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Tünde Alapi Róbert Berkecz István Ilisz

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COMBINED EXPERIMENTAL AND DFT STUDY OF LITHIUM-INDIUM-OXIDE STRUCTURE AND VIBRATIONAL PROPERTIES

Robert Vigi¹, Ljubica Đačanin Far², Svetlana Lukić-Petrović¹, and Tamara Ivetić¹

¹University of Novi Sad, Faculty of Sciences, Department of Physics, Trg Dositeja
Obradovića 3, 21000 Novi Sad, Serbia

²University of Belgrade, Vinča Institute of Nuclear Sciences, Center of Excellence for
Photoconversion, P.O. Box 522, 11000 Belgrade, Serbia
e-mail: robert1995@uns.ac.rs

Abstract

A promising lithium-indium-oxide (LiInO₂) wide band-gap semiconductor for scintillating detection, photoluminescence, and photocatalysis [1-3] was prepared by a mechanochemical solid-state synthetic procedure that can be found elsewhere [3]. Its structure and morphology were investigated by using X-ray diffraction (XRD), scanning electron microscopy (SEM), and Raman spectroscopy. SEM images show agglomerates of relatively uniform size of around 300 nm spherical-shaped particles of LiInO₂ powder, while the XRD pattern confirmed the formation of the nanocrystalline tetragonal structure with $I4_1/amd$ space group (no. 141) symmetry. Detailed vibration analysis, together with the assignments of the band modes, was performed through the best-fit match of the experimental and density functional theory (DFT) calculated Raman spectrum. Geometry optimizations and vibrational frequencies calculations were conducted using B97-1 functional correlation [4] and LanL2DZ was used as a basis set.

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