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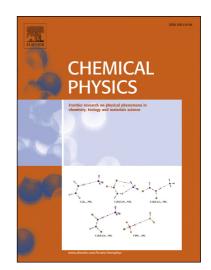
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Spectroscopic properties of ErAl₃(BO₃)₄ single crystal

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Abstract

Single crystal of $ErAl_3(BO_3)_4$ was grown and its structure was studied. Polarized absorption spectra of $ErAl_3(BO_3)_4$ single crystal were measured in the spectral range 1670-330 nm (6000-30000 cm⁻¹). The Judd-Ofelt spectroscopic parameters have been determined: Ω_2 =4.87·10⁻²⁰ cm², Ω_4 =2.49·10⁻²⁰ cm², Ω_6 =2.72·10⁻²⁰ cm². These parameters have been used to calculate the radiative transition probabilities, the multiplet luminescence branching ratios and the fluorescence life times of the manifolds. The luminescence spectra due to transitions ${}^2H_{11/2} \rightarrow {}^4I_{15/2}$ (526 nm), ${}^4S_{3/2} \rightarrow {}^4I_{15/2}$ (548 nm), ${}^4F_{9/2} \rightarrow {}^4I_{15/2}$ (662 nm), ${}^2H_{11/2} \rightarrow {}^4I_{13/2} + {}^4I_{9/2} \rightarrow {}^4I_{15/2}$ (800 nm) and ${}^4S_{3/2} \rightarrow {}^4I_{13/2}$ (855 nm) were recorded in α , σ and π polarizations. The most intensive luminescence belonged to ${}^4S_{3/2} \rightarrow {}^4I_{15/2}$ transition (548 nm).

PACS: 78.20.-e, 78.40.Ha

Keywords: rare earth compounds; Er^{3+} ; spectroscopic properties; Judd-Ofelt analysis; luminescence.

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