

Shell-model interpretation of the β^- decay of $^{212}\text{Bi}^g$, $^{212}\text{Bi}^{m1}$, and $^{212}\text{Bi}^{m2}$

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The β^- decay of the ^{212}Bi ground state ($T_{1/2} = 61$ min) and the isomers $^{212}\text{Bi}^{m1}$ ($T_{1/2} = 27$ min), and $^{212}\text{Bi}^{m2}$ ($T_{1/2} = 8$ min) are considered in the spherical shell model. The wave functions and energy spectra of all even and odd states in the Kuo-Herling model space were calculated for both ^{212}Bi and ^{212}Po . Possible candidates for the two ^{212}Bi isomers were located from these results and the allowed and first-forbidden β^- decays rates were formed using a recently developed parametrization of first-forbidden decays in the lead region. It was concluded that $^{212}\text{Bi}^{m1}$ is the yrast 8^- state and $^{212}\text{Bi}^{m2}$ the yrast 18^- state of ^{212}Bi . The decay rates calculated for these two β^- emitters are in excellent agreement with experiment. The predictions for the decay of the ground state of ^{212}Bi are not in good agreement with experiment, illustrating the difficulty of calculating weak first-forbidden decay modes reliably.

I. INTRODUCTION

^{212}Bi has three isomers which have been observed to decay at least in part by β^- emission leading to ^{212}Po . The ground state with $T_{1/2} = 61$ min, $Q(\beta^-) = 2252 \pm 4$ keV, and $J^\pi = 1^-$ (Ref. [1]) is designated $^{212}\text{Bi}^g$. An isomer, $^{212}\text{Bi}^{m1}$, with $T_{1/2} = 26.5 \pm 1.0$ min (Refs. [2,3]) and at an excitation energy E_x of 250 keV [2] has been speculated [2] to have $J^\pi = 9^-$. An 8 \pm 1-min activity, $^{212}\text{Bi}^{m2}$, with $E_x > 670$ keV decays by β^- (and possibly subsequent γ) emission to a ^{212}Po level at 2921 ± 15 keV which itself is isomeric with $T_{1/2} = 45.1 \pm 0.6$ s (Refs. [1-3]) and early on was speculated to have $J^\pi = 16^+$ (Ref. [4]) and 18^+ (Ref. [5]).

The existence of these three ^{212}Bi isomers and the ^{212}Po isomer has been known for some time yet the answers to the following questions have not been answered: (1) What is the situation which causes the isomerism of the 27- and 8-min activities? (2) Specifically, what are the spin-parity values of these two activities? (3) Can the β^- decay rates of the three ^{212}Bi activities be understood quantitatively? These three questions pose an interesting and challenging spectroscopic problem which is the subject of this study.

To address these questions shell-model calculations for both even- and odd-parity states were performed for both ^{212}Bi and ^{212}Po in the full Kuo-Herling particle space [6] which consists of the six proton orbits and seven neutron orbits above an inert (assumed) ^{208}Pb core. Beta-decay rates were then calculated using a recently developed parametrization of first-forbidden decays which is applicable to $A \sim 208$ nuclei [7].

II. RESULTS

A. Energy spectra

The calculations were performed with the shell-model code OXBASH [8]. The interaction used is a slightly mod-

ified version of the Kuo-Herling "bare + 1p1h" (KHP) interaction [6,9]. The modifications leading to the interaction KHP_e consisted mainly of changes to the proton-neutron interaction with the aim of reproducing the low-lying spectrum of ^{210}Bi . The changes and an appraisal of the interaction are described elsewhere [10].

A rather complete knowledge of the predicted energy spectra of ^{212}Bi and ^{212}Po is necessary for a thorough exploration of the questions posed in the Introduction. The necessary predictions are listed in Tables I and II. These results are used to construct the theoretical decays schemes shown in Fig. 1. The only experimental data included in the left two-thirds of the figure are the $T_{1/2}$ and β^- branching ratio for the ^{212}Bi isomers and the experimental $\log f_0 t$ values for ^{212g}Bi decay [1]. The experimental data on the right is that of Poletti *et al.* [11] which was obtained in a $^{208}\text{Pb}(^9\text{Be}, \alpha n \gamma)^{212}\text{Po}$ fusion-evaporation study. A similar $^{210}\text{Pb}(\alpha, 2n \gamma)^{212}\text{Po}$ study by Sugawara *et al.* [12] gave similar and consistent results. The unadorned J^π assignments in the experimental decays scheme are assumed here to be firm, although Poletti *et al.* and Sugawara *et al.* do not claim definite assignments for those level above 1500 keV. The J^π values in square brackets are speculations for J of Poletti *et al.* In four cases the shell-model calculation has been invoked to add an even-parity choice to these speculations. Rather obvious identifications with shell-model states can be inferred for all five of the speculated assignments by reference to Table II. The fusion-evaporation studies ended at the 14^+ state at 2885 MeV.

B. Calculation of the β^- decay rates

The $Q(\beta^-)$ value for a specific branch is taken as

$$Q(\beta^-) = 2252 + E_x(^{212}\text{Bi}) - E_x(^{212}\text{Po}) \text{ keV} \quad (1)$$

with the excitation energies taken from experiment when possible and from the KHP_e predictions otherwise. With the $Q(\beta^-)$ values pertaining for the three ^{212}Bi isomers,

the only conceivable β -decay branches are either Gamow-Teller ($\Delta J \leq 1$, $\pi_i \pi_f = +$) or first forbidden ($\Delta J \leq 2$, $\pi_i \pi_f = -$); and if first forbidden, $\Delta J \leq 1$. The latter restriction follows from the restriction on the strength of unique ($\Delta J = 2$) first-forbidden decays which is stated in the form [13]: *First-forbidden unique decays have $\log f_1 t \geq 8.5$.*

The Gamow-Teller (GT) beta moment is defined by [14]

$$B(\text{GT}) = 6166/f_0 t \text{ [GT(allowed) decay]}, \quad (2)$$

where f_0 is the Fermi integral and t the half-life of the decay branch in question [15,7]. For first-forbidden decays, comparison to experiment is made here via the averaged shape factor [16]

$$\overline{C(W)} = 9195 \times 10^5 / f_0 t = B_1^{(0)} + B_1^{(1)} \text{ fm}^2, \quad (3)$$

where advantage is taken of the fact that the decay is incoherent in the rank R of the contributing operators and, as stated above, the rank-two contribution is negligible. The rank-zero ($R0$) and rank-one ($R1$) components of first-forbidden decays can be formulated in terms of two $R0$ matrix elements M_0^S and M_0^T — the spacelike and timelike components of the axial current — and two

$R1$ matrix elements M_1^x and M_1^u — the $E1$ -like and $R1$ spin-dipole operators. Although the $R2$ contribution was calculated for all decay branches of interest, it was always negligible and need not be considered further. The $B_1^{(R)}$ were calculated using the full rigour and accuracy of the Behrens-Bühring formalism [16], however, the ζ approximation is useful for displaying the physics involved in the $R0$ and $R1$ decays [15,7]:

$$\begin{aligned} B_1^{(0)} &= [M_1^{(0)}]^2 = [\epsilon_{\text{mec}} M_0^T + a_S M_0^S]^2, \\ B_1^{(1)} &= [M_1^{(1)}]^2 = [a_u M_1^u - a_x M_1^x]^2, \end{aligned} \quad (\zeta \text{ approximation}) \quad (4)$$

where ϵ_{mec} is the meson-exchange-current (mec) enhancement factor evaluated for the lead region [7] as $\epsilon_{\text{mec}} = 2.01 \pm 0.05$. In the present calculations a value of 2.0 is used. The a_α are positive, largely kinematical and insensitive to nuclear structure. The ζ approximation has errors of $\leq 4\%$ for $R0$ and $\leq 10\%$ for $R1$.

The β matrix elements are calculated via

$$\begin{aligned} M_R^\alpha &= \sum_{j_i j_f} \mathcal{M}_R^\alpha(j_i j_f) = \sum_{j_i j_f} D_R(j_i j_f) M_R^\alpha(j_i j_f, \text{eff}) \\ &= \sum_{j_i j_f} D_R(j_i j_f) q_\alpha(j_i j_f) M_R^\alpha(j_i j_f) \end{aligned} \quad (5)$$

TABLE I. KHP_e predictions for the energy spectrum of ²¹²Bi. The index k orders states of a given J^π in energy. All levels are shown for $E_x < 1$ MeV while only yrast and yrare ($k = 1-2$) levels are shown for $E_x > 1$ MeV.

E_x (keV)	J^π	k	E_x (keV)	J^π	k	E_x (keV)	J^π	k	E_x (keV)	J^π	k
0	1 ⁻	1	911	6 ⁻	2	1586	6 ⁺	1	2519	16 ⁺	1
186	2 ⁻	1	936	4 ⁻	3	1599	7 ⁺	1	2543	18 ⁻	2
220	0 ⁻	1	968	4 ⁻	4	1615	6 ⁺	2	2561	18 ⁺	1
263	3 ⁻	1	991	6 ⁻	3	1620	5 ⁺	2	2568	21 ⁺	1
281	9 ⁻	1	1074	11 ⁻	1	1629	15 ⁻	2	2581	16 ⁺	2
303	8 ⁻	1	1080	9 ⁻	2	1633	10 ⁺	2	2586	0 ⁺	2
319	4 ⁻	1	1103	10 ⁻	2	1636	2 ⁺	2	2647	17 ⁺	2
345	6 ⁻	1	1188	11 ⁻	2	1637	7 ⁺	2	2669	19 ⁺	1
346	5 ⁻	1	1250	12 ⁻	1	1661	9 ⁺	1	2689	19 ⁺	2
347	1 ⁻	2	1261	0 ⁻	2	1664	8 ⁺	2	2712	20 ⁺	2
363	7 ⁻	1	1292	12 ⁺	1	1668	17 ⁻	1	2735	18 ⁺	2
436	10 ⁻	1	1293	10 ⁺	1	1680	12 ⁺	2	3079	21 ⁺	2
511	1 ⁻	3	1323	4 ⁺	1	1760	11 ⁺	2	3138	22 ⁺	2
654	2 ⁻	2	1325	11 ⁺	1	1812	1 ⁺	2	3285	24 ⁻	1
703	2 ⁻	3	1358	1 ⁺	1	1825	9 ⁺	2	3408	22 ⁻	1
708	3 ⁻	2	1367	13 ⁻	1	1905	13 ⁺	1	3595	23 ⁻	1
709	1 ⁻	4	1413	12 ⁻	2	2015	16 ⁻	2	3610	20 ⁻	1
731	7 ⁻	2	1429	2 ⁺	1	2143	13 ⁺	2	3683	24 ⁻	2
782	8 ⁻	2	1434	3 ⁺	2	2196	19 ⁻	1	3697	19 ⁻	2
784	1 ⁻	5	1461	15 ⁻	1	2204	17 ⁻	2	3698	21 ⁻	1
826	3 ⁻	3	1496	18 ⁻	1	2240	14 ⁺	1	3701	20 ⁻	2
835	5 ⁻	2	1500	5 ⁺	1	2276	0 ⁺	1	3877	22 ⁻	2
863	7 ⁻	3	1521	4 ⁺	2	2360	17 ⁺	1	3924	21 ⁻	2
865	3 ⁺	1	1522	13 ⁻	2	2368	15 ⁺	1	4022	23 ⁻	2
870	4 ⁻	2	1532	14 ⁻	1	2392	20 ⁺	1	5073	24 ⁺	1
879	2 ⁻	4	1547	14 ⁻	2	2406	14 ⁺	2	5096	23 ⁺	1
899	5 ⁻	3	1557	8 ⁺	1	2439	22 ⁺	1	5192	23 ⁺	2
904	8 ⁻	3	1581	16 ⁻	1	2510	15 ⁺	2	5329	24 ⁺	2

using Woods-Saxon (WS) wave functions. In Eq. (5) α labels a matrix element of rank R , $M_R^\alpha(j_i j_f)$ is a single-particle matrix element for the transition $j_i \rightarrow j_f$ in the impulse approximation, and the quenching factor $q_\alpha(j_i j_f)$ corrects $M_R^\alpha(j_i j_f)$ for the finite size of the model space and some effects of the nuclear medium. The $D_R(j_i j_f)$ are the one-body-transition densities which are the result of the shell-model calculations performed with the code OXBASH [8]. The quenching factor $q_{GT}(j_i j_f)$ for allowed decays was taken to be state independent and equal to 0.66 which is a value suggested by results obtained by Towner [17]. For first-forbidden decays recent results [18] were used which are associated with the parametrization for $A = 205$ – 212 nuclei [7] already alluded to.

With Fig. 1 as a reference we are now ready to consider the three ^{212}Bi activities. The information given in Fig. 1 and not as yet touched on will become clear as the discussions develop. We start with the 27-min $^{212}\text{Bi}^{m2}$ activity.

C. $^{212}\text{Bi}^{m2}$ decay

As pointed out by Poletti *et al.* [11] an isomer in either ^{212}Bi or ^{212}Po with a half-life in the 10-min region must in all probability be one which has no available

states to decay to by electromagnetic radiation with $L < 4$. From Fig. 1 it is clear that there are no predicted ^{212}Bi states with $E_x > 670$ keV which so qualify. In order to explain the $^{212}\text{Bi}^{m2}$ activity (with $E_x > 670$ keV) it is necessary to assume that either the ordering of the levels differs from the KHP_e predictions or that a highly unlikely retardation of one or more $L < 4$ electromagnetic transitions is occurring. The former alternative is considered much more likely on general grounds and calculations of electromagnetic transition strengths with the KHP_e wave functions supports this view. There are three obvious possibilities for isomers. The yrast 15^- and 18^- states are predicted to be separated by only 35 keV. This is well within the rms deviation of 62 keV for the ten states of ^{212}Po for which a correspondence between theory and experiment is indicated in Fig. 1. With the 18_1^- state below that of 15_1^- , it would have only $L \geq 5$ transitions available to lower ^{212}Bi states. A second possibility is that the 22_1^+ state lies below the 19_1^- and 20_1^+ states in which case all decays of the 22_1^+ state to lower states would have $L \geq 4$. This is considered less likely because the energy separation of the 19_1^- and 22_1^+ states is predicted to be 243 keV. Likewise, if the 15_1^- state were lowered by 211 keV relative to the $J^\pi = 12^+$, 12^- , and 13^- states shown in Fig. 1, it would have only $L \geq 4$ electromagnetic transitions available. These two pos-

TABLE II. KHP_e predictions for the energy spectrum of ^{212}Po . The index k orders states of a given J^π in energy. Only yrast and yrare ($k = 1$ – 2) levels are shown with the exception of some levels with $k > 2$ which are of present interest.

E_x (keV)	J^π	k	E_x (keV)	J^π	k	E_x (keV)	J^π	k	E_x (keV)	J^π	k
0	0^+	1	2422	7^+	2	2942	7^-	2	3822	16^-	2
788	2^+	1	2443	8^+	5	2955	12^+	2	3841	18^-	2
1172	4^+	1	2465	7^+	3	2969	8^-	2	3842	20^-	1
1356	6^+	1	2474	10^+	2	3032	16^+	2	3849	15^-	2
1405	8^+	1	2477	8^+	6	3039	13^+	1	3877	17^-	2
1429	2^+	2	2494	7^+	4	3056	1^-	1	3958	21^-	2
1515	1^+	1	2515	3^-	2	3083	14^+	2	3964	20^-	2
1543	0^+	2	2527	9^+	2	3098	21^-	1	4103	23^+	1
1669	4^+	2	2549	9^+	3	3102	13^+	2	4135	24^+	1
1727	6^+	2	2575	2^-	2	3124	18^+	2	4421	22^+	1
1729	8^+	2	2586	7^-	1	3138	12^-	2	4718	21^+	1
1732	10^+	1	2607	12^-	1	3149	15^+	1	4751	20^+	1
1858	3^+	1	2625	6^-	1	3187	15^+	2	4847	20^+	2
1904	8^+	3	2650	9^-	1	3204	1^-	2	4883	21^+	2
1952	3^-	1	2659	13^-	1	3245	17^+	1	4887	22^+	2
2003	4^+	3	2665	11^-	2	3263	19^-	1	4900	19^+	1
2084	6^+	3	2714	4^-	2	3303	13^-	2	4994	19^+	2
2089	5^+	1	2753	12^+	1	3375	17^+	2	5049	23^+	2
2111	1^+	2	2780	8^-	1	3492	17^-	1	5687	25^-	1
2164	8^+	4	2796	11^+	1	3509	15^-	1	6044	24^-	1
2172	7^+	1	2815	10^-	1	3540	14^-	1	6146	23^-	1
2190	9^+	1	2850	5^-	2	3599	18^-	1	6251	23^-	2
2223	4^+	4	2855	14^+	1	3632	16^-	1	6287	24^-	2
2282	11^-	1	2873	16^+	1	3653	18^+	3	6376	22^-	1
2287	2^-	1	2877	9^-	2	3714	0^-	1	6490	22^-	2
2289	3^+	2	2912	11^+	2	3719	19^-	2	6666	25^-	2
2307	4^-	1	2913	18^+	1	3773	17^+	3	7745	26^+	1
2403	5^-	1	2919	6^-	2	3784	0^-	2	8020	24^+	2
2411	5^+	2	2931	10^-	2	3802	14^-	2	8417	25^+	1

sibilities are not considered further because, as it turns out, the 18_1^- possibility provides a natural explanation for the activity [19]. The $J = 17, 18,$ and 19 states which are predicted to be energetically available for the β^- decay of the ^{212}Bi 18_1^- level are shown in Fig. 1.

The $J^\pi = ^{212}\text{Bi}$ 18_1^- state is found to be 98% $\pi 0h_{9/2}\nu 1g_{9/2}^2 0i_{11/2}$. First consider the first-forbidden decays of this state. There are two 17^+ and two 18^+ ^{212}Po states energetically available. The rank-one ($R1$) decays to all four are weak due to cancellation between the contributions of M_1^u and M_1^x to the $R1$ beta moment — see Eq. (4). This destructive interference is characteristic of $R1$ decays for $A = 209\text{--}212$ nuclei [7]. In the present instance it is due to the fact that the relative phases of the $\nu 1g_{9/2} \rightarrow \pi h_{9/2}$ and $\nu 1g_{9/2} \rightarrow \pi 1f_{7/2}$ single-particle $R1$ transitions are both even. The two KHP_e 18^+ states are largely an orthogonal mixture of $\pi 0h_{9/2}^2 \nu 1g_{9/2} 0i_{11/2}$ and $\pi 0h_{9/2} 1f_{7/2} \nu 1g_{9/2} 0i_{11/2}$, respectively. The $R1$ contributions to the two decays is predicted to be only 2.3%

and 0.8% of the $R0$ decays. There are two contributions to the $R0$ matrix elements $\nu 1g_{9/2} \rightarrow \pi h_{9/2}$ and $\nu 1g_{7/2} \rightarrow \pi 1f_{7/2}$. These are in phase for 18_1^+ and out of phase for 18_2^+ with the result that the $\log f_0 t$ values for the decay to these two states are 5.68 and 6.38, respectively, and — using the KHP_e predictions for the relative energies — β^- branching ratios to these two states of 7.8% and 0.6% are calculated for $T_{1/2} = 8$ min.

Now consider the allowed decays. The $J^\pi = 17^-, 18^-,$ and 19^- yrast states are predicted to be 80%, 86%, and 86% $\pi 0h_{9/2} 0i_{13/2} \nu 1g_{9/2}^2$, respectively. The Gamow-Teller beta moment $B(\text{GT})$ is calculated to be 0.0009, 0.0039, and 1.15 for these three final states so that the decays to the 17^- and 18^- states are predicted to be negligible compared to that to the 19^- state. The much larger value of $B(\text{GT})$ for the 19^- state is due to its being a “stretched” state as is the ^{212}Bi 18_1^- state. Thus for the transition between these two states the coupling coefficients associated with the only contributing single-particle transition, $\nu 0i_{11/2} \rightarrow \pi 0i_{13/2}$, are maximal. The f_0 value for the transition to the 19_1^- state is calculated to be 12.3 giving a predicted partial half-life of 440 s. This is in excellent agreement with the value of 520 s expected for a 92% branch (with 8% going via first-forbidden decays). Note that it would only take a 30-keV decrease in the relative excitation energies of the initial and final states to lower f_0 enough so as to exactly reproduce the $T_{1/2} = 520\text{-s}$ expectation.

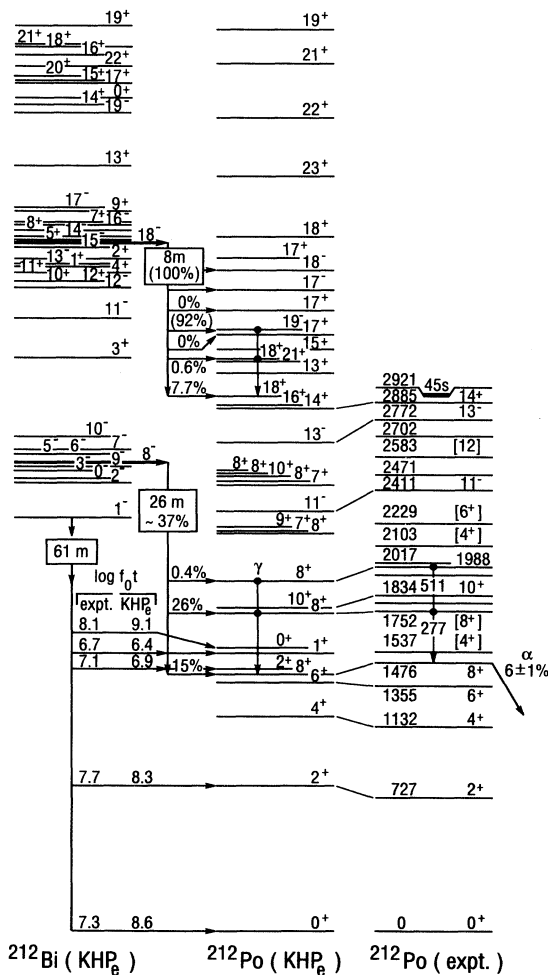


FIG. 1. Selected energy levels of ^{212}Bi and ^{212}Po and schematic of the β^- decays of ^{212}Bi . See the text for details.

D. $^{212}\text{Bi}^{m1}$ decay

From the relative intensities observed for the α decay of the ^{212}Bi 27-min isomer and the β^- -delayed α emission of the ^{212}Po 1476-keV 8_1^+ state [2], and with the aid of the α branching ratio for the 8_1^+ state of $6 \pm 1\%$ (Ref. [20]), a β^- branching ratio of $\sim 37\%$ can be deduced for the 27-min isomer [21]. In their study of this isomer, Lemmert *et al.* [3] observed the γ -ray cascades corresponding to $8_1^+ \rightarrow 6_1^+ \rightarrow 4_1^+ \rightarrow 2_1^+ \rightarrow 0_1^+$ in coincidence with the α decay of the ^{212}Po ground state (see Fig. 1). They also observed a 275-keV γ transition which is quite probably the same as the 277-keV $1752 \rightarrow 1476$ transition observed by Poletti *et al.* [11]. Since the 1834-keV level is identified as the yrast 10^+ state, the only reasonable identification of the 1752-keV level with the KHP_e states of Fig. 1 is with the 8_2^+ state.

There are two obvious possibilities for a 27-min isomer at $E_x = 250$ keV in ^{212}Bi ; the yrast 9^- or 8^- state, depending on which lies lower. The 9^- possibility was the only one suggested previously. The reason for this preference was that analogy was made to ^{210}Bi and in that nucleus the yrast 9^- and 8^- states lie at 271 and 583 keV, respectively. The compression of the spectrum of low-lying states of the even Bi isotopes as A increases from 210 to 214 has been remarked on previously [22]. This is another illustration. The odds as to which of the two states lies lowest seems about even. The KHP_e prediction is that the 9^- state is lowest by 22 keV. However, there are actually four 8^- states predicted below the second 9^- state and it would take only trivial changes in

the interactions between them to lower the 8_1^- state by 22 keV. The decays of both the 9^- and 8^- possibilities were considered. Note that both states have only first-forbidden decay channels available.

The predicted β^- decay of the 9^- state was found to be in severe disagreement with the experimental facts. First of all, the major β^- branch (4.9%) is predicted to be that to the 10_1^+ state and this decay mode would be accompanied by a 358-keV γ ray which was not observed experimentally [3]. The remaining β^- branching ratios add up to 0.24% and so the disagreement with experiment ($\sim 37\%$ β^- branching) is considerable for any branch which could be associated with a 275-keV γ -ray transition. The weakness of the $R1$ contributions to the decays is due to the same sort of destructive interference as seen for the 8-min isomer. The weakness of the $R0$ contribution to the $9_1^- \rightarrow 9_1^+$ transition is due to very small values for the relevant $D_0(j_i j_f)$.

The predicted β^- decay of the ^{212}Bi 8_1^- state is a quite different story. The $R1$ decays to the available 7^+ and 9^+ states exhibit the same destructive interference between M_1^x and M_1^u as the other decays considered so far. The decay branches to these states and to the 8_4^+ and 8_5^+ states are predicted to be negligible ($< 0.2\%$ in total). However the decays to the three lowest 8^+ states are predicted to sum to a β^- branching ratio of 41% in nice agreement with the experimental value of $\sim 37\%$. The predicted $\log f_0 t$ values for these three decays are 6.23, 5.51, and 6.76. They are, respectively, 99%, 99%, and 94% $R0$. For all three branches the $R0$ contribution from $\nu 1g_{9/2} \rightarrow \pi 0h_{9/2}$ is dominant. For 8_1^+ and 8_2^+ all five possible single-particle transitions contribute in phase. Thus, these two states share the role of the ‘‘pygmy’’ resonance of the initial state as recently discussed [7] (also see Sec. III). As shown in Fig. 1, the predicted branching ratio of 26% into the 8_2^+ state provides an explanation for the 275-keV γ -ray transition observed by Lemmert *et al.* [3].

E. $^{212}\text{Bi}^g$ decay

Beta decay of the ^{212}Bi ground state to the first two 0^+ and 2^+ states and the first 1^+ state was considered. The resulting $\log f_0 t$ values are compared to experiment in Fig. 1. The agreement is seen to be very poor. This is not surprising. All five transitions are predominantly $R1$ and with one exception the overlap of the initial and final states is very poor, i.e., the $D_0(j_i j_f)$ are small. In addition, the contributions of M_1^x and M_1^u are destructive. The one exception is the decay to 2_2^+ which has a strong $\nu 1g_{9/2} \rightarrow \pi 0h_{9/2}$ component and is dominated by M_1^u . It is gratifying that the agreement for the latter decay is reasonably good. There are two general reasons why $R1$ decays with $\log f_0 t \gtrsim 7.0$ can often not be predicted well. First, for such small rates the results are usually very sensitive to the wave functions — often to a point beyond which the wave functions can be expected to give meaningful results. Second, at this level the next order of terms in the Behrens-Bühring expansion begin to become non-negligible [16]. Note that all the first-

forbidden decays considered for the decay of $^{212}\text{Bi}^{m1}$ and $^{212}\text{Bi}^{m2}$ have $\log f_0 t < 6.8$. In summary we should be satisfied that the three weakest decays are predicted as such and the agreement is not too bad for the two strongest branches.

III. SUMMARY

It was expected that the Kuo-Herling model space for $Z \geq 82$, $A > 208$ will give a good description of the yrast and yrare states of ^{212}Bi and ^{212}Po and that an examination of the spectra of these states would reveal the possible candidates for $^{212}\text{Bi}^{m1}$ and $^{212}\text{Bi}^{m2}$. Accordingly the spectra were calculated and indeed several possibilities were obvious for each isomer. Calculations of the β^- decay rates were then performed to test the proposed candidates and obvious choices for both isomers were found. The success of this approach is due to two factors: (1) The availability of shell-model results for all ^{212}Bi and ^{212}Po states of the Kuo-Herling model space — heretofore only the even-parity spectrum of ^{212}Po was known [23]. (2) The ability to calculate first-forbidden decay rates with good accuracy [7].

The three isomers studied exhibit quite different decay properties which can be characterized by the proposed principal decay mode. The 8-min high-spin isomer $^{212}\text{Bi}^{m2}$ is predicted to decay by an allowed transition to the yrast 19^- state of ^{212}Po . The transition is quite fast due to the stretched nature of the ^{212}Bi 18_1^- and ^{212}Po 19_1^- states. It is proposed that the 19_1^- state then decays by an $E1$ transition to the 18_1^+ state of ^{212}Po as indicated in Fig. 1. Arguments and experimental evidence advanced by Poletti *et al.* [11] and Kudo *et al.* [24] strongly support the identification of the 18_1^+ state as the 45-s isomer of ^{212}Po . (However, it would be reassuring to see evidence for formation of the ^{212}Po 16_1^+ and 18_1^+ states in a fusion-evaporation study such as that of Poletti *et al.* [11].) If the ^{212}Po isomer is actually the 16_1^+ state, then the proposed decay of the 19_1^- state would be $19_1^- \xrightarrow{E1} 18_1^+ \xrightarrow{E2} 16_1^+$. In Table III are listed KHP_e predictions for some $E2$ and $E4$ transitions. These are given because of their potential use in obtaining an understanding of the ^{212}Po states involved in the decay of the 18_1^+ state. They also display adequate agreement with experiment with reasonable values of the effective charges.

The 27-min $^{212}\text{Bi}^{m1}$ isomer at an excitation energy of 250 keV is predicted to be the 8_1^- state and to decay mainly by $R0$ first-forbidden components to the first two 8^+ states of ^{212}Po . As discussed in detail previously [7], the initial (final) state in first-forbidden beta decay has $R0$ and $R1$ particle-hole giant resonances at high excitation ($\sim 1\hbar\omega$) in the final (initial) nucleus. This is well known and follows from the repulsive nature of the particle-hole interaction. What is less well known is that the attractive particle-particle interaction gives rise to another (albeit weaker) resonance (termed ‘‘pygmy’’ [7]) situated at low excitation energy. First-forbidden decay rates are fastest when the initial and final states are each others ‘‘pygmy’’ resonance. In the present instance the

TABLE III. KHP_e predictions for $E\lambda$ ($\lambda = J_f - J_i$) transitions between even-parity states of ^{212}Po . $B(E\lambda)$ is given by $(e_p M_p + e_n M_n)^2 / (2J_i + 1) e^2 \text{fm}^{2\lambda}$ where e_p and e_n are the effective proton and neutron charges in units of e . The values of $B(E\lambda)$ given in the next-to-last column are for $e_p = 2$, $e_n = 1$ and are in Weisskopf units (W.u.). For $A = 212$, the $B(E2)$ and $B(E4)$ values corresponding to 1 W.u. are $75.09 e^2 \text{fm}^4$ and $1.00 \times 10^5 e^2 \text{fm}^8$, respectively. The transitions were calculated with WS radial wave functions. The experimental (expt) $B(E2)$ values are from Ref. [11] and the $B(E4)$ value is from Ref. [24].

J_i	J_f	M_p ($e \text{fm}^\lambda$)	M_n ($e \text{fm}^\lambda$)	$B(E\lambda)$ (KHP _e)	$B(E\lambda)$ (expt)
2	0	15.29	27.39	8.95	
4	2	21.27	42.70	10.75	
6	4	19.23	45.30	7.19	13.5 ± 3.7
8	6	14.33	36.95	3.37	3.95 ± 0.05
10	8	7.02	17.17	0.62	2.2 ± 0.6
12	10	19.19	11.93	1.35	
14	12	36.94	61.30	8.39	
16	14	29.10	48.83	4.62	
18	16	12.11	26.15	0.91	
18	14	421.80	1168.00	1.09	$2.3^{+4.6}_{-1.4}$

8_1^+ and 8_2^+ states of ^{212}Po share the role of the ‘‘pygmy’’ resonance of the ^{212}Bi 8_1^- state. Hence the fast decays.

The ^{212}Bi ground state decays 64% by β^- emission and 55% of the β^- decay is to the ^{212}Po ground state [1]. The predicted decay for this branch displays pathological cancellations between contributing terms. For this branch Eq. (4) gives

$$\begin{aligned} B_1^{(1)} &= [9.15M_1^u - 26.88M_1^x]^2 \\ &= [4.44 - 4.45]^2 \text{ fm}^2 \text{ (}\zeta \text{ approximation)}, \end{aligned} \quad (6)$$

i.e., almost complete cancellation. Recall that for first-forbidden transitions the β^- shape factor is [16]

$$C(W_e) = k(1 + aW_e + b/W_e + cW_e^2), \quad (7)$$

where Eq. (4) is a good approximation for k . In the present case what little of k survives the cancellation shown in Eq. (6) is itself almost exactly cancelled by the ka term so that the rate is given by the kc term (for $R1$ decays $kb = 0$). Hence the very large $\log f_0 t$ prediction for this transition. As might be imagined, reasonable changes in the wave functions and in the $q_\alpha(j_i j_f)$ and contributions from higher-order terms in the Behrens-Bühring expansion can bring the predicted rate into agreement with experiment.

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