

Theoretical Studies on the Gyromagnetic Factors for Nd^{3+} in Scheelites-Type ABO_4 Compounds

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The gyromagnetic factors for Nd^{3+} in scheelite-type ABO_4 compounds ($A = \text{Cd}, \text{Ca}, \text{Pb}, \text{Ba}$; $B = \text{Mo}, \text{W}$) are theoretically studied by the perturbation formulas of the anisotropic g factors g_{\parallel} and g_{\perp} for a $4f^3$ ion in tetragonal symmetry. In these formulas, the contributions to the g factors due to the second-order perturbation terms and the admixtures of various energy levels are taken into account. The relevant crystal-field parameters are determined by the superposition model and the local geometrical relationship of the A^{2+} sites occupied by the impurity Nd^{3+} . The obtained g factors agree reasonably with the observed values. The discrepancies between theory and experiment are discussed.

Key words: EPR; Crystal-fields and Spin Hamiltonian; Nd^{3+} ; Scheelite-type ABO_4 Compounds ($A = \text{Cd}, \text{Ca}, \text{Pb}, \text{Ba}$; $B = \text{Mo}, \text{W}$).