

PHASE EQUILIBRIA AND CRYSTAL CHEMISTRY OF THE CsF–LuF₃ SYSTEM

D. Avignant*, J. Metin, D. Chatonier and J. C. Cousseins

Université de Clermont-Ferrand II, Laboratoire de Chimie des Solides, B.P. 45,
63170 Aubière (France)

Carrying out an exploratory program intended for the research of efficient matrices for use as crystalline host materials for U.V. emission in solid state lasers realization, a recent investigation of the CsF–LuF₃ system by differential thermal analysis and X-ray diffraction has been undertaken. This study has allowed us to characterize three compounds with formula Cs₃LuF₆, CsLu₂F₇ and CsLu₃F₁₀ as well as various metastable homogeneous domains lying in the high temperatures region for concentrations between 75 and 86 moles per cent of LuF₃ approximatively.

The crystal structure of a metastable non stoichiometric phase Cs_{1-x}Lu₃F_{10-x} which represents a new structural type characterized by a three-dimensional framework of corner and edge-shared pentagonal bipyramids surrounding the rare earth has been determined from single crystal X-ray diffraction data. Relationships with other structural types, namely RbIn₃F₁₀ [1], β-RbLu₃F₁₀ [2], CsYb₃F₁₀ [3] and U₃O₈ will be discussed.

As more becomes known about crystal chemical of rare earths with small ionic radii we should be able to predict new structural types arising from intergrowth, the existence of which will be discussed at the light of preliminary results.

- 1 J.C. CHAMPARNAUD-MESJARD, D. MERCURIO and B. FRIT, *J. Inorg. Nucl. Chem.*, 39, 947 (1977).
- 2 A. ARBUS, M.T. FOURNIER, J.C. COUSSEINS and A. VEDRINE, *Acta Cryst.*, B38, 75 (1982).
- 3 A. ALEONARD, M.Th. ROUX and B. LAMBERT, *J. Solid State Chem.*, 42, 80 (1982).