

## Description and prediction of Long-lived Isomers in the $_{101}\text{Md}$ Isotopic Chain

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As a part of our continuing investigations on level structures of the odd-odd deformed nuclei, including the occurrence and characterization of long-lived isomers (LLI) therein, we have lately focused on nuclei of the heavy actinide region. In this report, we briefly present the results of such investigations for the even mass [A=246 (2) 258] isotopes of first transfermium species, namely Mendelevium (Z=101). A similar exercise on this topic was carried out over 30 years ago [1]. A factual summary of the first identifying experiments on each  $Z > 100$  nuclide (including all the Md isotopes) and the later experiments confirming their discovery, has been recently presented by Thoennessen [2]. The first identification of a Z=101 nuclide ( $^{256}\text{Md}$ ) was reported from Berkeley in 1955. Subsequently isotopes with A=254-258 and with A=248-253 were discovered [2] in 1970 and in 1973 respectively. However, even after 45 years, level structures (even for gs) remain elusive in each of these nuclei. Presently available data on LLI in even-A Md isotopes from the current data bases [3] is summarized in Table 1. A critical examination of these data leads us to following observations.

**Table 1:** ENSDF/XUNDL (August 2018) based listing of long-lived isomers in even-A (246-258)  $_{101}\text{Md}$  isotopes.

$^A\text{Md}$	$t_{1/2}$	$E_x$	$J^\pi$
246(a)	0.9 s	0+x	?
246(b)	4.4 s	0+y	?
248	7 s	0.0	?
250	52 s	0.0	(7 <sup>-</sup> )
252	2.3 m	0.0	(1 <sup>+</sup> )
254(a)	10 m	0+x	?
254(b)	28 m	0+y	?
256	77.7 m	0.0	(1 <sup>-</sup> )
258(a)	51.5 d	0	(8 <sup>-</sup> )
258(b)	57 m	0+x	(1 <sup>-</sup> )

a) In 5 (out of 10) known LLI, no  $J^\pi$  (and hence no configuration) has been mentioned.

b) For each of the other 5 LLI, a tentative  $J^\pi$  has been *suggested* from ‘analogy to (A-1) isotope/isotone and GM rule ordering’ [3], with no experimental or theoretical inputs.

c) LLI pairs are known in only 3 (out of 7) isotopes under consideration; the relative energy spacing is not known (or even estimated) in any of them. Whereas eight 2qp options are listed [3] for low lying levels in  $^{254}\text{Md}$ , for low-spin  $^{258}\text{Md}$  four different configurations and even a different  $J^\pi$  are suggested [4].

We present here the characterization of known isomer pairs, and prediction of LLI in other cases. Firstly we report our results on the characterization [6] of known isomer pair in  $^{254}\text{Md}$ . Next we present a detailed step wise calculation of level energies leading to prediction of higher-lying high-spin isomer in  $^{252}\text{Md}$ . Finally we present the case of  $^{250}\text{Md}$  wherein detailed calculations lead to a revision of the gs spin-parity.

We evaluate the bandhead energies of all the physically admissible 2qp GM doublets within specified energy domain using our well-tested 3-step Two Quasiparticle Rotor Model {TQRM} [5,6] employing the expression:

$$E(I, K) = E_0 + (E_p + E_n) + E_{rot} + \langle V_{np} \rangle;$$

$$E_{rot} \approx -\frac{\hbar^2}{2I} (2\Omega_{\leftarrow}) \delta_{K,K^-};$$

$$\langle V_{np} \rangle = -\left[\frac{1}{2} - \delta_{\Sigma,0}\right] E_{GM} + (-)^I E_N \delta_{K,0}. \quad (1)$$

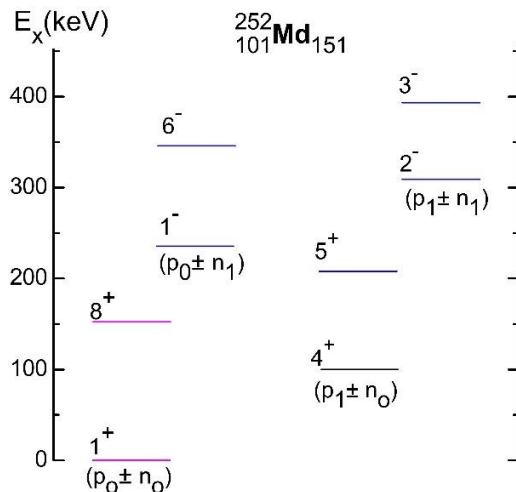
**$^{254}\text{Md}$ :** The bandhead energies of the physically admissible 2qp states were evaluated using our TQRM analysis. Critical examination of the results rules out high-spin assignment to both the isomers and instead leads us to the following assignments for the LLI pair with the 10 m isomer being designated as the gs of  $^{254}\text{Md}$  [6]:

10 m gs    1<sup>-</sup>0{p:1/2<sup>-</sup>[521]⊗n:1/2<sup>+</sup>[620]}  
28 m 0+x    3<sup>-</sup>3{p:7/2<sup>-</sup>[514]⊗n:1/2<sup>+</sup>[620]}

**<sup>252</sup>Md:** The latest data sheets (NDS) [3] lists only a 2.3 m isomer as <sup>252</sup>Md gs with possible  $1^+\{7/2-[514]\otimes n:9/2-[734]\}$  assignment. In Table 2, we list all the physically admissible 2qp bands for  $(E_p+E_n)\leq 350$  keV and the TQRM calculated energies of these 8 lowest <sup>252</sup>Md bands are plotted in Fig 1. Our analysis thus predicts the existence of a high-spin  $K^\pi=8^+$  LLI at  $\sim 150$  keV and possibly another shorter lived  $K^\pi=4^+$  isomer at  $\sim 100$  keV.

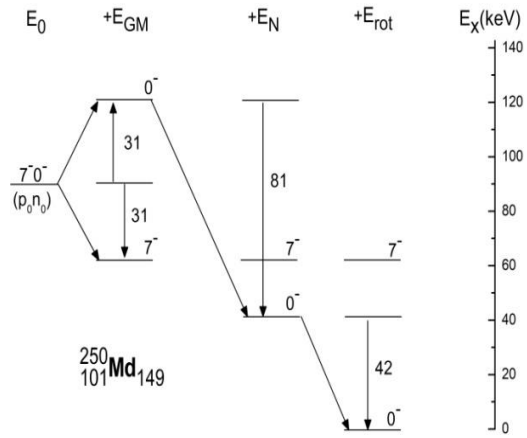
**Table 2:** Physically admissible 2qp GM doublet bands ( $K_T$  &  $K_S$ ) in <sup>252</sup>Md for  $(E_p+E_n)\leq 350$  keV.

$p_i$	$n_j$ $E_p$	$n_0$ 0		$n_1$ 200	
		$9/2-[734\uparrow]$	$8^+$	$5/2^+[622\uparrow]$	$6^-$
$p_0$	0	$1^+$	$8^+$	$1^-$	$6^-$
$7/2$	$[514\downarrow]$				
$p_1$	55	$4^+$	$5^+$	$2^-$	$3^-$
$1/2$	$[521\downarrow]$				



**Fig 1:** Calculated bandhead energies of the low-lying eight 2qp states in <sup>252</sup>Md using notations of Table 2.

**<sup>250</sup>Md:** As given in the latest NDS [4], straightforward coupling of the 101<sup>st</sup> p orbital  $7/2-[514]$  and the 149<sup>th</sup> n-orbital  $7/2^+[624]$  yields  $K^\pi=7^-$  &  $0^-$  as GM doublet bands, with GM rule placing spins-parallel  $K_T=7^-$  lower (and hence <sup>250</sup>Md gs). However, as evident from our Eq. (1), there are 2 additional energy contributions specifically for the  $K^\pi=0^-$  band. As detailed in



**Fig. 2:** The term-wise contribution to the bandhead energies of the gs GM pair showing the placement of  $K_S=0^-$  below the  $K_T=7^-$  state.

our reference [7] and illustrated in Fig. 2 here, these contributions bring the  $K^\pi=0^-$  level below the  $K_T=7^-$  level thus making the  $K^\pi=0^-$  as <sup>250</sup>Md gs. The  $K^\pi=7^-$  level appears as a higher-lying ( $\sim 60$  keV) high-spin longer-lived isomer.

Possible occurrence of isomer pairs in the other two Md isotopes, namely <sup>248</sup>Md and <sup>256</sup>Md are being investigated. For example, the latest (2017) NDS assigns  $I^\pi K=1^0\{7/2-[514]\otimes n:7/2^+[613]\}$  to the 78 m <sup>256</sup>Md gs with no mention of the location of its  $K^\pi=7^-$  GM counterpart. This surely constitutes a longer lived higher-lying high-spin state in <sup>256</sup>Md, a mention of which is made in NUBASE 2016 [3].

**References**

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