## Crystal-Field Energy Levels of Trivalent Erbium Ion in Cubic Symmetry

Said Laachir<sup>a</sup>, Mohamed Moussetad<sup>b</sup>, Rahma Adhiri<sup>b</sup>, and Ahmed Fahli<sup>a</sup>

<sup>a</sup> UFR S.I.R.I, Université Hassan II-Mohammedia Faculté des Sciences, Ben M'sik, B.P 7955 Casablanca, Morocco

<sup>b</sup> L.P.S.C.M, Université Hassan II-Mohammedia Faculté des Sciences, Ben M'sik, B.P 7955 Casablanca, Morocco

Reprint requests to S. L.; E-mail: saidlaachir@yahoo.fr

Z. Naturforsch. **66a**, 457–460 (2011); received December 16, 2010

This paper describes a scheme for the numerical calculation of crystal field (CF) energy levels and at the same time wave functions of the trivalent erbium ion in cubic symmetry. The 16-fold degenerate term  ${}^{4}I_{15/2}$  of the trivalent erbium ion splits into three Stark quartets  $\Gamma_{8}$  and two different doublets  $\Gamma_{6}$  and  $\Gamma_{7}$  (irreducible representations). The CF energy matrix of the  $\mathrm{Er}^{3+}$  ion has been constructed and calculated from the complete diagonalization method, and the corresponding wave functions were used to calculate the ground state g-values. This method is outlined and illustrated by the examples of the Si:Er for comparison. The calculated g-factors are g = 6.8 and g = 6.0 for  $\Gamma_{6}$  and  $\Gamma_{7}$ , respectively.

Key words: Crystal Field; Energy Level; Erbium; Rare Earth.