

Phase Diagram of the System LiF-GdF₃-YF₃

I.M. Ranieri¹, A.H.A. Bressiani², S.P. Morato¹, S.L. Baldochi¹,
K. Shimamura³, T. Fukuda³

1- Center for Lasers and Applications, IPEN-CNEN/SP, CP 11049, Butantã,
05422-970, São Paulo, SP, Brazil.

2- Department of Materials Engineering, IPEN-CNEN/SP, CP 11049, Butantã,
05422-970, São Paulo, SP, Brazil.

3- Institute for Materials Research, Tohoku University,
Sendai, 980-8577, Japan.

Corresponding author email: iranieri@net.ipen.br

LiGdF₄ crystals have an important application as laser active media when doped with rare earth ions [1]. The LiF-GdF₃ phase diagram presents two invariant points: a eutectic at 25 mol% GdF₃ and 698 °C, and a peritectic at 34 mol% GdF₃ and 750 °C. GLF is the unique intermediary compound. In a previous paper [2] it was investigated the codoping of GLF crystals with yttrium, and good quality crystals could be obtained using the peritectic composition of the system LiF-GdF₃ as starting composition.

In this work the phase diagrams of the system LiF- Gd_(1-x)Y_xF₃ (x= 0,5 e 0,75) have been constructed using differential thermal analysis. The measurements were performed with samples weighing around 50 mg placed in open platinum crucibles, under a flux of purified helium, with a heating rate of 10 °C/min. The phase diagrams were determined up to compositions of 40 mol% LiF: 60 mol% Gd_(1-x)Y_xF₃.

To determine the phases present in the phase diagrams, some samples were melted under a flux of hydrogen fluoride gas and cooled at rates of 10-20 °C. One of the samples was observed using a scanning electron microscope with temperature and from 100 to 600°C, in steps of 100°C. The patterns were analyzed by the program DBWS-9807a [3] which makes the refinement of the structural parameters using the Rietveld method. The variations of the parameters with the temperature were fitted to third-order polynomials. The parameters behave in a way which is compatible to the thermal properties of the compounds. The measured lattice parameters of Cr:LiCAF are larger than those of pure and Ce:Na-doped crystals, which showed very similar values. The same was not observed for LiSAF host. The measured lattice parameters of Ce:Na:LiSAF are larger than the observed values of pure and Cr-doped LiSAF.

References:

- [1] Payne, S.A.; Chase, L.L.; Smith, L.K.; Kway, W.L.; Newkirk, H.W. "Laser performance of LiSrAlF₆:Cr³⁺", J. Appl. Phys. 66, 1051 (1989).
- [2] Liu, Z.; Izumida, S.; Ohtake, H.; Sarukura, N.; Shimamura, K.; Mujilatu, N.; Baldochi, S. L.; Fukuda, T. "High pulse-energy, all-solid-state, ultraviolet laser oscillator using large Czochralski-grown Ce:LiCAF crystal" Jpn. J. Appl. Phys., 37, L1318 (1998).
- [3] Young, R.A.; Larson, A.C.; Paiva-Santos, C.O. User's guide to program DBWS-9807a for Rietveld analysis of x-ray and neutron powder diffraction patterns. School of Physics, Georgia Institute of Technology, Atlanta, (1999).