

First observation of excited states in ^{184}Pb : spectroscopy beyond the neutron mid-shell

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Abstract. Excited states have been identified for the first time in ^{184}Pb , the first even-even Pb isotope beyond the $82 < N < 126$ mid-shell, using the recoil-decay tagging (RDT) technique. A collective band built on the first-excited 2^+ state has been observed. This resembles those seen in $^{186,188}\text{Pb}$ and the Hg isotones, and can thus be associated with a prolate-deformed shape. Variable moment of inertia (VMI) fits of the prolate 0^+ level energies in $^{184,186,188}\text{Pb}$ indicate that the minimum appears at $N = 103$, the same neutron number at which the corresponding minimum in Hg isotopes is observed.

PACS. 21.60.Ev Collective models – 23.20.-g Electromagnetic transitions – 25.70.gh Compound nucleus – 23.60.+e α decay

1 Introduction

There now exists a large body of experimental information supporting the coexistence of different shapes at low excitation energies in isotopes of lead. Low-lying 0_2^+ levels associated with weakly-deformed oblate structures have been observed to coexist with spherical ground states in several Pb isotopes with $N \geq 106$ [1,2]. These states are thought to result from 2 particle-2 hole proton excitations across the $Z = 82$ shell gap to $1h_{9/2}$ orbitals [3]. A rotational structure built on such a state has been seen in ^{196}Pb [4] and $M1$ shears bands associated with weakly-deformed oblate shapes are known in several Pb isotopes near $A = 198$ [5–7]. In $^{186,188}\text{Pb}$, low-lying deformed rotational structures have been observed at $I > 2\hbar$ by Heese *et al.* [8] and Baxter *et al.* [9]. The bands, which are similar to those in the corresponding isotones $^{184,186}\text{Hg}$, are associated with predicted [10] prolate-deformed minima. In Hg isotopes the excitation energy of the prolate-deformed intruder band decreases with decreasing neutron number, reaching a minimum near the mid-shell at $N = 103$ [11]. This trend has been reproduced using Nilsson-Strutinsky

calculations [10], which also predict a minimum close to $N = 103$ for Pb isotopes. In the present work γ rays have been studied for the first time in $^{184}\text{Pb}_{102}$. This represents an extension of our knowledge of excited states in even-even Pb isotopes to beyond the $82 < N < 126$ mid-shell.

2 Experimental Details

The $^{148}\text{Sm}(^{40}\text{Ca},4n)^{184}\text{Pb}$ reaction was employed with a target of thickness $500\mu\text{g}/\text{cm}^2$ and 99.94% enrichment. The beam, consisting of 95% ^{40}Ca and 5% ^{40}Ar , was delivered by the $K = 130$ MeV cyclotron at the Accelerator Laboratory of the University of Jyväskylä (JYFL) at a bombarding energy of 195 MeV. Prompt γ rays from the target were detected by the JUROSPHERE array which consisted of 11 TESSA-type [12] and 14 EURO-GAM Phase I [13] Compton-suppressed germanium detectors (absolute efficiency $\sim 1.6\%$ at 1.3 MeV). Recoiling fusion-evaporation residues were magnetically separated in-flight from the background of primary beam and nuclei produced by fission using the RITU gas-filled recoil separator [14]. The recoils were then implanted into a 80mm (horizontal) \times 35mm (vertical) Si detector covering approximately 70% of the image at the focal plane. The detector consisted of 16 strips of 5mm width, each position-sensitive in the vertical direction with a position

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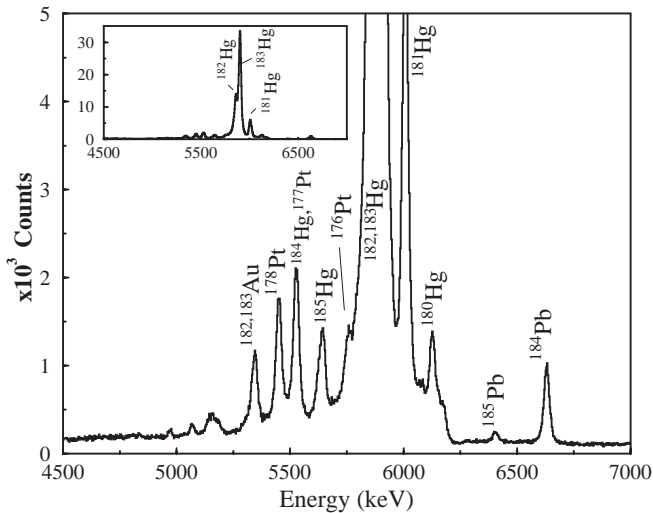


Fig. 1. Singles α -particle energy spectrum for the $^{40}\text{Ca} + ^{148}\text{Sm}$ reaction. The prominent peaks are assigned to known activities. An α -particle energy of 6626(6) keV was measured for ^{184}Pb using 9000(100) decays

resolution of $400\mu\text{m}$. For the recoils and their α decays, the position, energy and time of detection was recorded. Figure 1 shows the energy spectrum for single α -decay events. The spectrum was calibrated internally using the following α -particle energies from literature [15]: ^{178}Pt $E_\alpha = 5446(3)$ keV, ^{177}Pt $E_\alpha = 5517(4)$ keV, ^{183}Hg $E_\alpha = 5904(4)$ keV, ^{181}Hg $E_\alpha = 6005(4)$ keV and ^{180}Hg $E_\alpha = 6119(5)$ keV. An α -particle energy of 6626(6) keV was obtained for ^{184}Pb from 9000(100) decays. This is consistent with the previously-measured value of 6632(10) keV [16].

3 Results

Figure 2(a) is a spectrum showing prompt γ rays in coincidence with all recoils detected at the RITU focal plane. It is dominated by those from Hg isotopes produced in $(^{40}\text{Ca}, \alpha\text{xn})$ and $(^{40}\text{Ar}, \text{xn})$ channels. Prompt γ rays emitted by ^{184}Pb residues were selected from those produced via other exit channels using the recoil-decay tagging (RDT) technique [17,18]. The events of interest corresponded to the detection of a recoil and a ^{184}Pb α decay at the same position in the Si detector within a time interval of 1.5 seconds (approximately three ^{184}Pb α -decay half-lives). Using these events and the method prescribed in [19], a half-life value of 0.60(9)s for ^{184}Pb was deduced which is consistent with the value of 0.55(6)s obtained in the previous measurement [16]. Figure 2(b) is a spectrum showing RDT-gated γ rays in ^{184}Pb . Four γ -ray transitions are clearly identified along with Pb X-rays. The dashed lines marked on the spectra illustrate that the transitions observed in Fig. 2(b) are not observed in Fig. 2(a) because of the large background of γ rays from other nuclei. The inset to Fig. 2(b) is a tentative level scheme deduced using this spectrum. The $I^\pi = (2^+)$ state is assigned an energy of 701.5 keV on the basis of 2^+ energy-

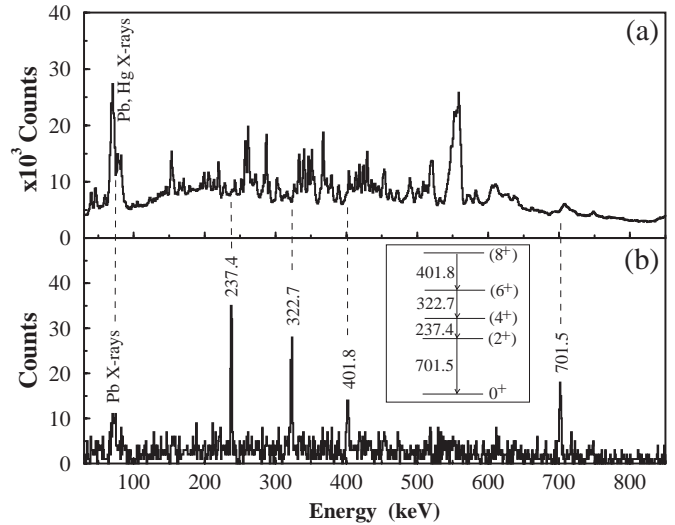


Fig. 2. **a** Singles energy spectrum of γ rays in coincidence with all fusion-evaporation residues detected in the RITU focal-plane Si detector, **b** singles energy spectrum showing γ -ray transitions in ^{184}Pb and Pb X-rays. The spectrum was generated using the RDT technique (see text). Inset: tentative decay scheme for ^{184}Pb

level systematics in Pb isotopes; the only intense peak in the region of the spectrum where the 2^+ to 0^+ transition is expected has an energy of 701.5 keV. The transitions of energies 237.4, 322.7 and 401.8 keV appear to form a rotational band similar to those built on 2^+ states in $^{186,188}\text{Pb}$. They are therefore tentatively assigned as the following transitions: (4^+) to (2^+) , (6^+) to (4^+) and (8^+) to (6^+) respectively. The relative intensities of the 701.5, 237.4, 322.7 and 401.8 keV transitions are 100(11), 98(12), 94(13) and 38(13) respectively.

4 Interpretation

Aligned angular momentum, i_x , as a function of rotational frequency is shown for $^{184,186,188}\text{Pb}$ in Fig. 3. The transition energies of $^{186,188}\text{Pb}$ were taken from [8,9]. The same variable moment of inertia reference parameters, $J_0 = 27 \hbar^2\text{MeV}^{-1}$ and $J_1 = 199 \hbar^4\text{MeV}^{-3}$, have been used for the three nuclei. It should be noted that there is evidence for quasi-particle alignment in ^{188}Pb at $\hbar\omega = 0.3$ MeV. Above $I = 2\hbar$ the alignment properties of the rotational bands in the three isotopes are similar, but not identical. The alignment of the intruder bands decreases with increasing neutron number, indicating a gradual decrease in collectivity. This is in contrast with the behaviour of the Hg isotones where no significant change in collectivity with neutron number is observed for $^{182,184,186}\text{Hg}$ [21]. The behaviour of the three Hg isotopes above 0.2 MeV is almost identical to that of ^{186}Pb in Fig. 3. The similarities between the Pb and Hg isotopes indicate that the low-lying bands in the Pb isotopes are built on the same prolate-deformed structures assigned to the Hg bands.

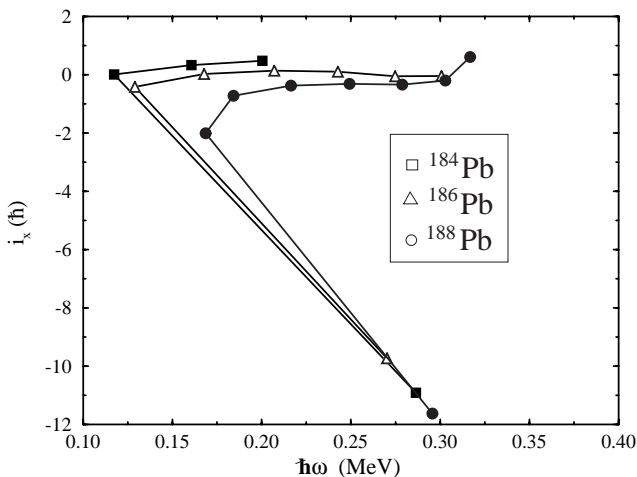


Fig. 3. Plots of aligned angular momentum, i_x , as a function of rotational frequency, $\hbar\omega$, for the ^{184}Pb (this work) and $^{186,188}\text{Pb}$ [8,9]. Rotational references with Harris parameters of $J_0 = 27 \hbar^2\text{MeV}^{-1}$ and $J_1 = 199 \hbar^4\text{MeV}^{-3}$ have been subtracted

Figure 4 shows plots of excitation energy as a function of $I(I+1)$ for the levels in ^{184}Pb from this work and $^{186,188}\text{Pb}$ from refs. [8,9]. The values of E^* at $0\hbar$ for these three nuclei were estimated using variable moment of inertia (VMI) [22] fits using levels in the intruder bands. Dashed lines connect these points to the experimental data points. The energies of the following levels were used in the fits: levels with $I^\pi = 4^+$ to 8^+ for ^{184}Pb , 8^+ to 14^+ for ^{186}Pb and 6^+ to 14^+ for ^{188}Pb . The 16^+ state in ^{188}Pb was not included in the fit because here there is evidence for quasi-particle alignment (see Fig. 3). Omission of either the 6^+ or 14^+ states, which may be displaced by mixing and quasi-particle alignment respectively, did

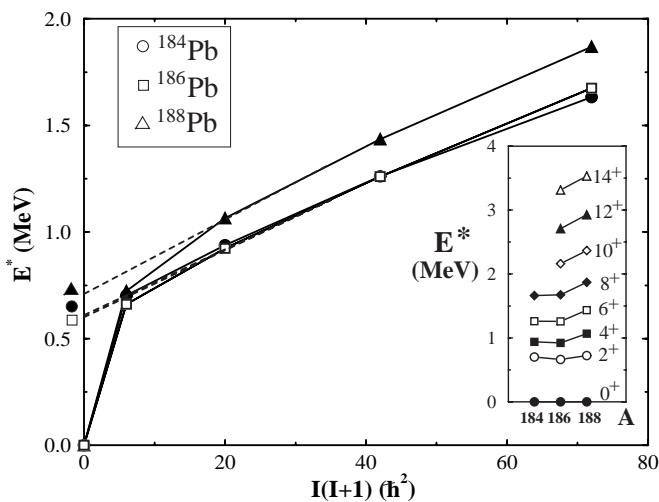


Fig. 4. Plots of excitation energy, E^* , as a function of spin, I , for $^{184,186,188}\text{Pb}$. The dashed lines are extrapolations to zero spin, where the values of E^* have been determined using VMI fits. Inset: yrast states in $^{184,186,188}\text{Pb}$

Table 1. Fitted VMI parameters and bandhead energies

| Nucleus | J_0 (keV^{-1}) | C (10^6keV^3) | Bandhead energy (keV) |
|-------------------|-----------------------------|-----------------------------|-----------------------|
| ^{184}Pb | 0.0281 | 2.19 | 610 |
| ^{186}Pb | 0.0284 | 2.80 | 600 |
| ^{188}Pb | 0.0244 | 2.24 | 710 |

not alter the estimated bandhead energy for ^{188}Pb . The estimated bandhead energies are given in Table 1, along with the fitted VMI parameters J_0 and C .

In order to compare the prolate bandhead energies for the three Pb isotopes it is necessary to ascertain the extent to which the low-spin (0^+ , 2^+) states are perturbed by mixing with other configurations. The mixing of states of different deformation in ^{188}Pb is discussed in [20]. The mixing of the 0^+ states in this nucleus serves to *increase* the prolate 0^+ level energy by ~ 50 keV. This perturbation does not effect the following arguments so the unmixed prolate bandhead energy obtained from the VMI fit is used. The high-lying, high-spin members of the prolate band in ^{188}Pb are unmixed and unperturbed so the estimated bandhead energy of 710 keV is reliable. The oblate-deformed 0^+ states have not been observed in $^{184,186}\text{Pb}$. The excitation energy of the oblate-deformed minimum is predicted [10] to increase suddenly as N decreases from 106 to 104, so the mixing between the oblate and prolate configurations is assumed to be weak for $^{184,186}\text{Pb}$.

In the present work the effect of mixing between the prolate-deformed and spherical configurations has been calculated. In ^{186}Pb a perturbation of the prolate bandhead of 2.7 keV is expected if the experimentally-observed 2^+ state is assumed to be prolate, while a 3.7 keV displacement is predicted if the 2^+ state is spherical. Both scenarios lead to the conclusion that the energy of the prolate-deformed 0^+ state is close to 600 keV.

Only 4 excited levels were observed in ^{184}Pb , therefore the 4^+ level was included in the VMI fit to estimate the energy of the prolate 0^+ state. In practice the 4^+ state will be mixed and therefore displaced, resulting in a perturbation of the VMI-fitted 0^+ state. Displacement of the 4^+ level in ^{186}Pb was calculated to be small (7 keV) so the displacement of the 4^+ state, and therefore the fitted 0^+ state, in ^{184}Pb is also expected to be small. Measurement of ^{185}Bi proton decay [23] has provided additional evidence that the mixing strength is small in ^{184}Pb . If the 4^+ state in ^{184}Pb is assumed to be prolate as in ^{186}Pb , the unmixed prolate 4^+ level will be higher in energy than the observed (mixed) 4^+ state and therefore the bandhead energy will be raised above the fitted value. The prolate-deformed 0^+ state may therefore be slightly higher than the VMI-fitted value of 610 keV. However, within the uncertainties of the fitting, the energies of the prolate 0^+ states in ^{184}Pb and ^{186}Pb are the same, regardless of which configurations are assumed for the 2^+ and 4^+ states. This indicates that the excitation energy of the prolate-deformed configuration has a minimum close to $N = 103$. The corresponding

minimum in Hg isotopes can be found at the same neutron number [11].

It would be of considerable interest to study the α -decay fine structure of ^{190}Po in order to measure directly the excitation energy of the prolate 0^+ state in ^{186}Pb and extract the mixing strength from the hindrance factor relative to decay to the ground state. Unfortunately a corresponding measurement for ^{188}Po appears to be beyond the reach of current experimental techniques.

5 Summary

Transitions have been identified for the first time in ^{184}Pb using the RDT technique. This represents γ -ray spectroscopy beyond the $82 < N < 126$ mid-shell in Pb isotopes. A low-lying rotational structure has been observed. The band behaves in a similar way to those observed in $^{186,188}\text{Pb}$ and the Hg isotones and is thus assigned a prolate-deformed configuration. VMI calculations of the bandhead energies for the rotational structures in $^{184,186,188}\text{Pb}$ reveal that the prolate-deformed minimum occurs at $N = 103$, as seen in the Hg isotopes.

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