

Luminescence and Thermal-Quenching Properties of Red-Emitting $\text{Ca}_2\text{Al}_2\text{SiO}_7:\text{Sm}^{3+}$ Phosphors

March 2020, [Journal of Electronic Materials](#) 49(6)

DOI: [10.1007/s11664-020-08086-x](https://doi.org/10.1007/s11664-020-08086-x)

Abstract

$\text{Ca}_2\text{Al}_2\text{SiO}_7:\text{x.Sm}^{3+}$ ($x = 0.5, 1.0, 1.5, 2.0, 2.5,$ and 3.5 mol.%) ($\text{CAS}:\text{x.Sm}^{3+}$) phosphors were synthesized by a solid-state reaction technique. The structure, photoluminescence properties and thermal stability of phosphors were investigated in detail. Results of X-ray diffraction show that $\text{CAS}:\text{x.Sm}^{3+}$ materials have a single-phased tetragonal structure, and an expansion of the unit cell relates to the increasing of Sm^{3+} concentration. Photoluminescence study displayed that the $\text{CAS}:\text{x.Sm}^{3+}$ phosphors reach the highest emission intensity at 1.5 mol.% Sm^{3+} and achieved the luminescence quenching phenomenon a higher concentration. The dominant interaction mechanism of the concentration quenching process is determined due to the dipole–dipole interaction, and the critical transfer distance (R_c) is 26.7 Å. The temperature dependence of photoluminescence spectra indicates that the $\text{Ca}_2\text{Al}_2\text{SiO}_7:\text{Sm}^{3+}$ (1.5 mol.%) phosphor possess good thermal stability, and that the activation energy is around 0.12 eV (968 cm^{-1}). Several characteristic vibrations in the 200–1000- cm^{-1} region were observed by Raman spectra, and the color chromaticity coordinates of the samples were also calculated and discussed.