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## New aspects of the low-energy structure of 211At

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Lifetimes of low-energy states in <sup>211</sup>At have been measured using the recoil-distance Doppler shift, Dopplershift attenuation, and fast-timing methods at the University of Cologne. The obtained reduced transition probabilities have been compared to two shell-model calculations, a large-scale shell-model calculation using the Kuo-Herling residual interaction and a calculation using a single-j approximation for protons in the  $0h_{9/2}$ orbital. The newly obtained reduced transition probabilities are described very well by a single-j calculation. This, together with the fact that the energy spectrum of <sup>211</sup>At is also well described, indicates that seniority can be regarded as a good quantum number in <sup>211</sup>At. While the single-j calculation can only describe states with a dominant  $0h_{9/2}^3$  configuration, the presence of other low-lying proton orbitals, like  $1f_{7/2}$  and  $0i_{13/2}$ , requires a multi-j calculation. The multi-j calculation using the Kuo-Herling interaction gives a satisfactory description of the nuclei in the region but significantly overestimates some of the ground-state transition probabilities, for example, the  $B(E2; 13/2_1^- \rightarrow 9/2_1^-)$  value in <sup>211</sup>At. This discrepancy has been attributed to the presence of higher-order particle-hole excitations in the wave function of the ground state, which are not accounted for by the Kuo-Herling interaction. The effects of those excitations on the transition rates, however, are weaker in  $^{211}$ At than they are in  $^{210}$ Po. On the other hand, a strong underestimation of the E2 strengths involving the  $7/2_1^-$  state is also observed, where one proton occupies the  $0f_{7/2}$  orbital. Therefore, a phenomenological modification to the  $\langle 0h_{9/2}, 0h_{9/2}|\hat{V}|0h_{9/2}, 1f_{7/2} \rangle_{J=2}$  two-body matrix element has been introduced which leads to a considerably better description of the structure of <sup>210</sup>Po and <sup>211</sup>At. However, the origin of this effect needs to be further investigated.

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