

Structural and electrical properties of bismuth magnesium tantalate pyrochlores.

Abstract

The subsolidus cubic pyrochlore phases in the $\text{Bi}_2\text{O}_3\text{-MgO-Ta}_2\text{O}_5$ (BMT) system were prepared with the proposed formula, $\text{Bi}_{3+(5/2)x}\text{Mg}_{2-x}\text{Ta}_3-(3/2)x\text{O}_{14-x}$ ($0.12 \leq x \leq 0.22$). Replacement of smaller cations, Mg^{2+} and Ta^{5+} by larger Bi^{3+} cations with considerable oxygen non-stoichiometry within structure was proposed. The synthesised samples were confirmed phase pure by X-ray powder diffraction and their refined lattice parameters were in the range of $10.5532(4)\text{-}10.5672(9)$ Å. The grain sizes of the samples determined by SEM analysis were in the range of $0.6\text{-}10.60$ μm and their average relative densities were more than 80%. Five infrared-active modes were also observed in their FTIR spectra due to their metal-oxygen bonds. The BMT pyrochlores were highly electrical resistive with high dielectric constants, ϵ' in the range of $\sim 70\text{-}85$; dielectric losses, $\tan \delta$ in the order of 10^{-3} at frequency 1 MHz and a negative temperature coefficient of permittivities, $\text{TC}\epsilon'$ of ~ -158 to -328 $\text{ppm}/^\circ\text{C}$.

Keyword: C. Dielectric properties; D. Tantalates; Pyrochlores; Ceramics.