

<sup>176</sup>Yb(pol t,α) 1979Lo09

Type	Author	History	Citation	Literature Cutoff Date
Full Evaluation	M. Shamsuzzoha Basunia		NDS 102, 719 (2004)	1-Jun-2004

E=17 MeV with polarization of ≈78% target: 96.43% enriched <sup>176</sup>Yb measured α spectra with Q3D magnetic spectrometer. Typical energy resolution FWHM=24 keV.

Measured angular distribution and analyzing powers in the angular range θ=10°–50°. Relative uncertainties are ≈5%; absolute uncertainties are about 20%.

<sup>175</sup>Tm Levels

E(level)	J <sup>π</sup> †	T <sub>1/2</sub>	S <sup>‡</sup>	Comments
≈0 <sup>#</sup>	(1/2 <sup>+</sup> & 3/2 <sup>+</sup> )	15.2 min 5	≤0.70	Probable doublet comprised of J=1/2 and J=3/2 members of the 1/2[411] rotational band. A small energy spacing between these levels is expected for typical values of the decoupling parameter in this region. S: the upper limit in spectroscopic factor assumes all the cross section is due to the 3/2 member of the band. J <sup>π</sup> : g.s. 1/2 <sup>+</sup> assignment from systematics in <sup>169</sup> , <sup>171</sup> Tm T <sub>1/2</sub> : from Adopted Levels.
≈126 <sup>#</sup>	(5/2 <sup>+</sup> & 7/2 <sup>+</sup> )		≤0.72	Probable doublet. As for the ground state, the 5/2 and 7/2 members of the band are expected to lie close to each other. S: for J=7/2. S=0.12 for J=5/2.
353 <sup>#</sup>	(9/2 <sup>+</sup> )		0.11	
439 <sup>@</sup>	(7/2 <sup>-</sup> )		0.081	
532 <sup>@</sup>	(9/2 <sup>-</sup> )		0.13	
611 <sup>&amp;</sup>	(3/2 <sup>+</sup> )		0.11	
645 <sup>@</sup>	(11/2 <sup>-</sup> )		2.1	
672 <sup>&amp;</sup>	(5/2 <sup>+</sup> )		0.72	
756 <sup>&amp;</sup>	(7/2 <sup>+</sup> )		0.18	
870 <sup>&amp;</sup>	(9/2 <sup>+</sup> )		0.11	
941	(15/2 <sup>-</sup> )			Tentative assignment: 15/2 <sup>-</sup> , 7/2[523]. Level is probably populated through a multistep process.
1004				
1072				
1146	(7/2 <sup>+</sup> , 3/2 <sup>+</sup> )		0.16	S: for J=7/2. S=0.072 for J=3/2.
1212	(7/2 <sup>-</sup> , 5/2 <sup>+</sup> )		0.065	S: for J=7/2. S=0.066 for J=5/2.
1365	(7/2 <sup>+</sup> , 3/2 <sup>+</sup> )		0.096	S: for J=7/2. S=0.047 for J=3/2.
1423 <sup>a</sup>	(5/2 <sup>+</sup> )		0.23	
1495	(7/2 <sup>-</sup> , 5/2 <sup>+</sup> )		0.048	S: for J=7/2. S=0.056 for J=5/2.
1583				
1706 <sup>b</sup>	(5/2 <sup>+</sup> )		0.11	
1744				
1810 <sup>b</sup>	(7/2 <sup>+</sup> )		0.42	
1916 <sup>c</sup>	(7/2 <sup>-</sup> )		0.24	
1985	(5/2 <sup>+</sup> )		0.16	Energy systematics of the 1/2[420] Nilsson orbital in Ho isotopes suggests a 5/2 <sup>+</sup> , 1/2[420] assignment.
2013				
2056				
2095 <sup>c</sup>	(11/2 <sup>-</sup> )		1.3	
2216				
2310				

† Comparison with DWBA calculations permitted spin and parity assignments of several rotational bands. Weakly populated band

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 $^{176}\text{Yb}(\text{pol t}, \alpha)$  **1979Lo09 (continued)**

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 $^{175}\text{Tm}$  Levels (continued)

members assignments rely on the Nilsson model and are based on energy patterns and comparison with neighboring nuclei.

‡ Experimental values of nuclear structure factors are given by  $(1/(2N))(d\sigma/d\Omega(\text{exp})/d\sigma/d\Omega(\text{DWBA}))$  with  $N=23$ .

# 1/2(411) band.

@ 7/2(523) band.

& 3/2(411) band.

<sup>a</sup> 5/2(402) band: tentative.

<sup>b</sup> 5/2(413) band.

<sup>c</sup> 5/2(532) band.