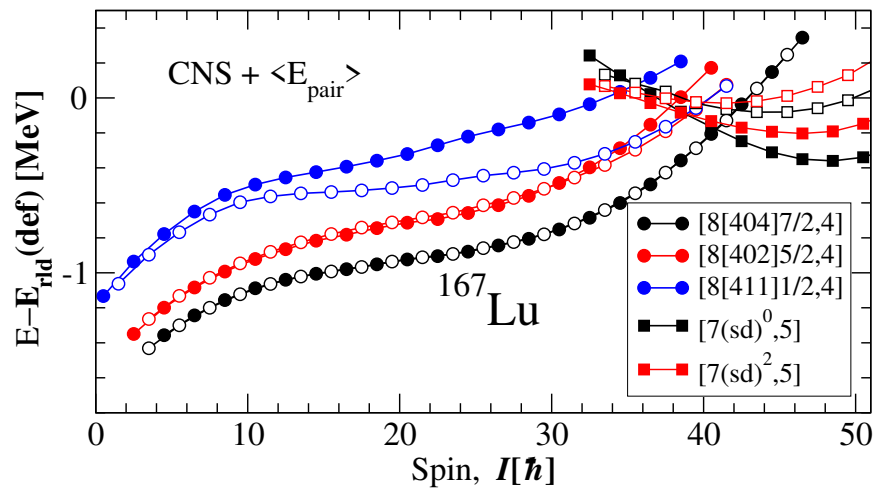
# Proton $N_{osc} = 4$ bands, smooth exp. and calc. CNS.



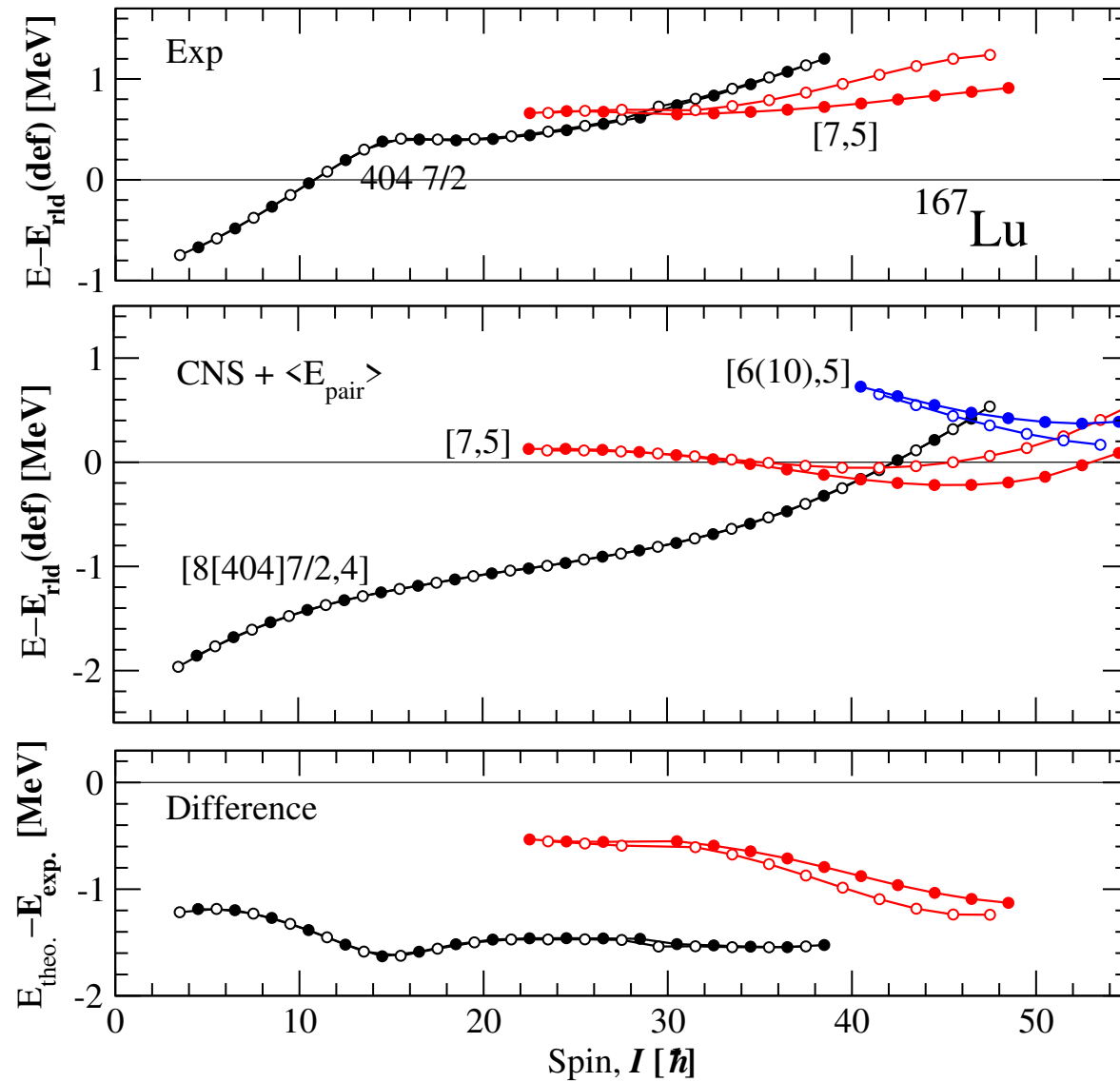
Same general structure in exp. and calc.

Small change in s.p. parameters to lower the 411 1/2 band.

Stronger pairing for 3 qp band, [7.5]  
⇒ band crossing at lower spin.

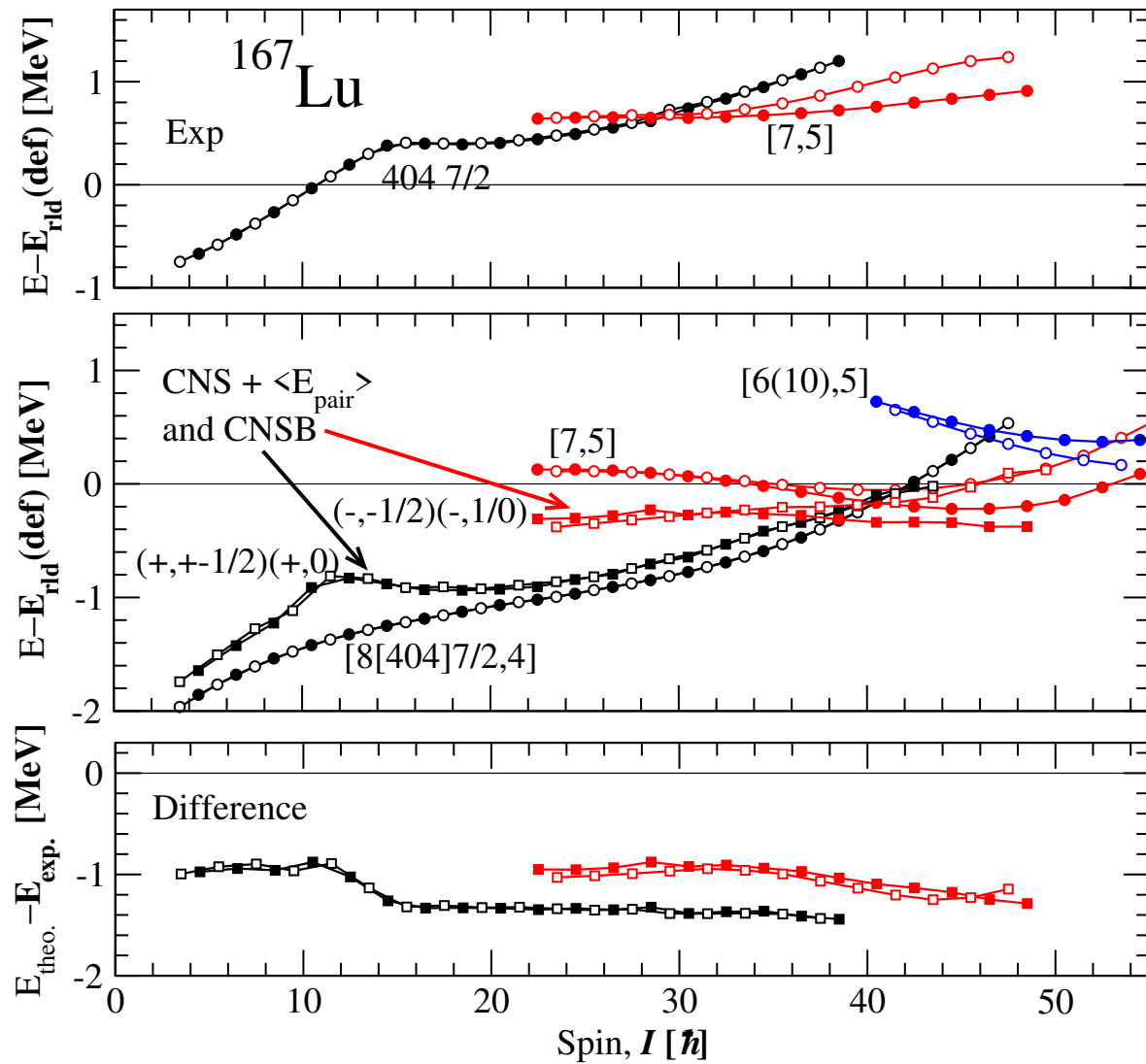


# Yrast states, $\pi = +$ .



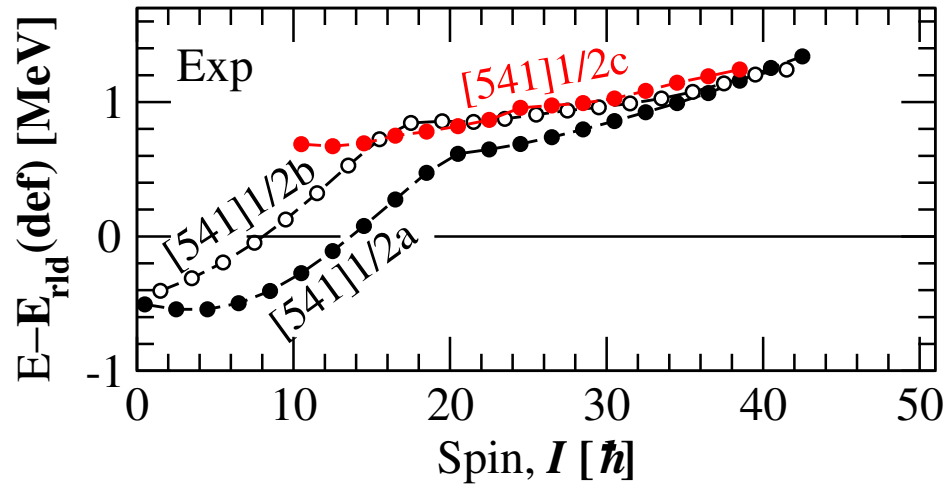
Large scattering of differences, cf. full CNSB calculations.

# Yrast states, $\pi = +$ , CNS vs. CNSB.

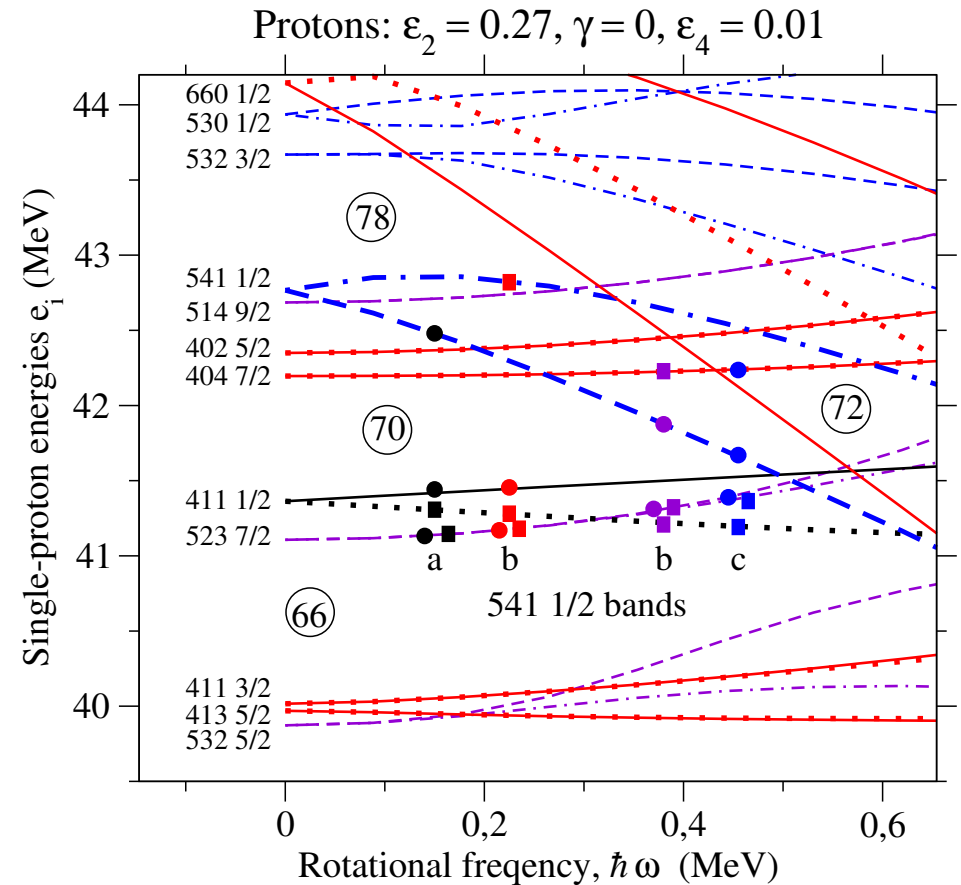


Differences within  $\pm 0.25$  MeV!

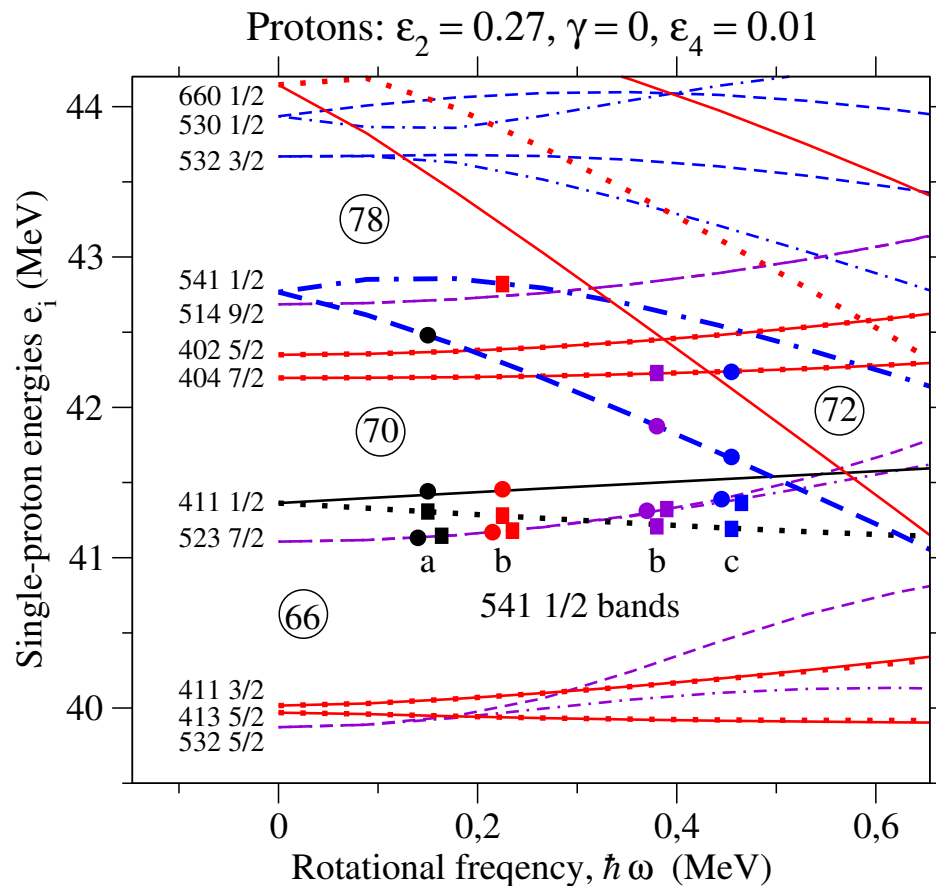
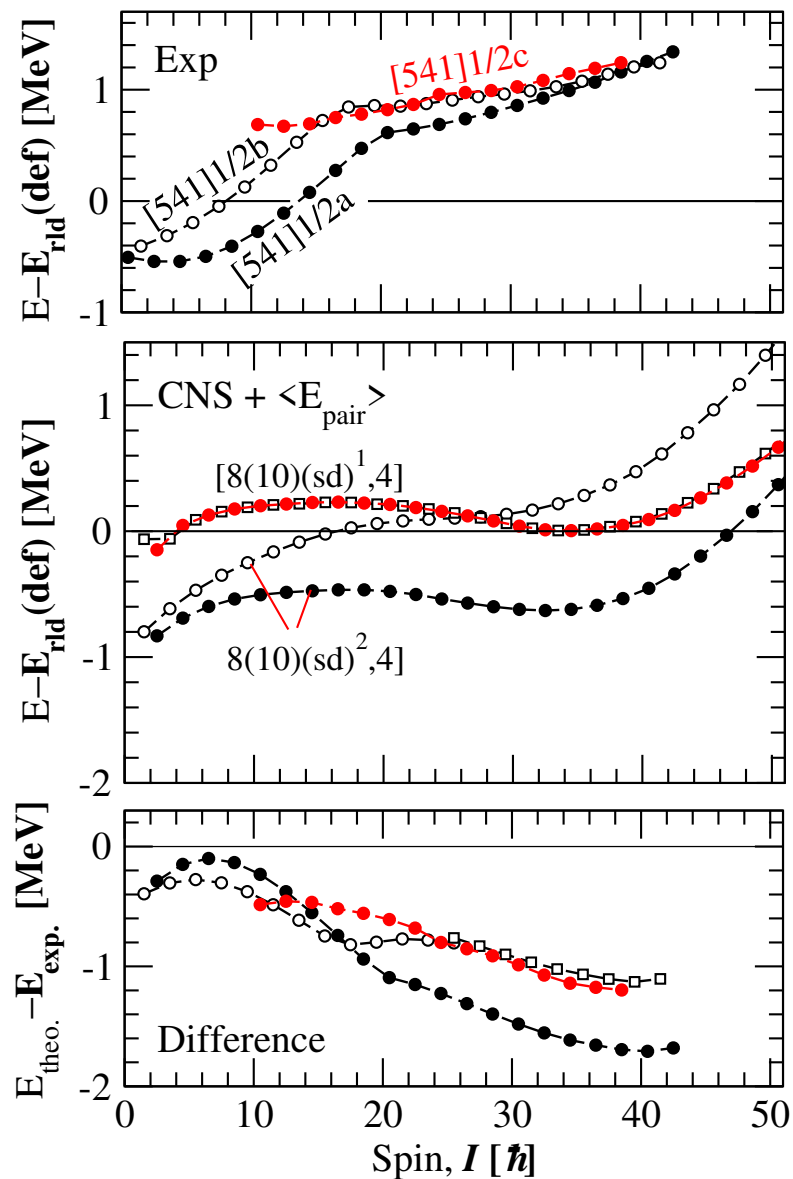
# 541 1/2 bands (from $fh$ shells).



Three bands related 541 1/2 orbital (from  $h_{9/2}f_{7/2}$  ( $fh$ ) shells above the  $Z = 82$  gap).  
Orbitals 13,14  $\rightarrow$  541 1/2 a,b,c



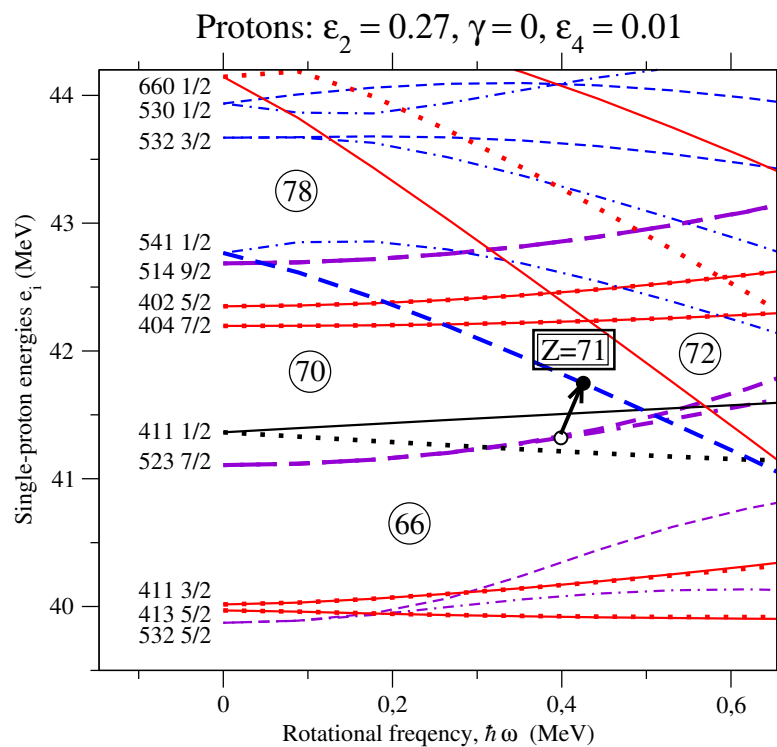
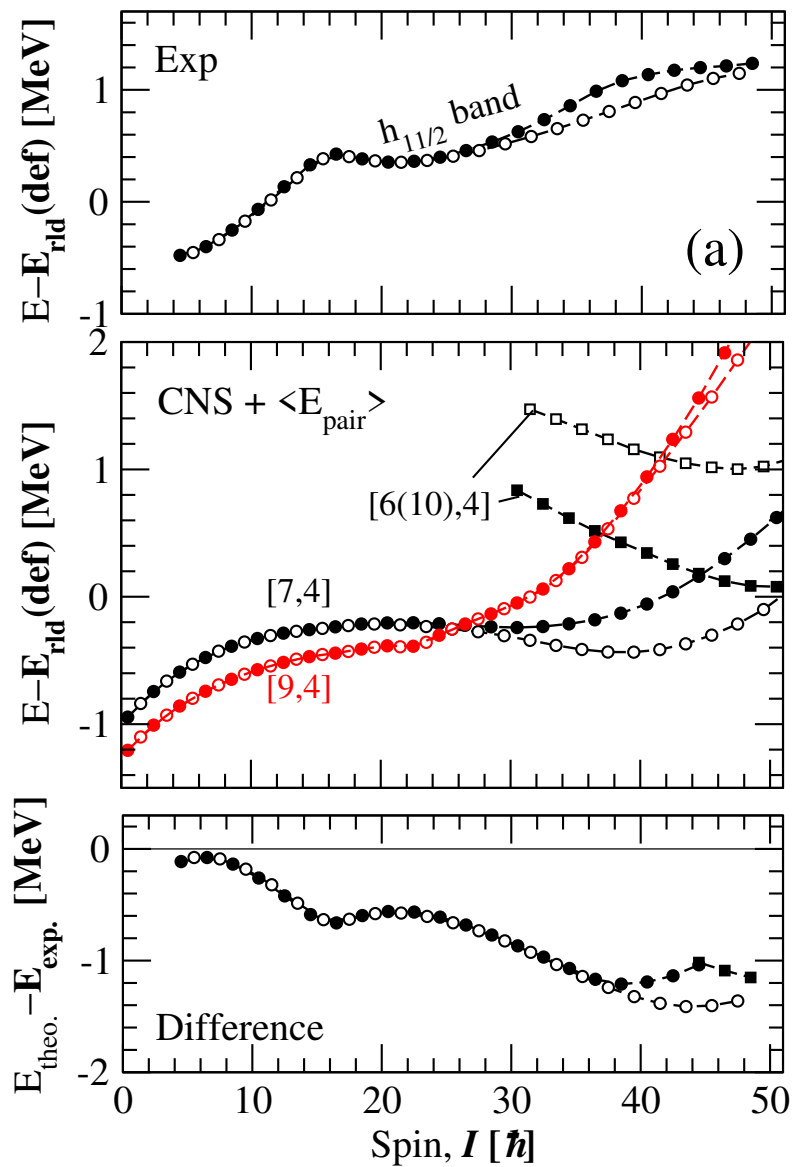
# Exp. vs. calculation for 541 1/2 bands.



With the 404 7/2 orbital closer to the 411 1/2 orbital, bands b,c would come closer to band a.

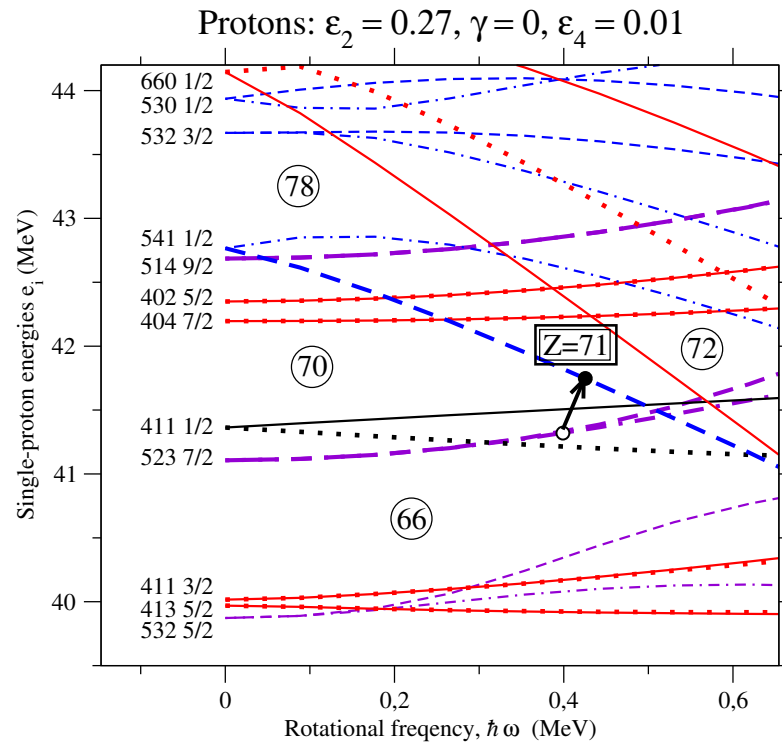
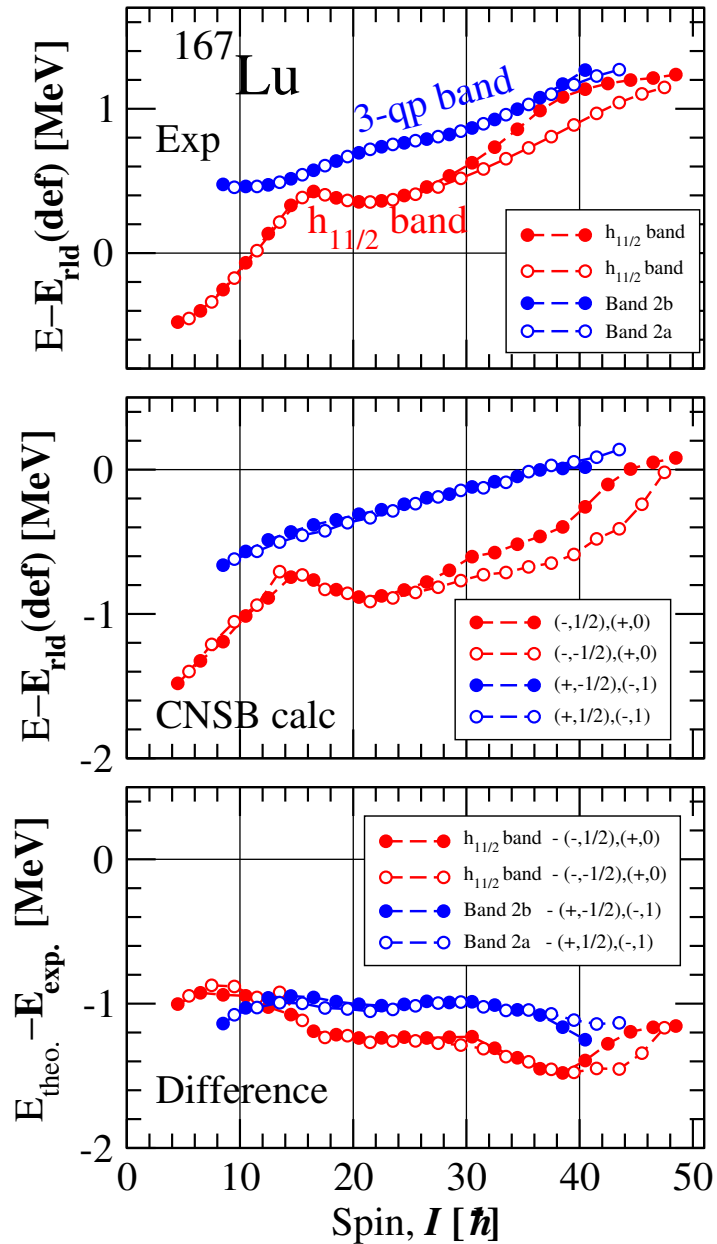
Decoupling will lower the calc. energies at low spin.

# $h_{11/2}$ band vs. CNS calc.





# $h_{11/2}$ and $\pi = -$ 3qp band vs. CNSB calc.



# Conclusions

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- Combination of CNS and CNSB powerful method to analyze nuclear level schemes.
- Assignments to observed rotational bands in  $^{167}\text{Lu}$ .
- Unpaired CNS calculations useful because of possibility to make detailed configuration assignments (instructive at all spin values).
- Improved and modified understanding compared with assignments in experimental paper (PRC 92, 064313). Those assignments are mainly based on 'paired' band crossings.
- Such paired band crossings appear to be important only below  $I \lesssim 30$  (AB and BC crossing for  $^{167}\text{Lu}$ ), cf. original article by Bengtsson and Frauendorf.
- High spin bands are mainly formed from the gradual alignment of the spin vectors of the valence particles.
- Rotational bands best characterized by the number of particles in  $j$ -shells or groups of  $j$ -shells, e.g.  $\pi(h_{11/2})^8$  conf's in  $^{167}\text{Lu}$  yrast at low spin while  $\pi(h_{11/2})^6$  become more favoured in energy at high spin.

# Observed $\mathcal{J}^{(2)}$ for 3 qp and TSD bands, $^{167}\text{Lu}$ .

