



First Identification of Collective Band Structure in Odd-odd ^{166}Re

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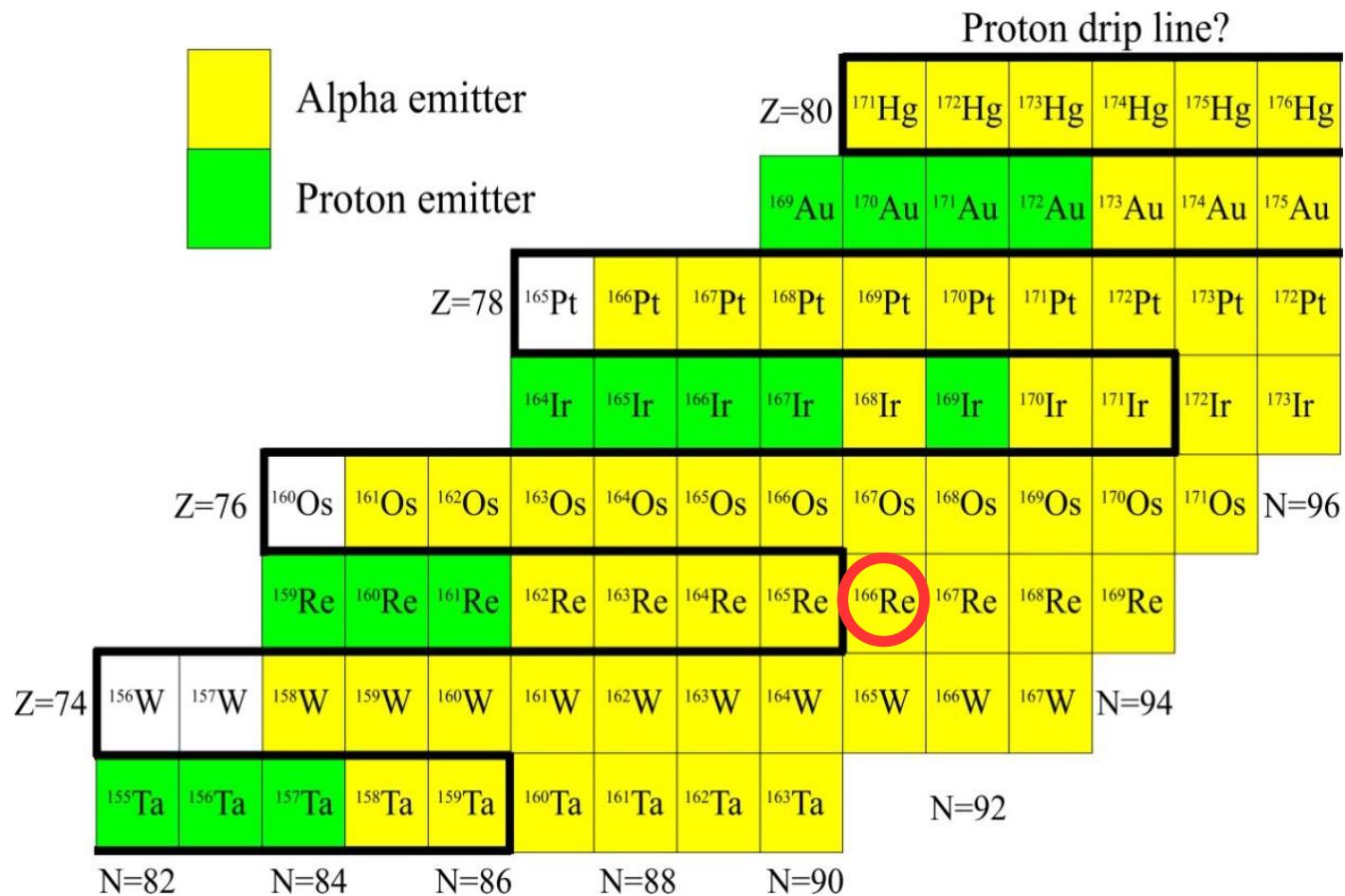
Tsinghua University, Beijing, China

JUROGAM/GREAT Collaboration

Outline of the Talk

- **Spectroscopy at the Proton Drip Line**
- **Experimental Setup**
- **Preliminary Results**
- **Conclusion**

Proton drip line



courtesy of RD Page, LISA presentation

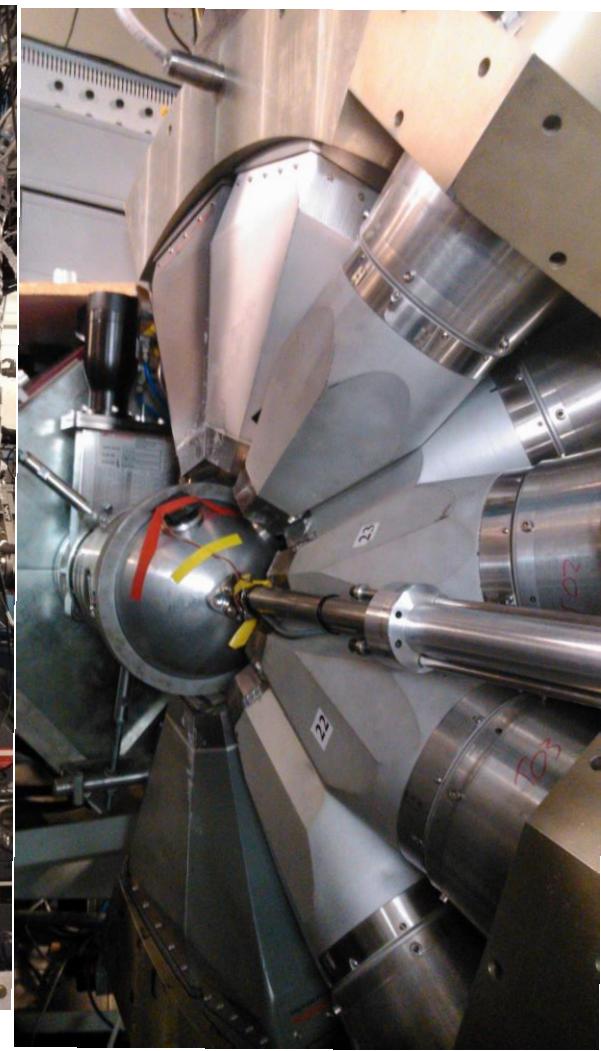
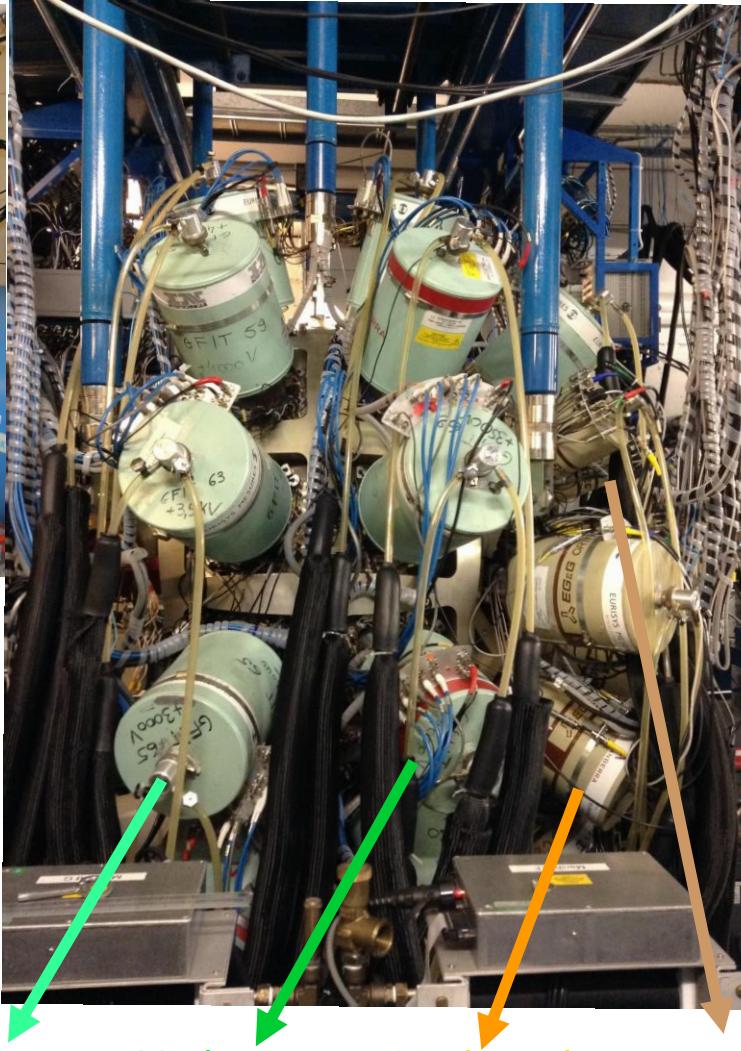
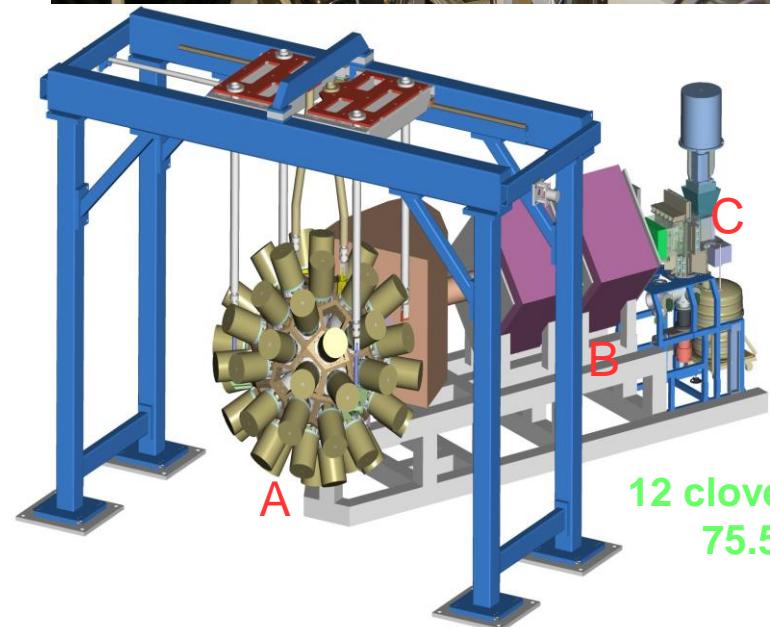
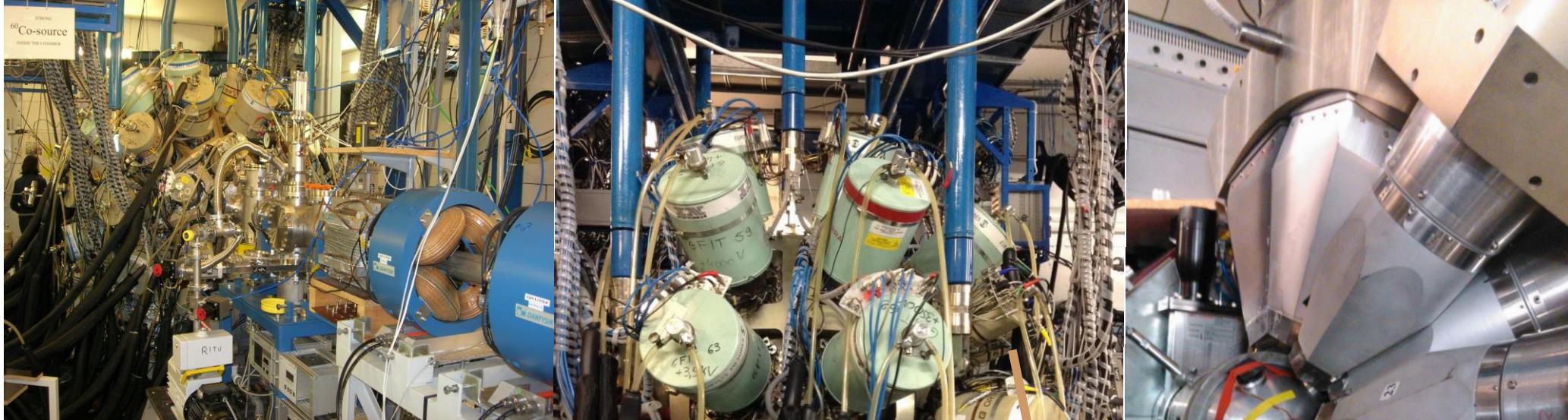
Identified collective bands

| | Compound | | | | | | | | |
|----|----------|-------|-------|-------|-------|-------|-------|-------|-------|
| 78 | Pt162 | Pt163 | Pt164 | Pt165 | Pt166 | Pt167 | Pt168 | Pt169 | Pt170 |
| 77 | Ir161 | Ir162 | Ir163 | Ir164 | Ir165 | Ir166 | Ir167 | Ir168 | Ir169 |
| 76 | Os160 | Os161 | Os162 | Os163 | Os164 | Os165 | Os166 | Os167 | Os168 |
| 75 | Re159 | Re160 | Re161 | Re162 | Re163 | Re164 | Re165 | Re166 | Re167 |
| 74 | W158 | W159 | W160 | W161 | W162 | W163 | W164 | W165 | W166 |
| 73 | Ta157 | Ta158 | Ta159 | Ta160 | Ta161 | Ta162 | Ta163 | Ta164 | Ta165 |
| 72 | Hf156 | Hf157 | Hf158 | Hf159 | Hf160 | Hf161 | Hf162 | Hf163 | Hf164 |
| 71 | Lu155 | Lu156 | Lu157 | Lu158 | Lu159 | Lu160 | Lu161 | Lu162 | Lu163 |
| | 84 | 86 | 88 | 90 | 92 | | | | |
| | 1p | 2p | 3p | 4p | 5p | 6p | 7p | | |

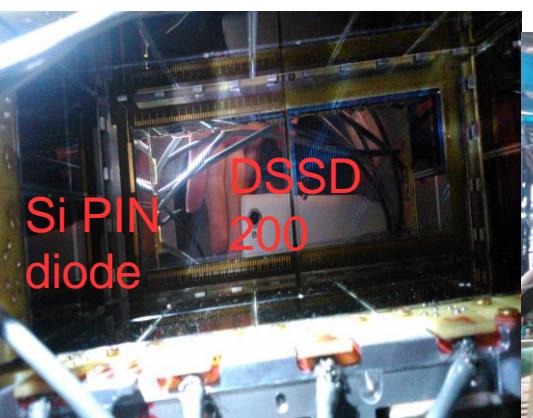
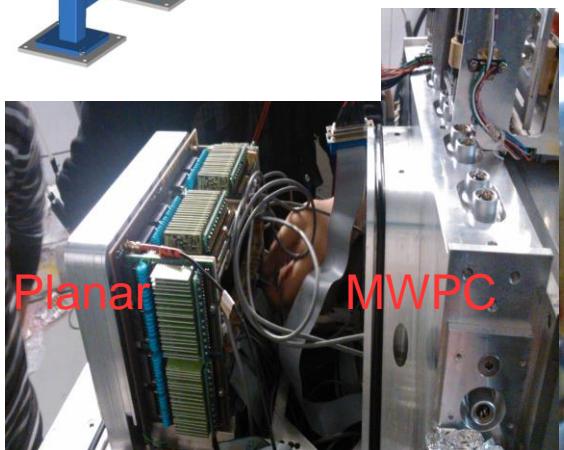
Experimental Setup

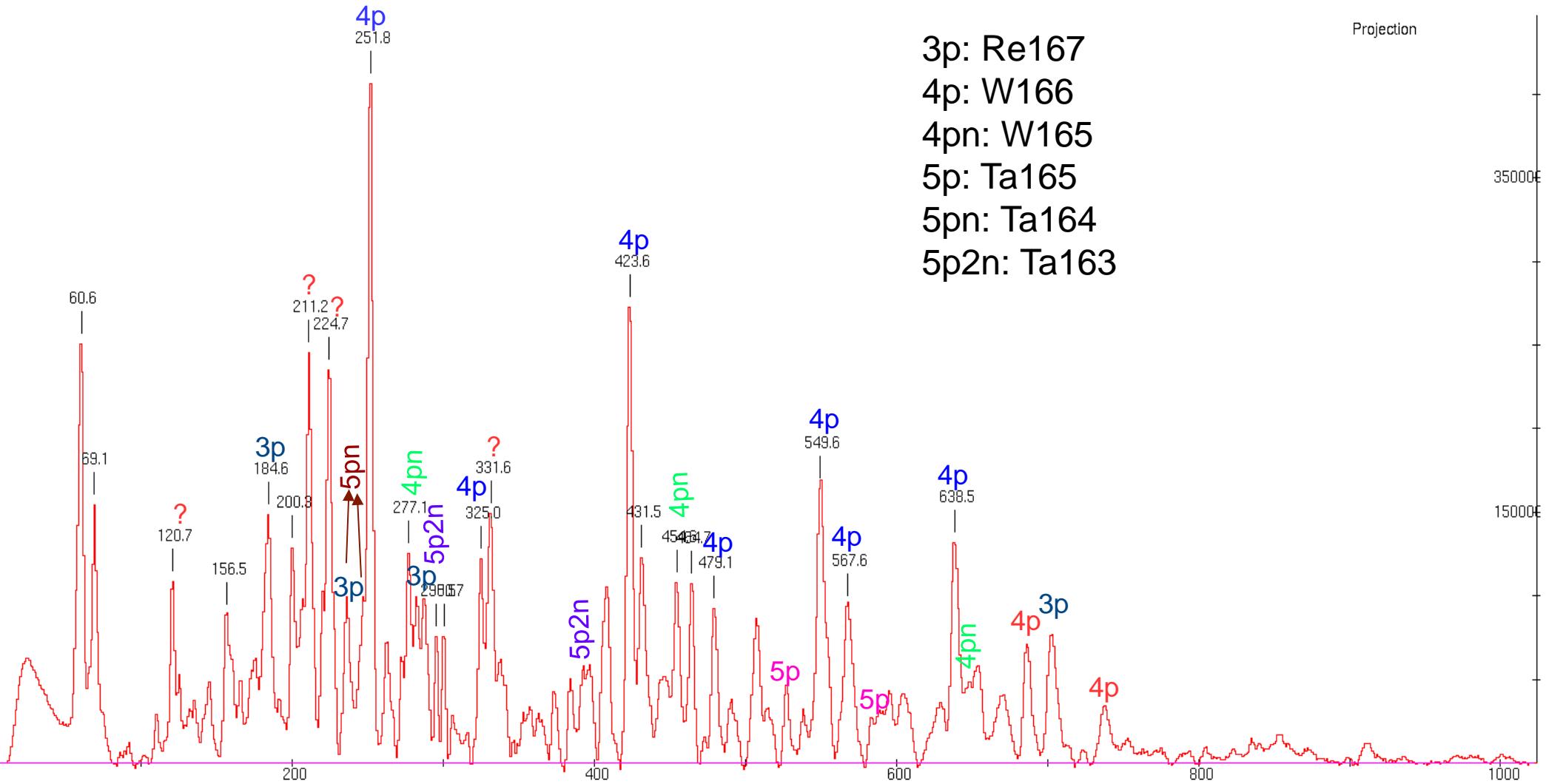
- Reaction: $^{92}\text{Mo}(^{78}\text{Kr}, 3\text{pn})^{166}\text{Re}$
- $E_{\text{beam}} = 380 \text{ MeV/u}$
- Accelerator: K130 cyclotron at the University of Jyväskylä
- Target: 0.6-mg/cm² ^{92}Mo with 1-mg/cm² Ta support
- DPUNS Plunger: 1-mg/cm² Mg degrader with the distances of 5, 100, 200, 500, 1000, 2000, 3000, 5000, 8000um
- Setup: **JUROGAM II(24 clovers + 15 tapered phase I) Germanium
RITU(Recoil Ion Transport Unit) gas-filled recoil separator
GREAT(Gamma Recoil Electron Alpha Tagging) Spectrometer
MWPC
Si PIN diode detectors
DSSD: 120x + 80y
3 clovers + planar(24x + 12y) in FP**
- Total photopeak efficiency: 4.2% at 1.3 MeV

^{60}Co -source
INSIDE THE CHAMBER



A: JUROGAM II
B: RITU
C: GREAT





Identification of ^{166}Re

Total projection in the Cube

Identification of ^{166}Re

The Known Excited States

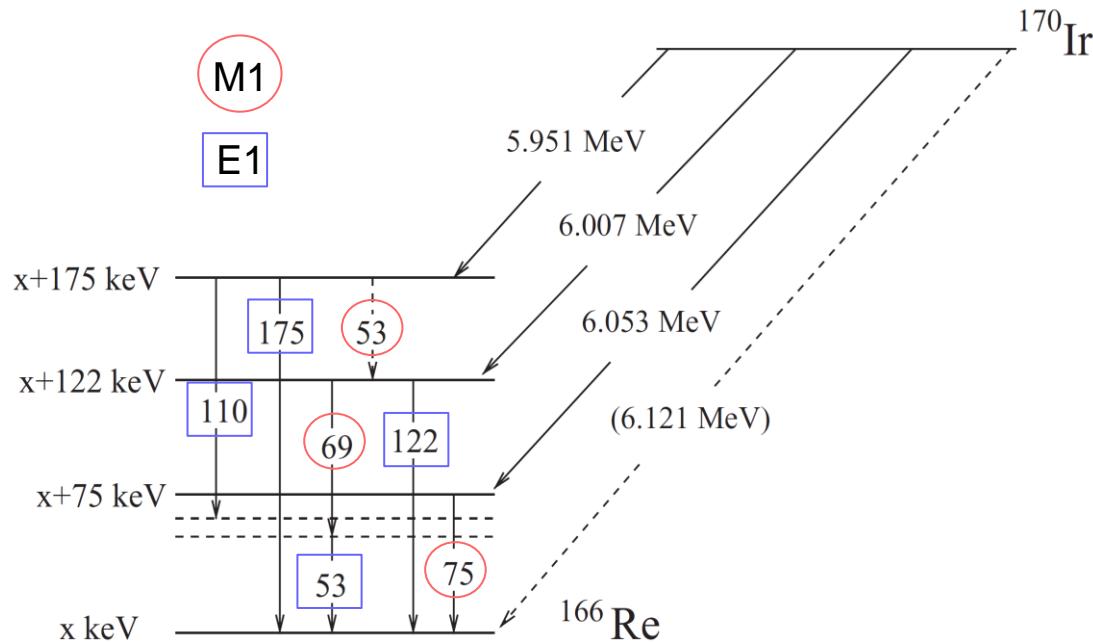


FIG. 8. Tentative level structure of ^{166}Re deduced from the α -decay study of ^{170}Ir .

B. Hadinia, *et al*, PRC **76**, 044312 (2007)

Level scheme of ^{166}Re

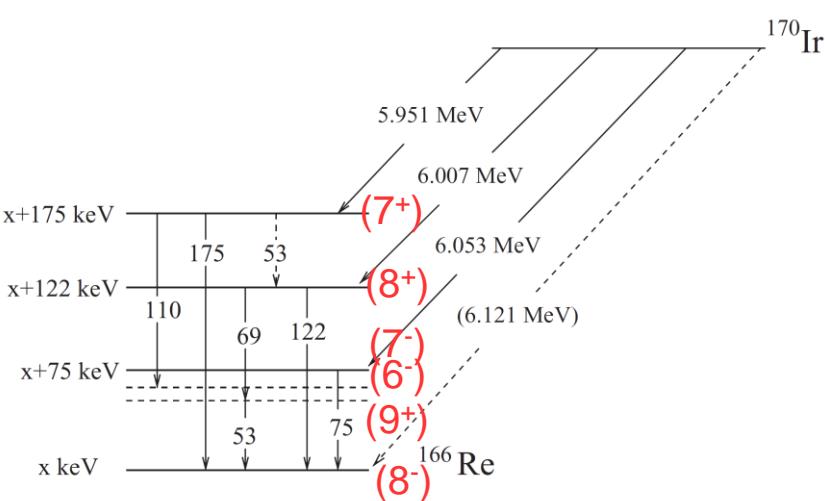
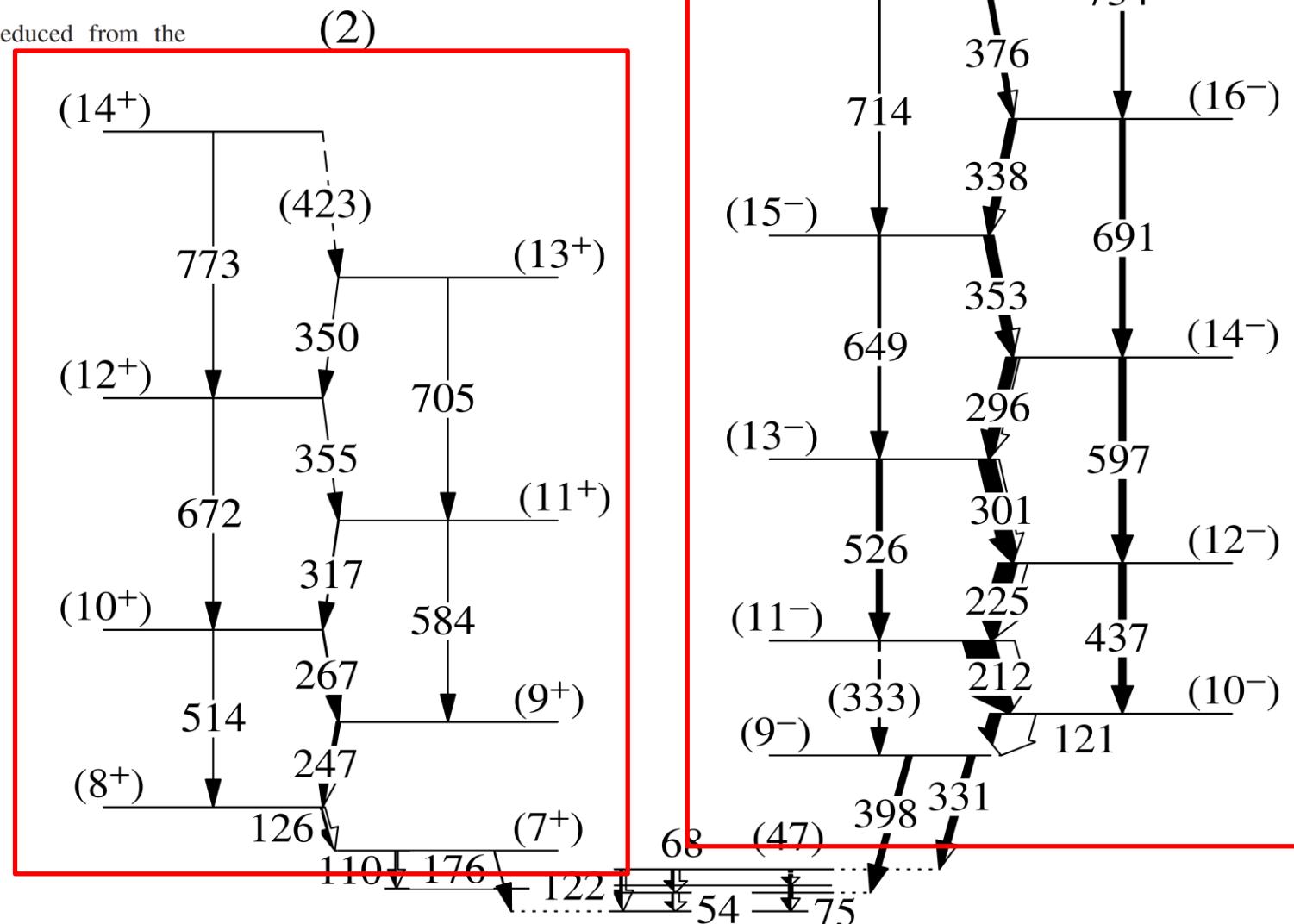
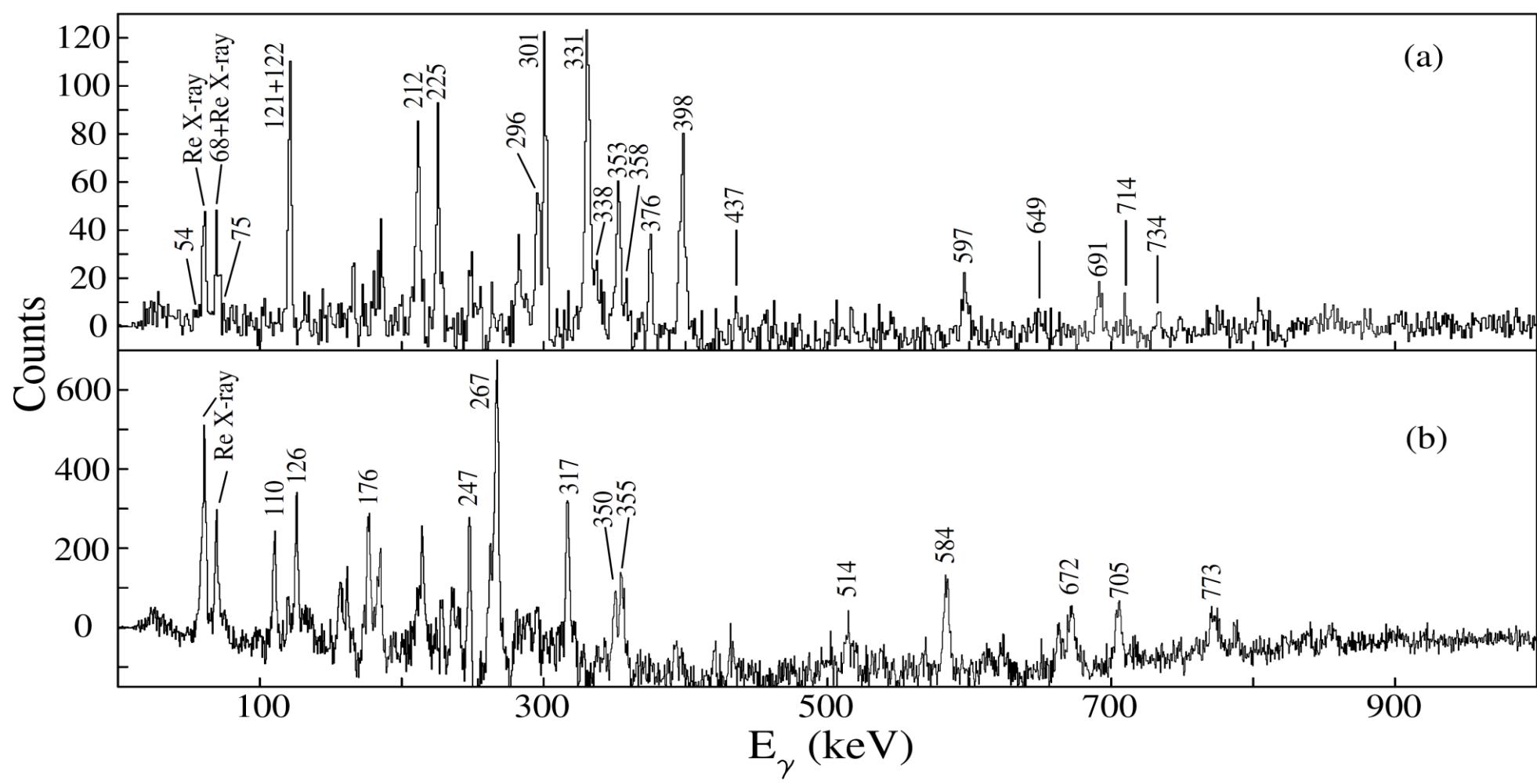
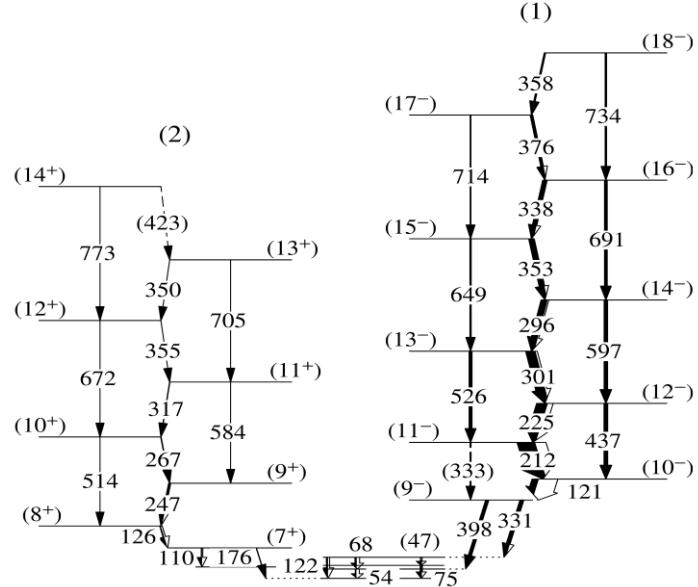
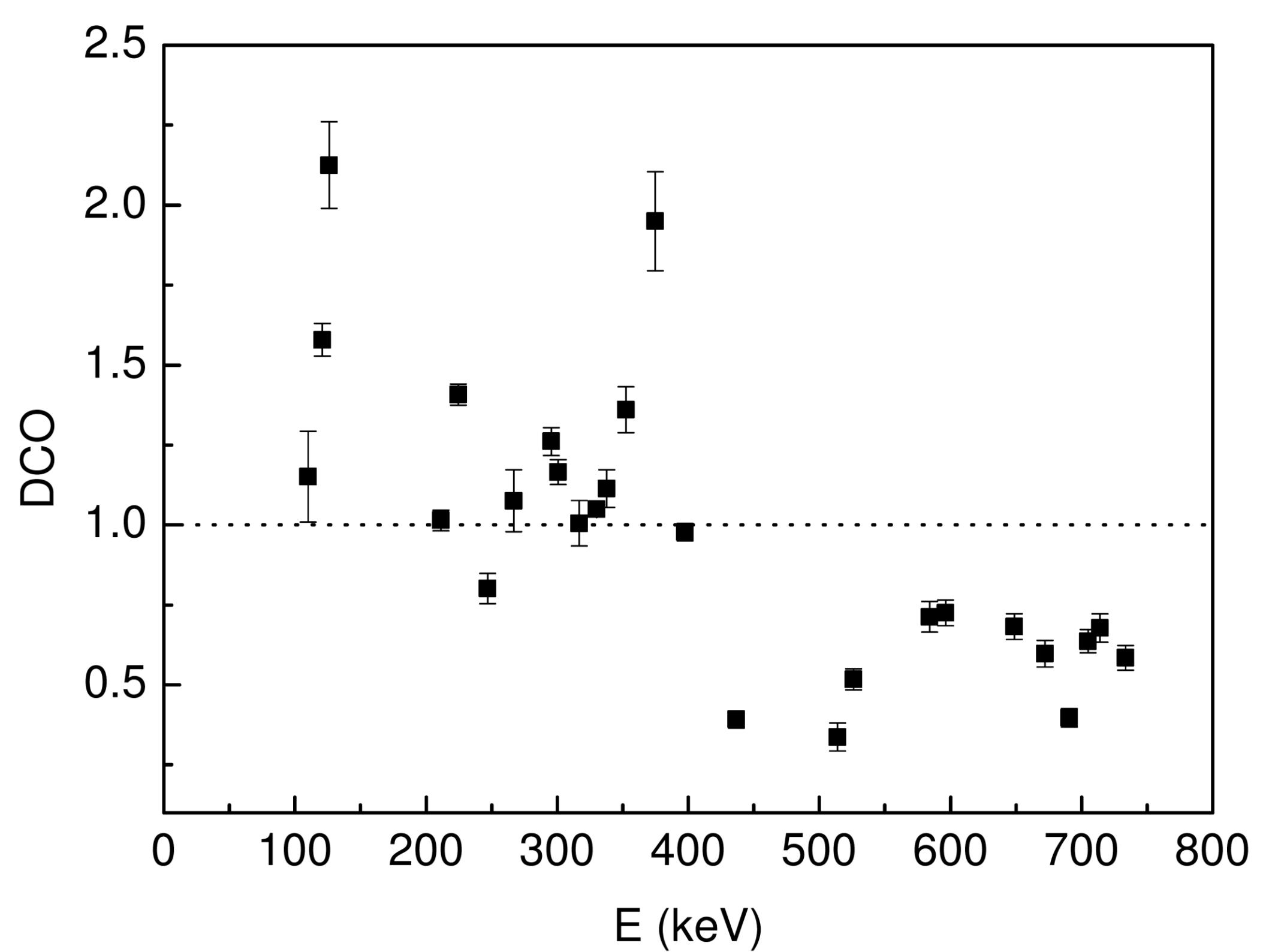


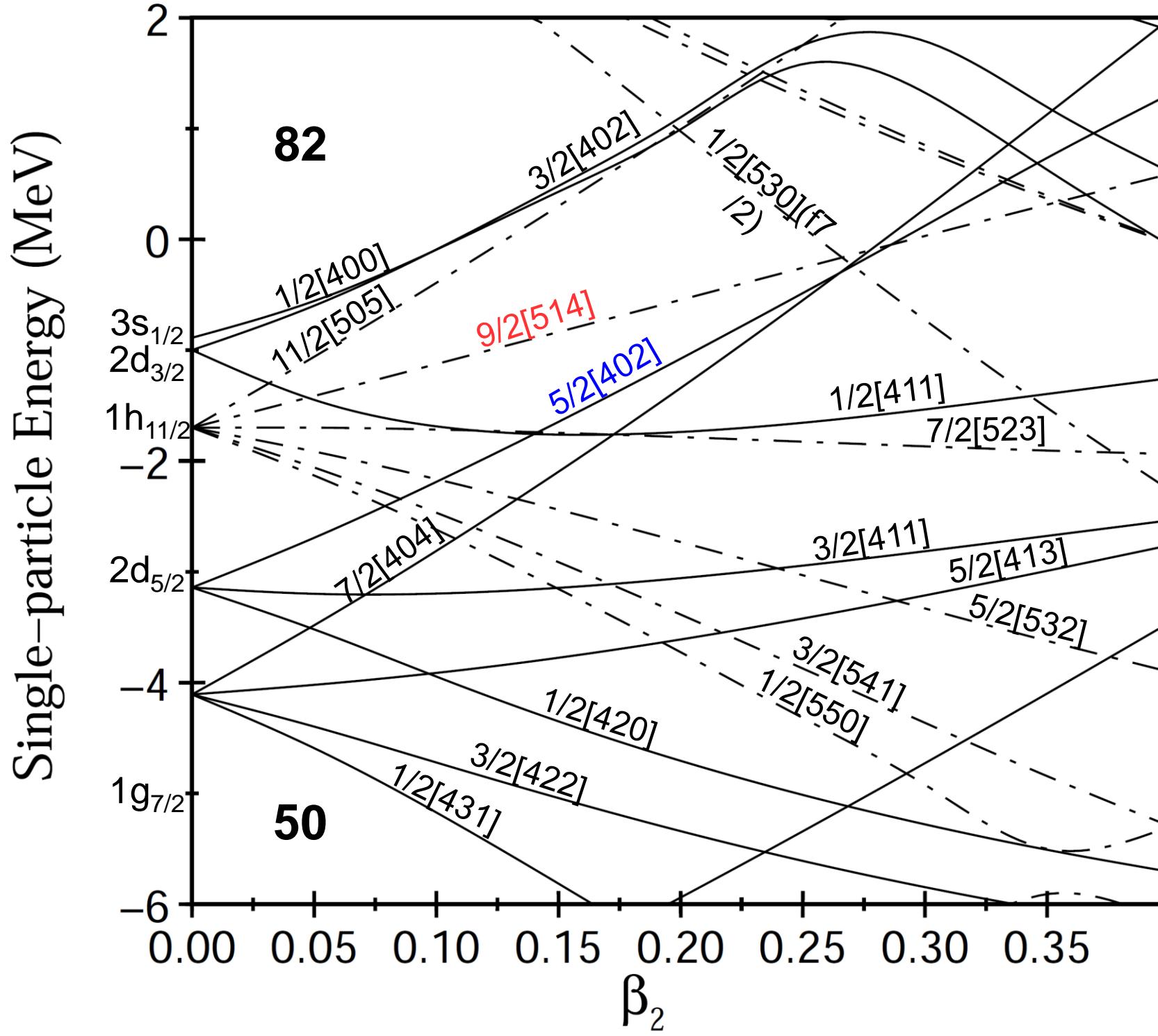
FIG. 8. Tentative level structure of ^{166}Re deduced from the α -decay study of ^{170}Ir .



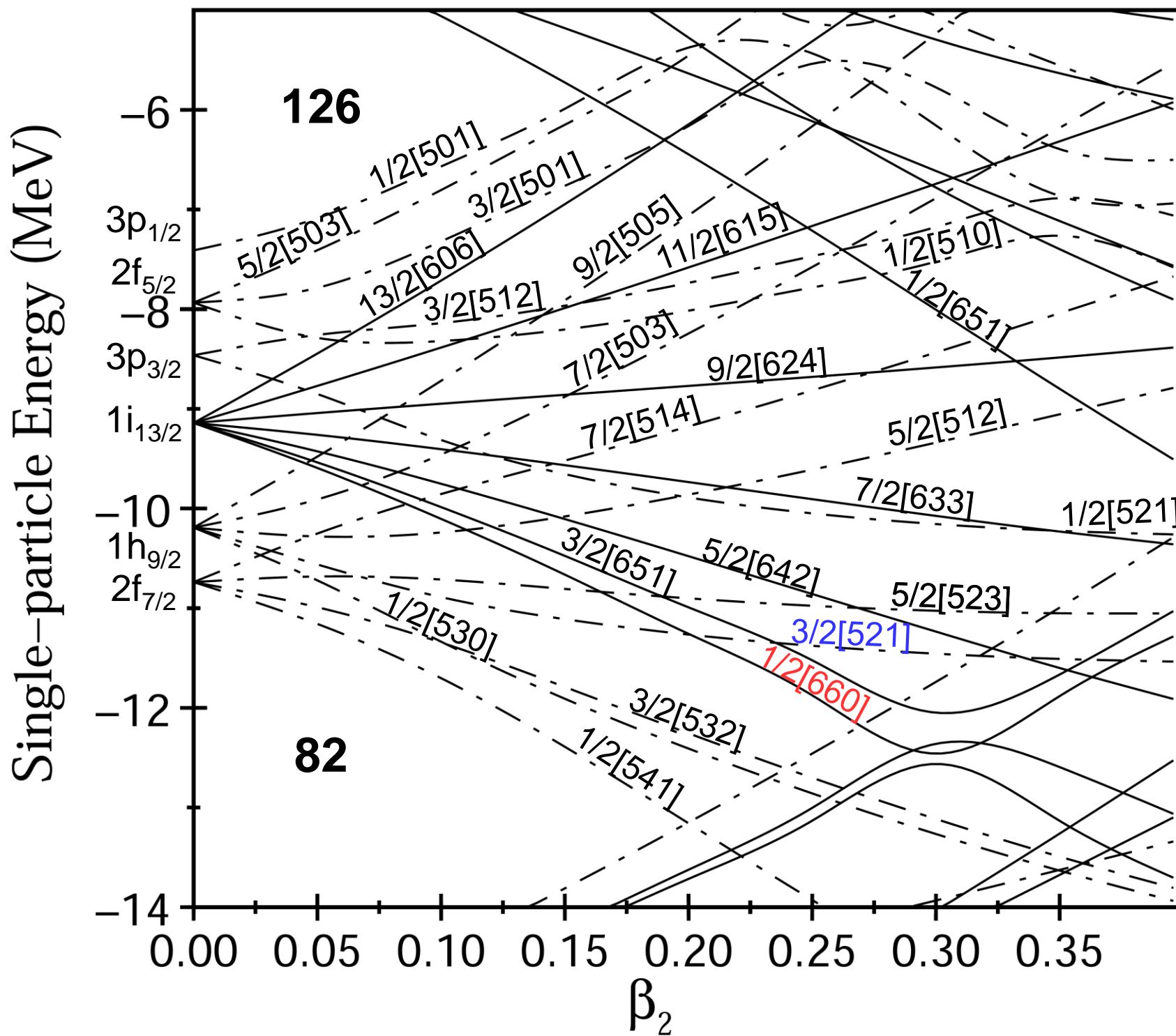


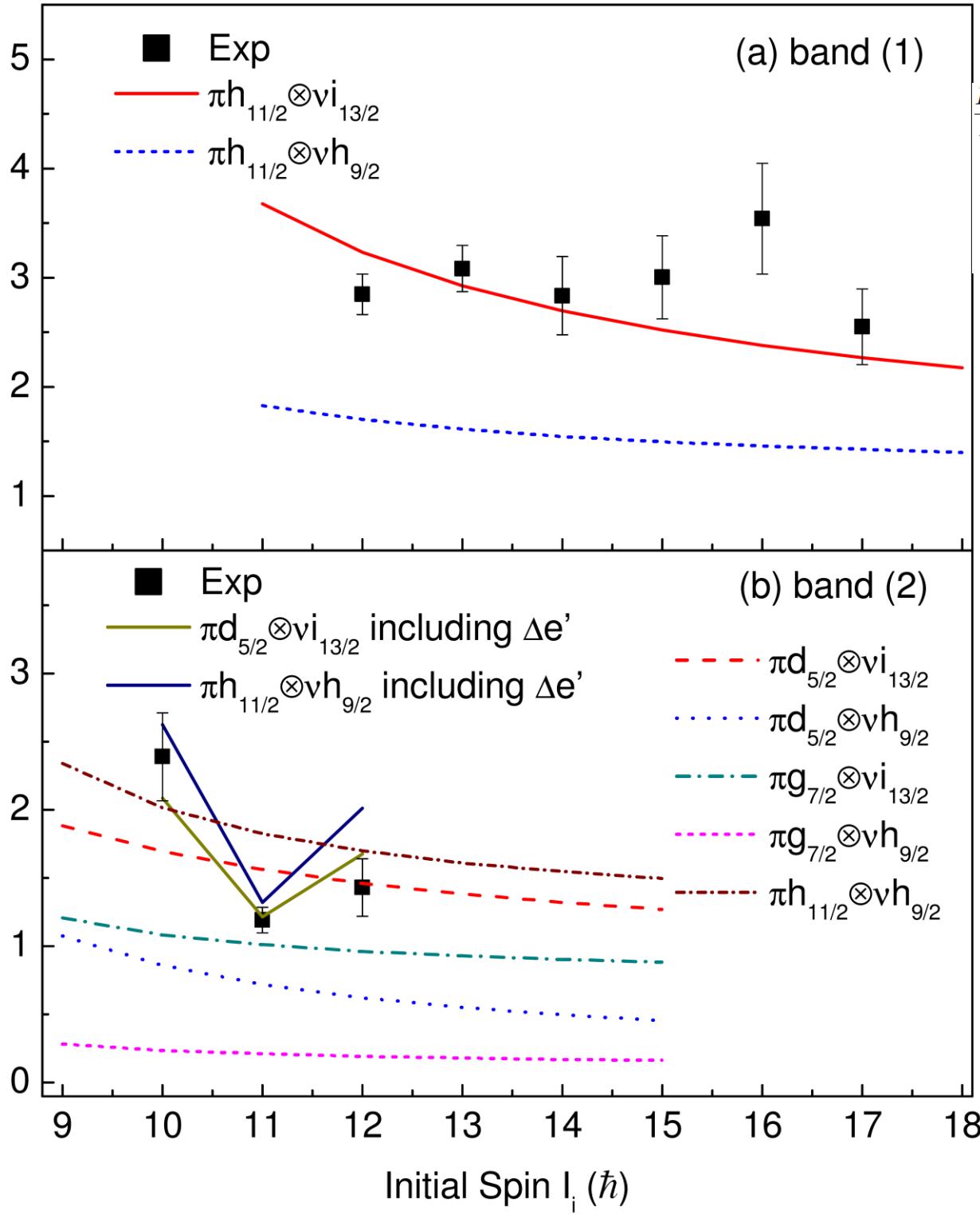


\Proton single particle levels : Universal Woods-Saxon potential



\Neutron single particle levels : Universal Woods-Saxon potential





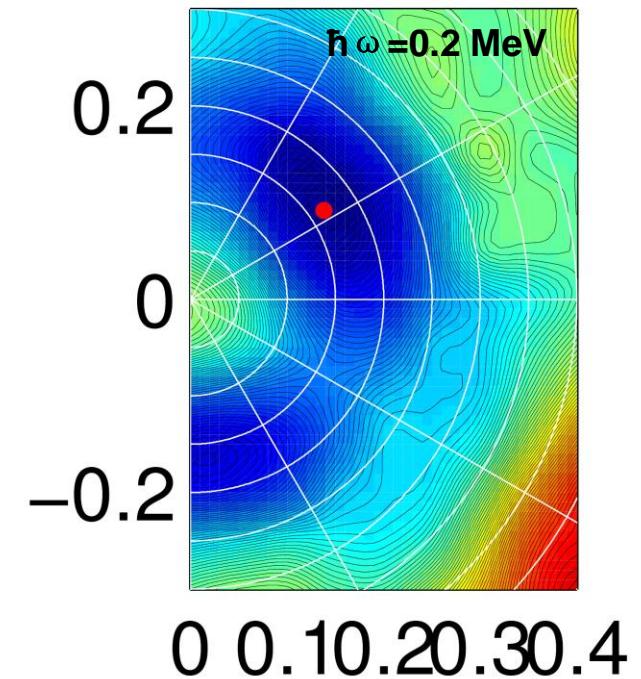
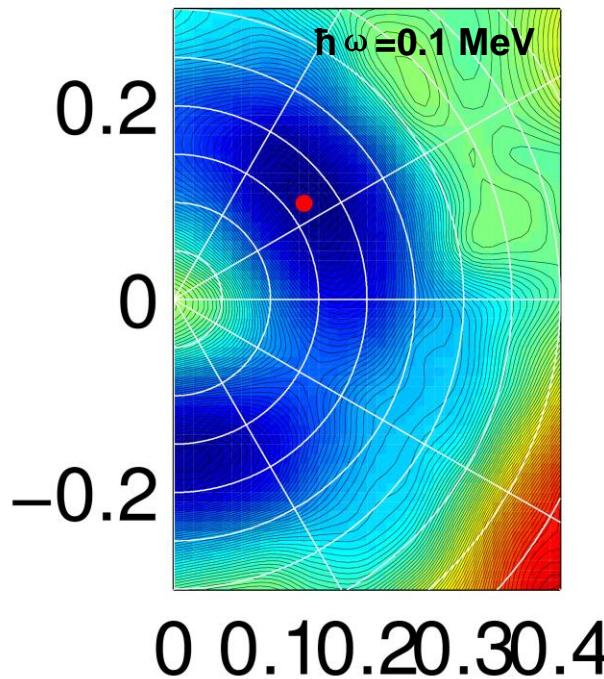
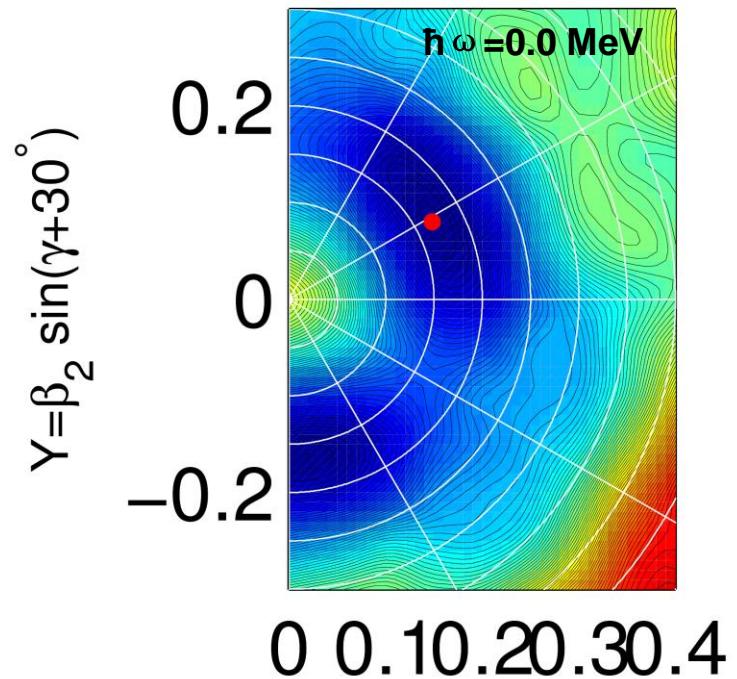
Theo.

$$\frac{B(M1; I \rightarrow I-1)}{B(E2; I \rightarrow I-2)} = \frac{12}{5Q_0^2 \cos^2(\gamma + 30^\circ)} \left[1 - \frac{K^2}{(I-1/2)^2}\right]^{-2} \frac{K^2}{I^2} \times \\ [(g_1 - g_R)(\sqrt{I^2 - K^2} - i_1)(1 \pm \frac{\Delta e'}{\hbar\omega}) - (g_2 - g_R)i_2]^2$$

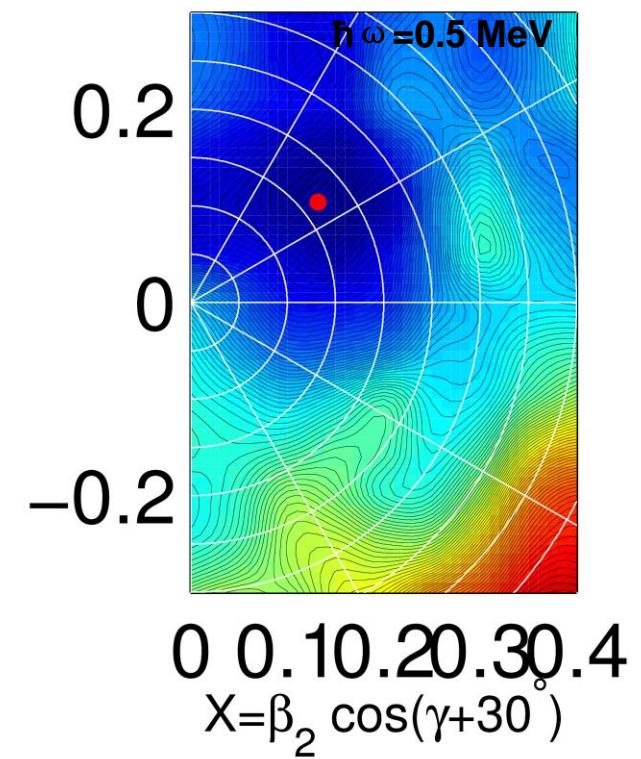
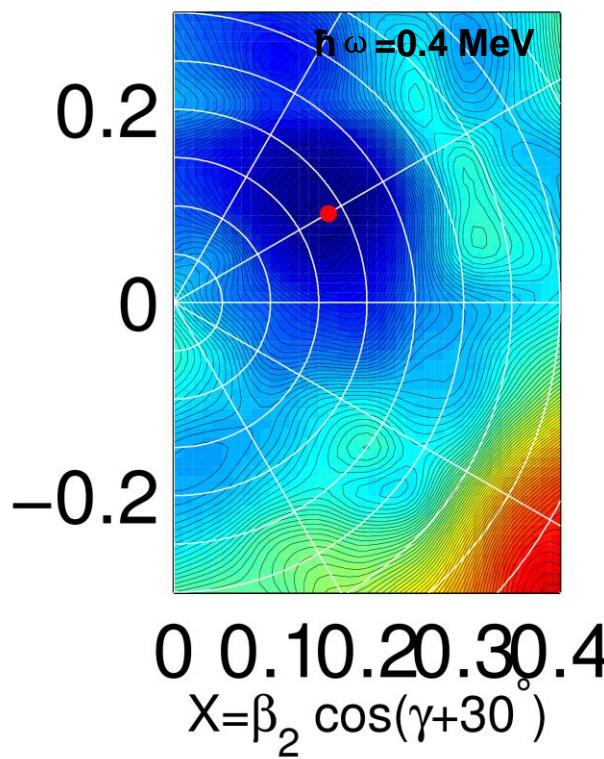
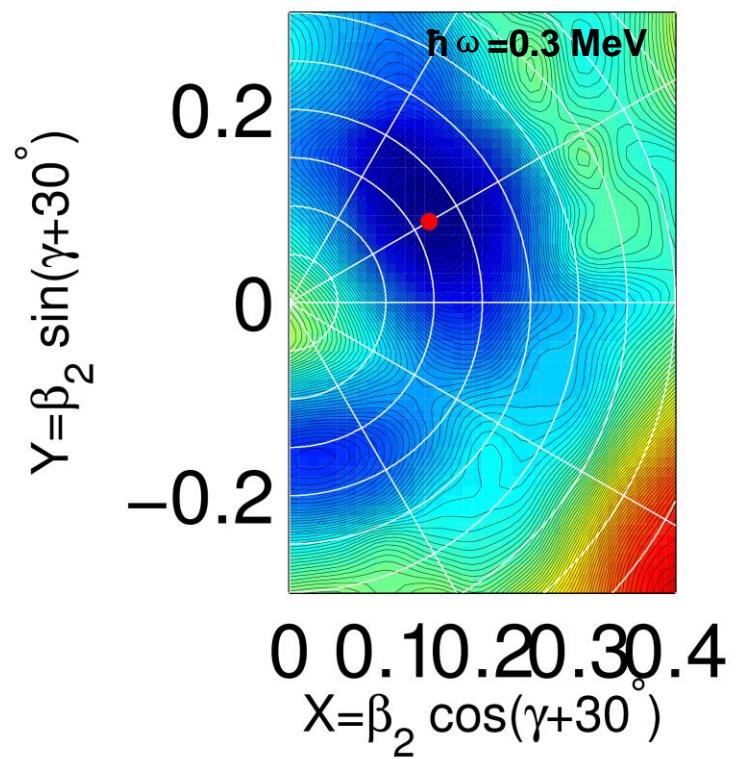
Semiclassical model of Dönau and Frauendorf

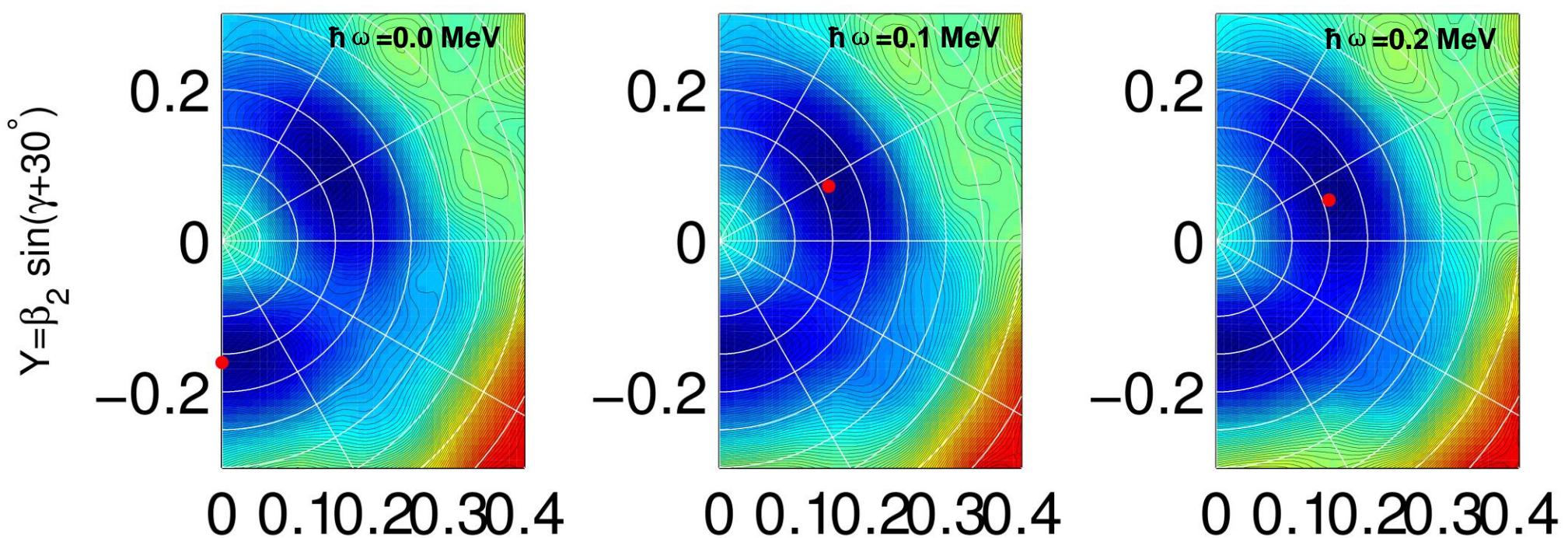
Exp.

$$\frac{B(M1; I \rightarrow I-1)}{B(E2 : I \rightarrow I-2)} = 0.697 \frac{1}{\lambda} \frac{E_\gamma^5(E2)}{E_\gamma^3(M1)} \frac{1}{1 + \delta^2} \left[\frac{\mu_N^2}{e^2 b^2} \right]$$

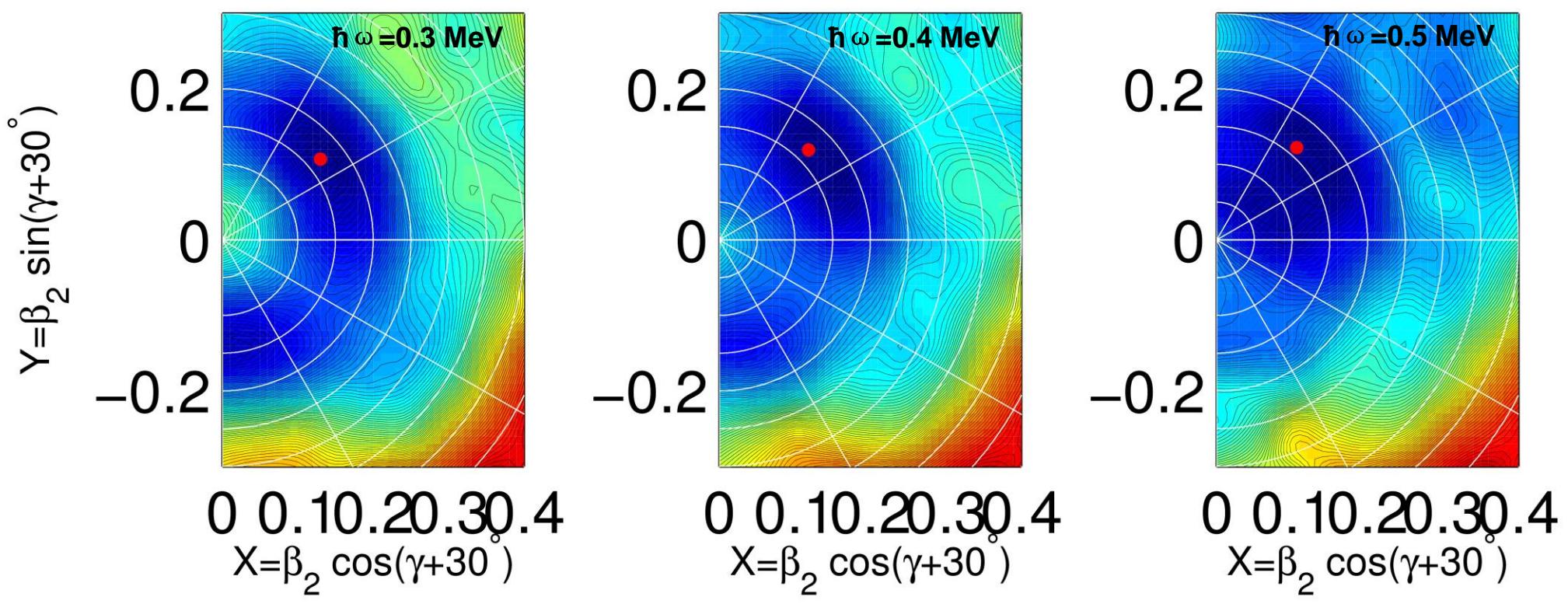


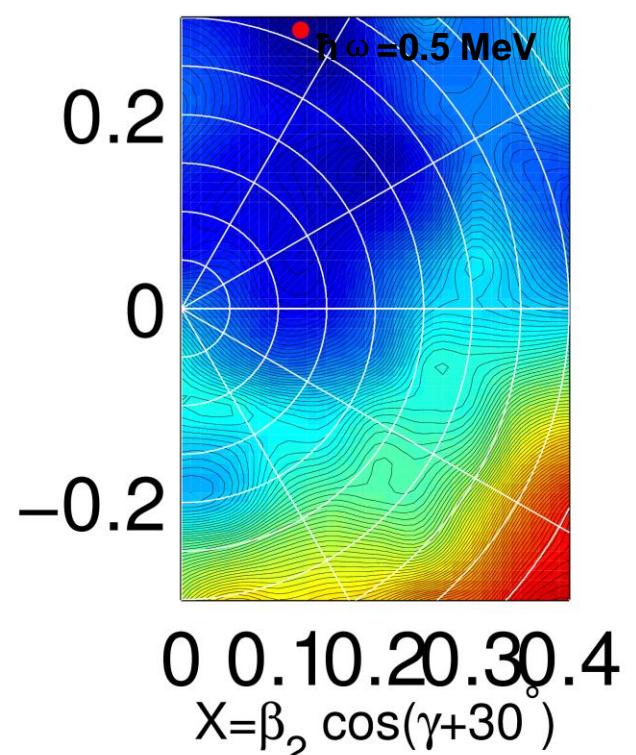
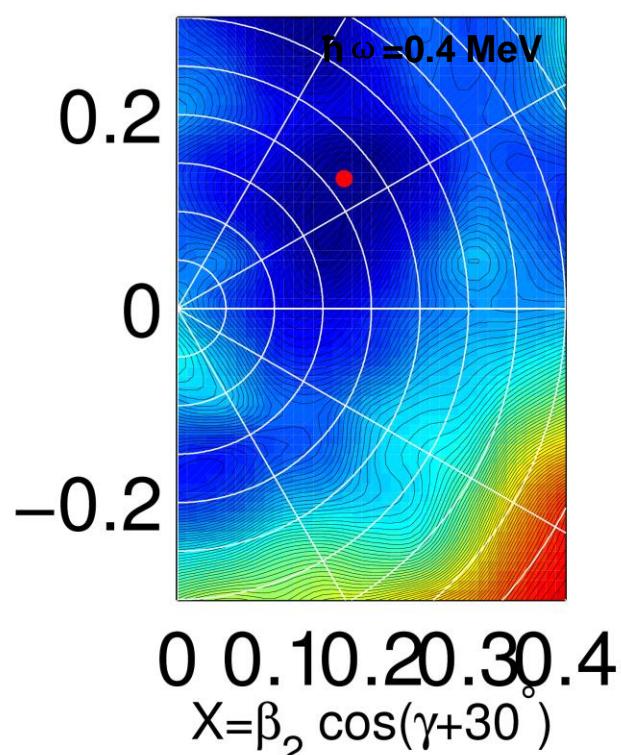
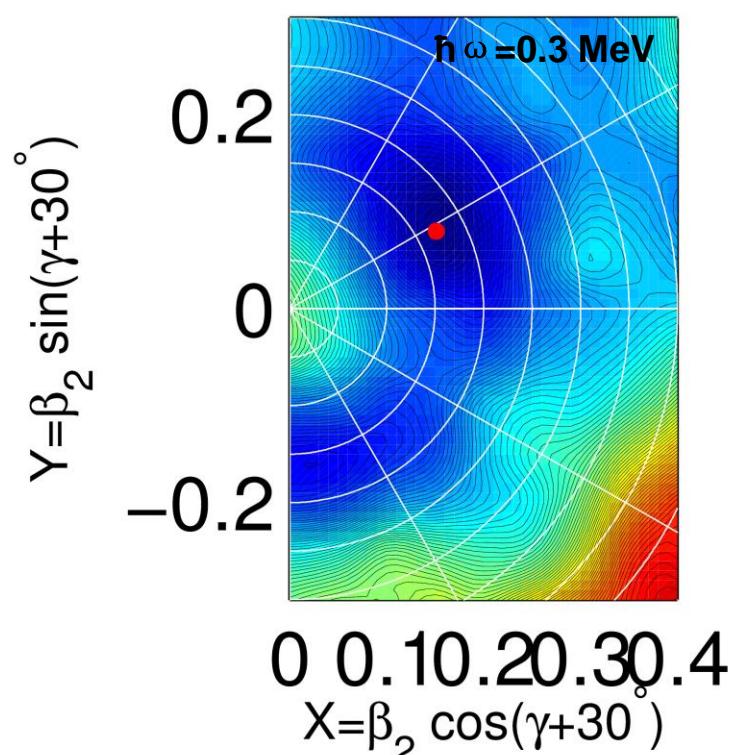
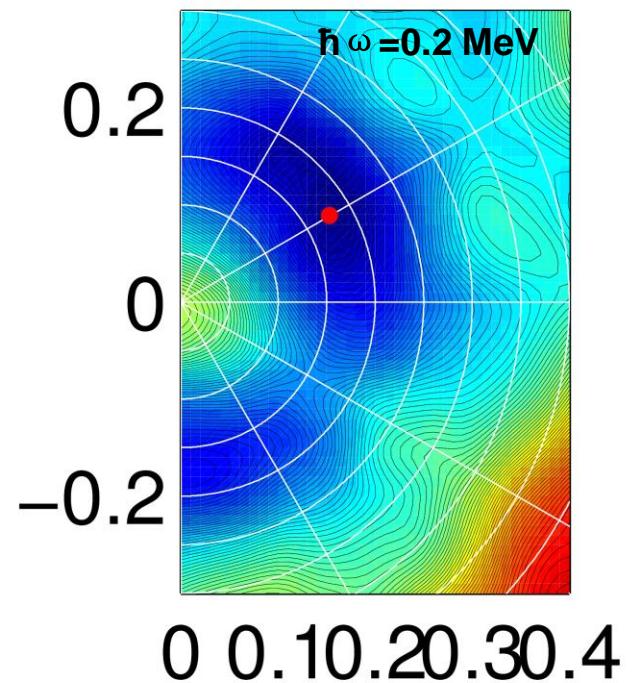
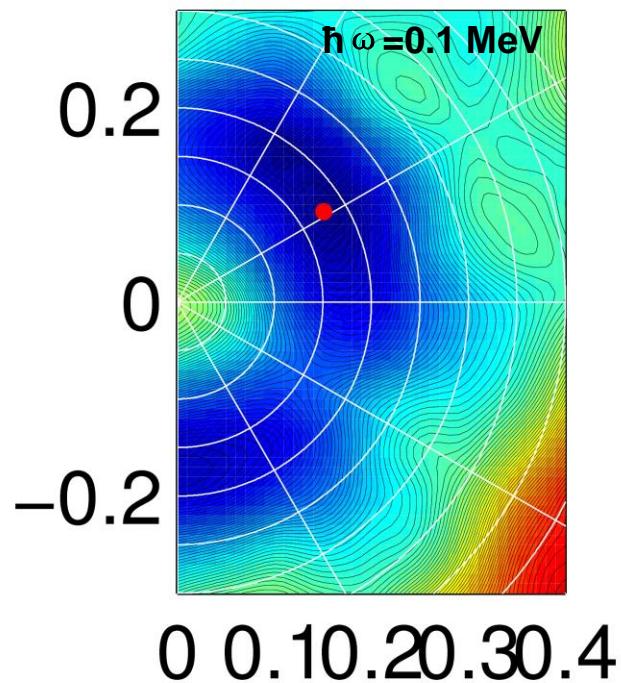
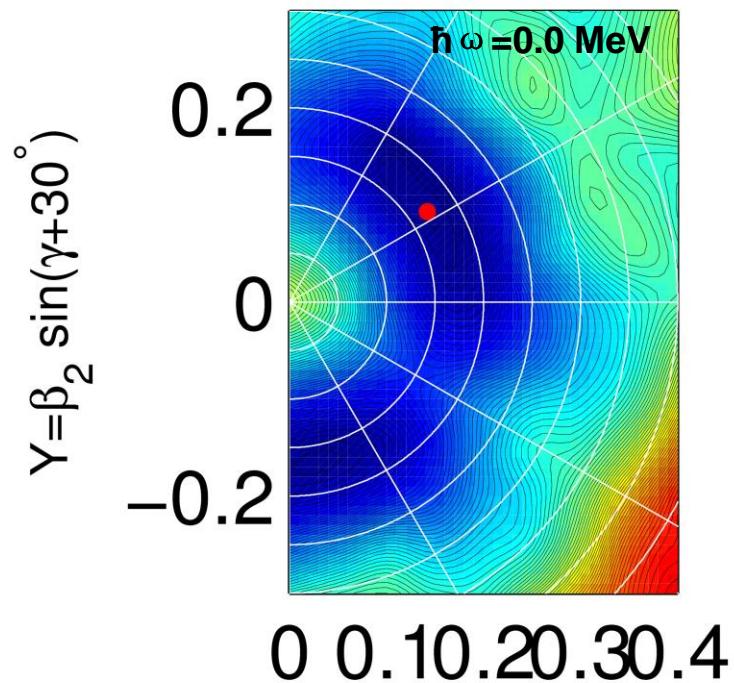
P: npps; N: ppps

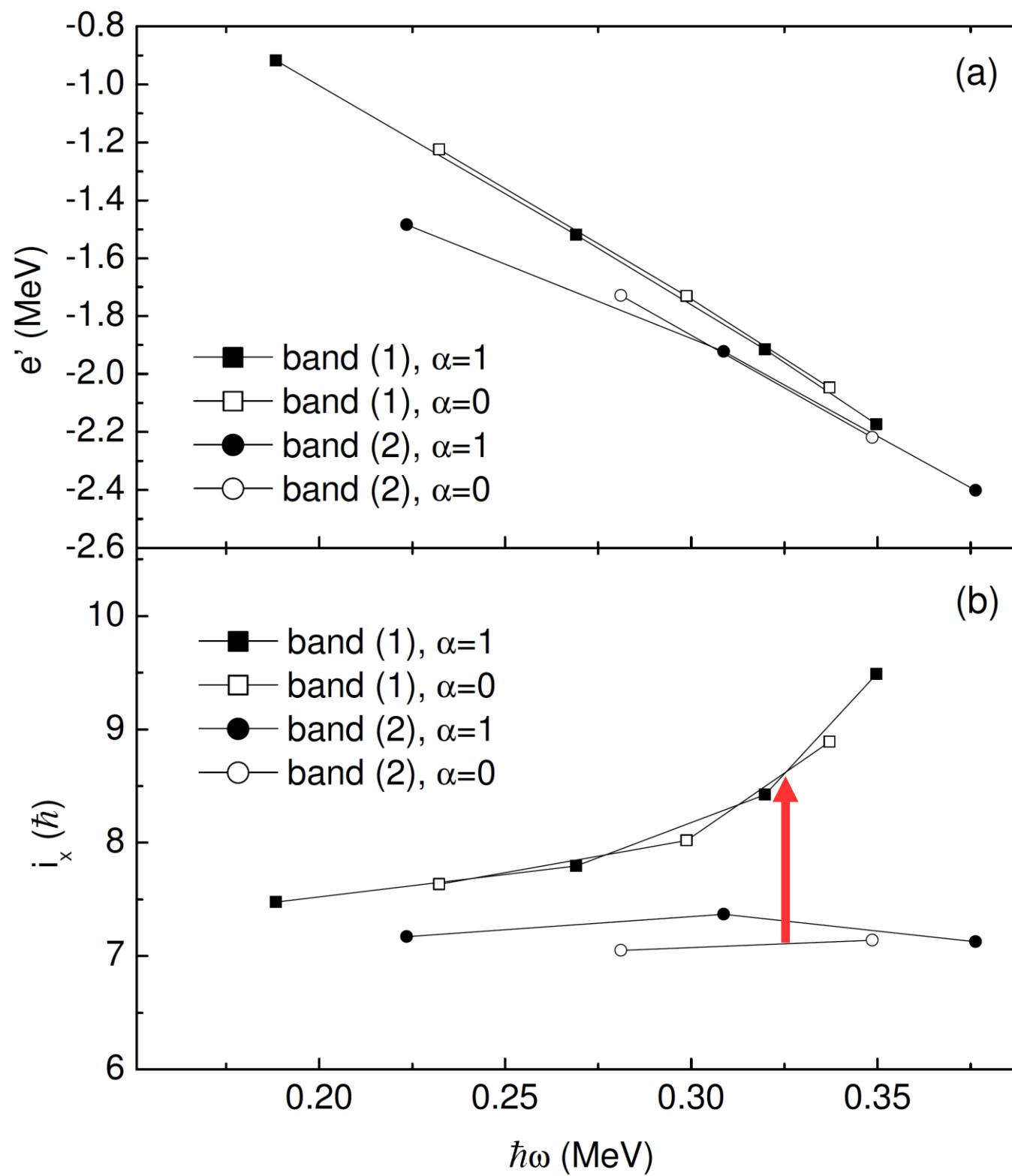




P: npns; N: npps







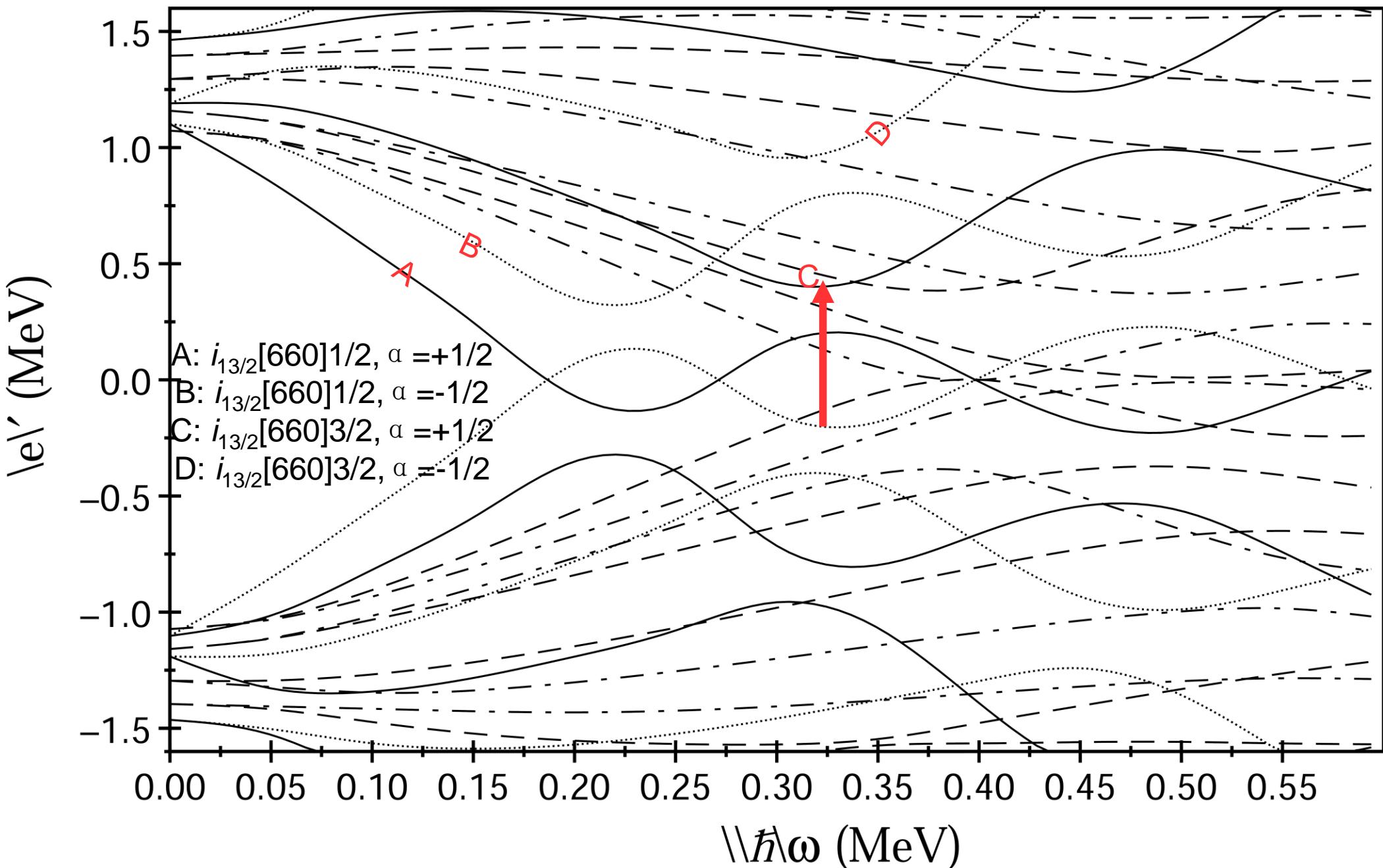
Harris
Parameters:
Band 1
 $J_0 = 13 \text{ } \hbar^2 \text{MeV}^{-1}$,
 $J_1 = 64 \text{ } \hbar^4 \text{MeV}^{-3}$

Band 2
 $J_0 = 4 \text{ } \hbar^2 \text{MeV}^{-1}$,
 $J_1 = 85 \text{ } \hbar^4 \text{MeV}^{-3}$

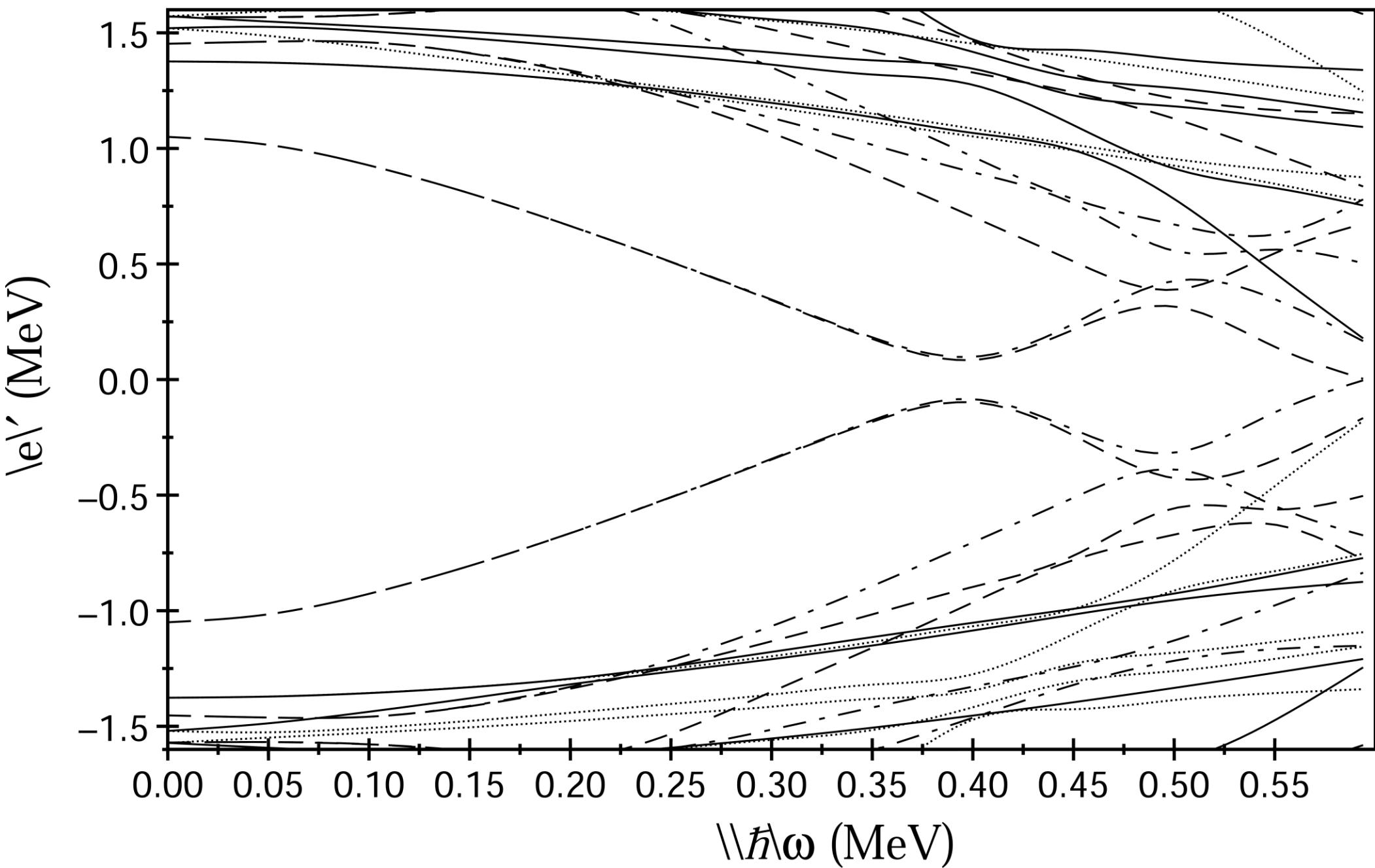
\Quasineutron levels :- Universal Woods-Saxon potential

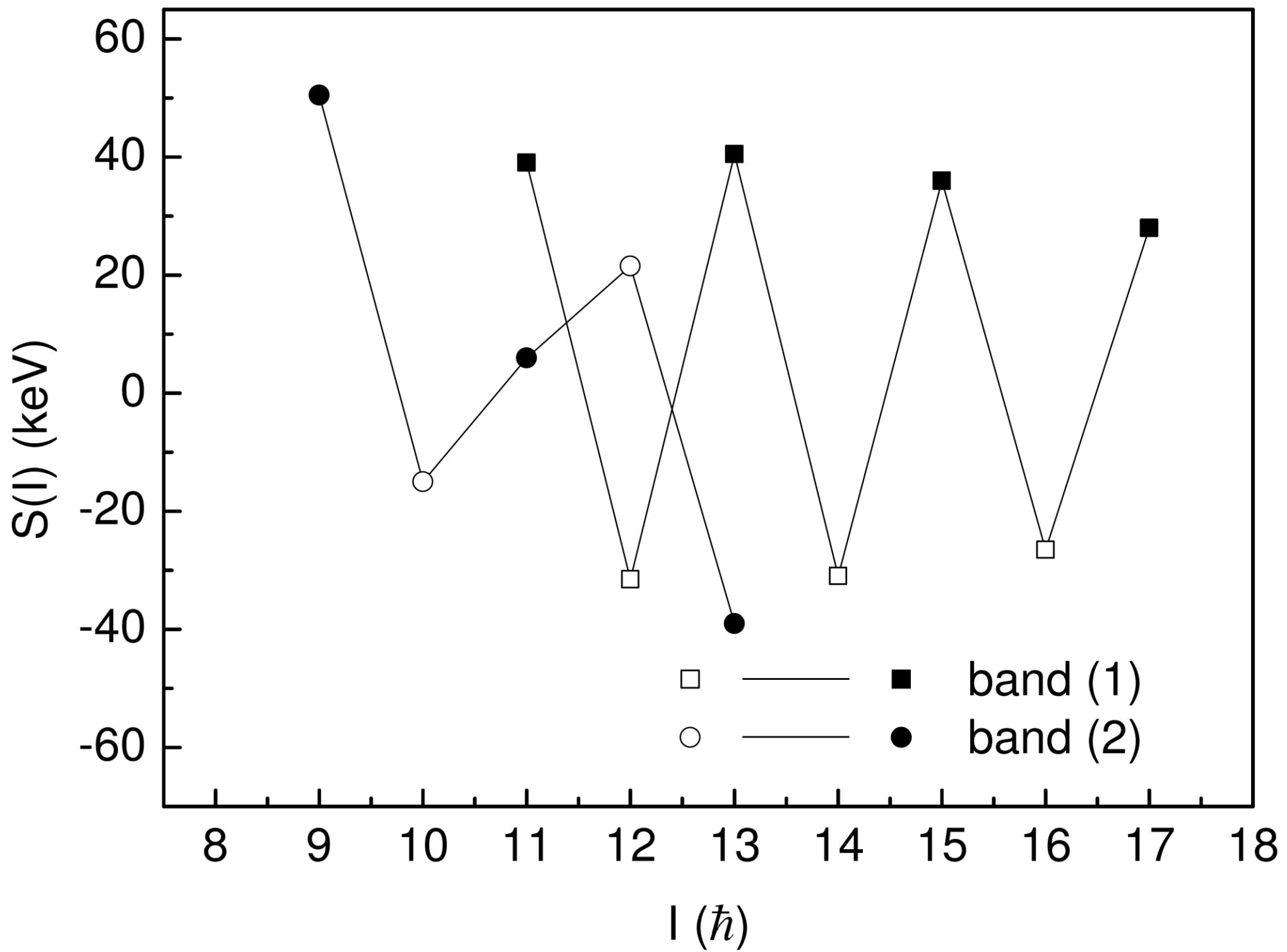
\N= 91, BETA2= 0.168, BETA4= 0.009, GAMMA= -1.6°, IMODEL= 2, DELTA0= 0.000

\(\langle\pi,\alpha\rangle\) : solid=\(+,+1/2\), dotted=\(+,-1/2\), dot-dash=\(-,+1/2\), dashed=\(-,-1/2\)



\Woods-Saxon Quasiproton levels : Universal Woods-Saxon potential
 $Z = 75$, $\text{BETA}2 = 0.168$, $\text{BETA}4 = 0.009$, $\text{GAMMA} = -1.6^\circ$, $\text{IMODEL} = 2$, $\text{DELTA}0 = 0.000$
 $(\pi, \alpha) :$ solid=(+, +1/2), dotted=(+, -1/2), dot-dash=(-, +1/2), dashed=(-, -1/2)





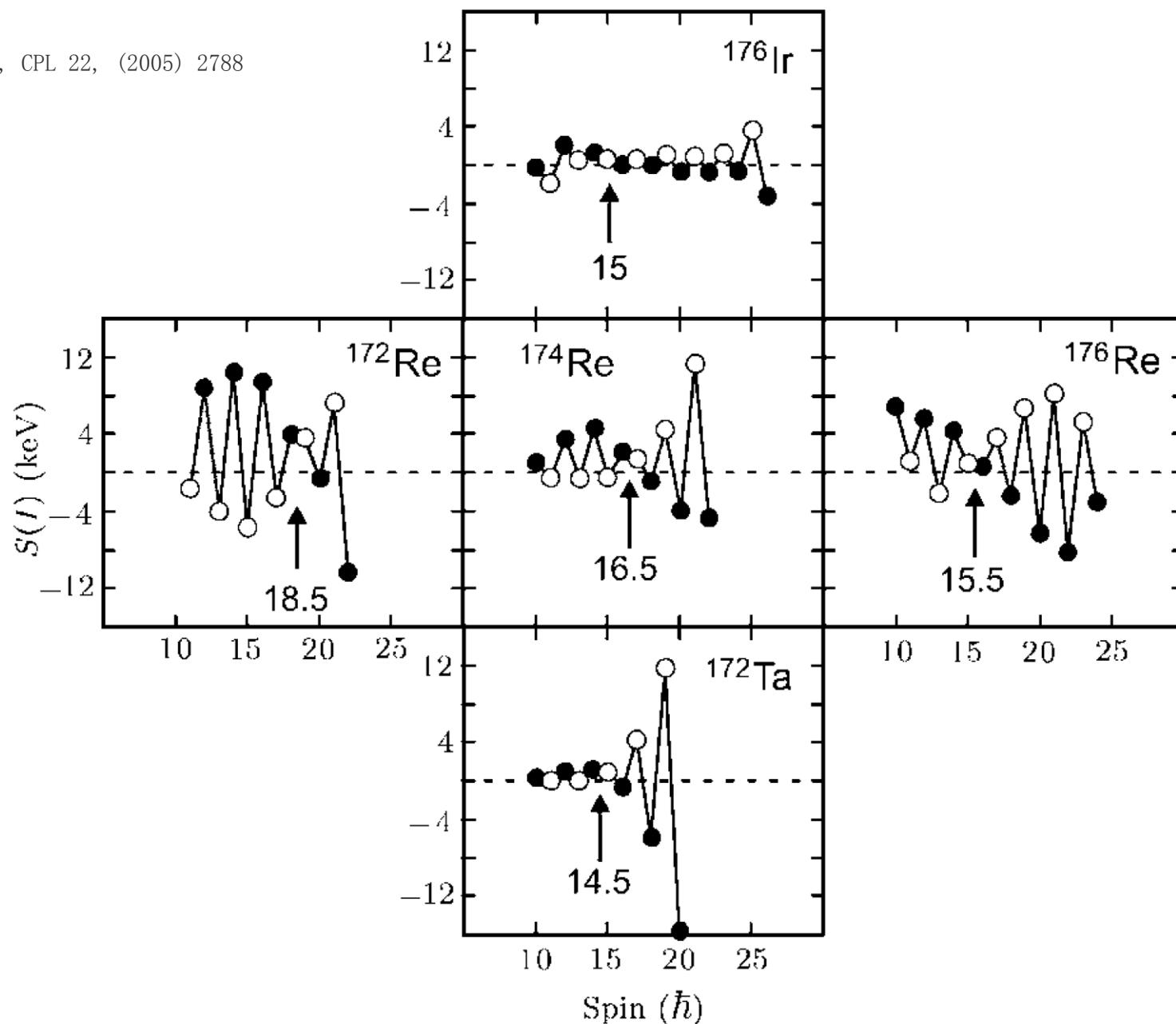


Fig. 4. Behaviour of signature splitting $S(I)$ versus I for band 1 and the $\pi h_{11/2} \otimes i_{13/2}$ bands in adjacent odd-odd nuclei. The arrow indicates the signature inversion spin.

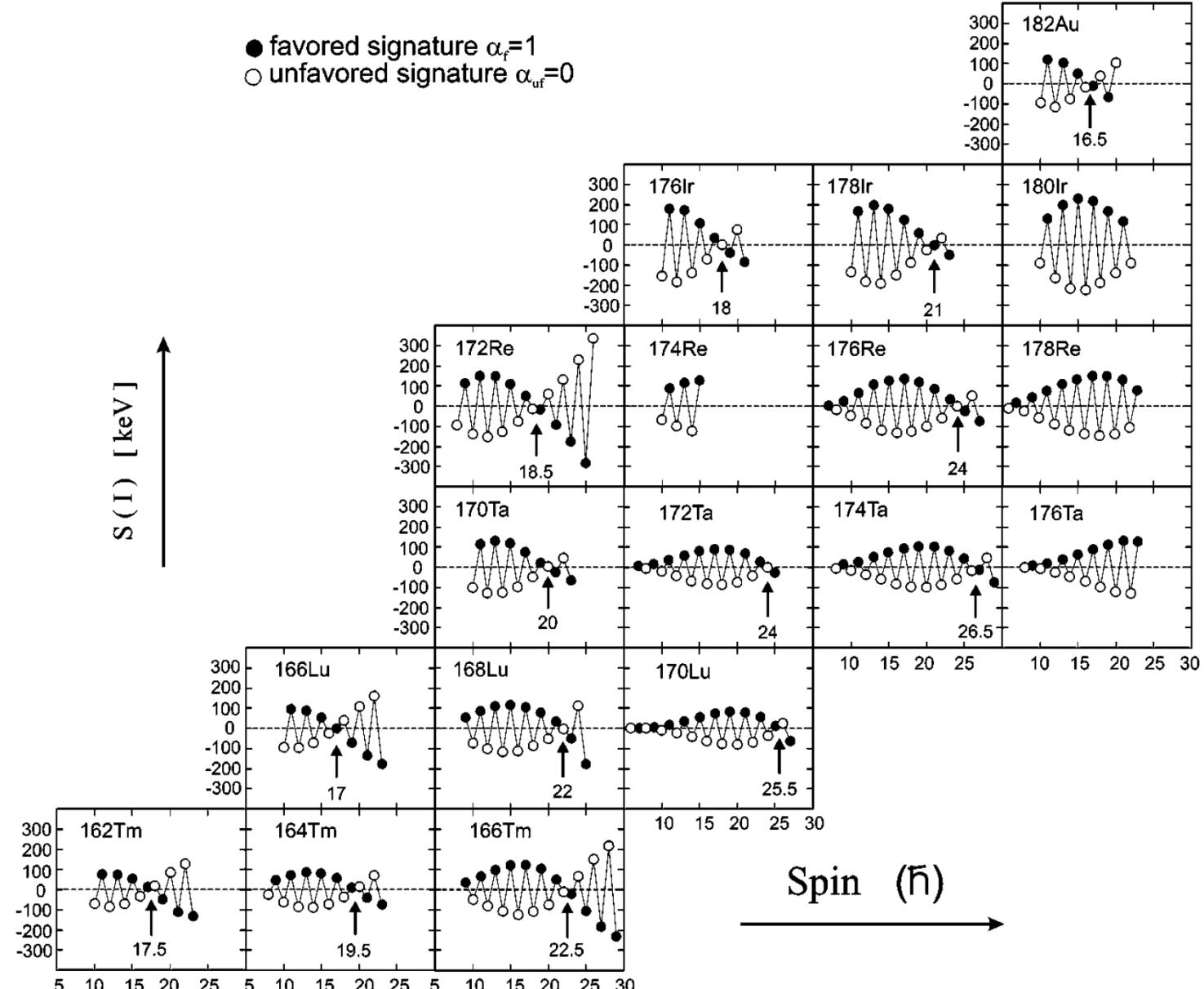


FIG. 9. A compilation of signature inversion for the $\pi h_{9/2} \otimes \nu i_{13/2}$ bands in $A=160\sim 180$ mass region. The filled (open) symbols correspond to the levels with favored signature $\alpha_f=1$ (unfavored signature $\alpha_{uf}=0$). The arrows indicate the signature crossing spins. The data sources are ^{182}Au [11], ^{176}Ir [10,49], ^{178}Ir [5,50], ^{180}Ir [32], ^{172}Re [this work], ^{174}Re [24], ^{176}Re [4], ^{178}Re [4,25], ^{170}Ta [41,44], ^{172}Ta [51], ^{174}Ta [13], ^{176}Ta [34], ^{166}Lu [52], ^{168}Lu [53,54], ^{170}Lu [43], $^{162,164}\text{Tm}$ [13], ^{166}Tm [55].

Conclusion

- ✓ First identification of two collective bands in odd-odd ^{166}Re
- ✓ The configurations for the two bands have been tentatively assigned and the deformation has been predicted by TRS calculations
- ✓ The backbending for band (1) may originate from the $i_{13/2}$ BC crossing
- ✓ Signature splitting and inversion was found in band (1), abnormal signature splitting was found in band (2).

Conclusion

- ✓ First identification of two collective bands in odd-odd ^{166}Re
- ✓ The configurations for the two bands have been tentatively assigned and the deformation has been predicted by TRS calculations
- ✓ The backbending for band (1) may originate from the $i_{13/2}$ BC crossing
- ✓ Signature splitting observed in bands (1) and (2) in agreement with TRS calc. (small triaxial deformation)
- ✓ Signature inversion observed in band (2). This phenomenon lacks consistent theoretical interpretation!!!

Thank you !