"Nanocomposites and nanomaterials"

Structure studies of Ag-doped As₂S₃ chalcogenide glasses

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Nanostructured chalcogenide glasses (ChGs) have numerous optical and photonics applications, due to their optical transparency in IR spectral region, chemical stability, photosensitivity phenomena, high linear and non-linear refractive index. ChGs have great potential for many applications including integrated and non-linear optics, information technology, etc.

In present work the results of investigations on the influence of Ag doping (10, 15, 20 wt.%) on structural properties of As₂S₃ glasses are presented. X-ray diffraction studies confirmed amorphous structure of glasses. Radial distribution functions have been obtained and analyzed, which testified about certain increase of the first coordination sphere radius after introduction of silver into As₂S₃. From obtained HH and HV-polarized Raman spectra of doped As₂S₃:Ag it can be seen that structure changes. The Raman spectra of these ChGs are dominated by strong band at 340 cm⁻¹ attributed to the symmetric As-(S)-As stretching vibrations in AsS_{3/2} pyramids. Besides this strong band at 340 cm⁻¹, there are additional features (shoulders) at 314 cm⁻¹ and can be assigned to the asymmetric stretching modes of AsS_{3/2} pyramids. Weak bands situated at 187, 221, 231 cm⁻¹ correspond to the presence of the non-stoichiometric molecular fragment of As₄S₄ nanophase. Doping of As₂S₃ glass with Ag causes a slight increase of the band intensity located at 187, 221, 231 cm⁻¹. The presence of excess sulfur is indicated by the weak band at 475 cm⁻¹, characteristic for the vibrations of S-S bonds, and it is increased with the Ag introduction. Deconvolution of obtained Raman spectra was performed for analysis of their compositional dependencies.

The main effect that is observed at introduction of silver into nanostructured As₂S₃ glass is the change of the relative concentration of the main and non-stoichiometric molecular structural units characteristic for As₂S₃ glasses.

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