

GROUND-STATE SPLITTINGS OF ACTINIDE IONS IN CaF₂

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The energy-level splittings of Am²⁺, Cm³⁺, Es²⁺, Pu³⁺, and U⁴⁺ in an octahedral site of CaF₂ were studied using restricted Hartree-Fock and spin-orbit configuration-interaction calculations. The CaF₂ host was modeled with a large finite cluster of ions which approximate the Madelung potential of the crystal lattice^a. The actinide dopants were treated with relativistic effective core potentials and Gaussian correlation consistent double-zeta plus polarization basis sets. The primary properties of interest are energy splittings and magnetic moments.

^aN. W. Winter, M. Ross and R. M. Pitzer *J. Phys. Chem.* **94**, 1172 (1990).