## Title

Shell Model Calculation of Po212



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## DISCLAIMER

Shell Model Calculation of $Po^{212}$.* Norman K. Glendenning and Kichinosuke Harada†, Lawrence Radiation Laboratory, Berkeley.--So far there are a few shell model calculations[1-3] on $Po^{212}$, but unfortunately their results are more or less different from each other, owing to the use of different residual forces. A more realistic shell model calculation based on Glendenning's method is done for the lower lying levels and the high-spin isomer of $Po^{212}$. In this calculation we use the residual p-p, n-n and p-n interactions which are so determined as to reproduce the experimental $Po^{210}$, $Pb^{210}$ and $Bi^{210}$ spectra, and adopt different size parameters for proton- and neutron-oscillator potentials. The results will be discussed and compared with those of the previous works.

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