## 11. Photoluminescence Properties of Eu3+ Doped Li3AlMo3O12 Red Phosphors

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## Abstract

Eu<sup>3+</sup> doped red phosphors Li<sub>3</sub>AlMo<sub>3</sub>O<sub>12</sub> were prepared by the combustion method, and their photo luminescent properties were studied. The photoluminescence spectra results indicate that the excitation spectra of the samples consist of a broad absorption band of 200-350 nm and some sharp lines around 350-420 nm which corresponds to the f-f transitions within 4f<sup>8</sup> configuration of Eu<sup>3+</sup>. The emission spectra consist of the orange region  $^5D_0 \rightarrow ^7F_1(593 \text{ nm})$  and the red region  $^5D_0 \rightarrow ^7F_2(617 \text{ nm})$ . Among these phosphors, Li<sub>3</sub>AlMo<sub>3</sub>O<sub>12</sub> synthesized at 750°C exhibits the strongest red emission under 393 nm excitation and appropriate CIE chromaticity coordinates (x=0.637, y=0.359) close to the NTSC standard value. Al<sup>3+</sup> was selected as a potential candidate to replace Eu<sup>3+</sup> in the lithium molybdate host lattices. It was discussed that there is very limited information regarding the impact of Al<sup>3+</sup> cations on the luminescent properties of molybdate based phosphors. The fact that Al<sup>3+</sup> has a very small atomic radius in comparison to those of Eu<sup>3+</sup> and that it has a maximum coordination number of 6 makes it an interesting and potentially good candidate for improving the molybdate phosphors. In addition, Al<sup>3+</sup> originating in Al(NO<sub>3</sub>)<sub>3</sub>.9H<sub>2</sub>O (the starting material used in this experiment) is much cheaper than any Eu<sup>3+</sup> starting material.

## 1. Introduction

Molybdate crystals are presently the subject of considerable interest and detailed evaluation because of their valuable structural, thermal, electrical, and spectroscopic properties

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