# ELECTRONIC STRUCTURE AND MAGNETIC PROPERTIES $Gd_{1-x}Tb_xNi_3$ COMPOUNDS

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The magnetic susceptibility, the crystal and electronic structure of the polycrystalline intermetallic  $\mathrm{Gd}_{1-x}\mathrm{Tb}_x\mathrm{Ni}_3$  compounds has been investigated. All investigated compounds crystallize in the rombohedral PuNi $_3$ type of crystal structure (space group R-3m). The effect of partial substitution Gd by Tb is reflected in the linear decrease of the lattice parameters in the whole range of x, whereas the Curie temperature determined from AC susceptibility measurements decreases from 114 K for GdNi $_3$  to 100 K when x=0.5 and is unchanged until full concentration of Tb (x=1). Additionally the presence of the helimagnetic phase when Gd is substituted by Tb was observed. The electronic structure was studied by using X-ray Photoelectron Spectroscopy (XPS). Both valence band and core level spectra are analyzed. The valence band spectra close to the Fermi level are dominated by the Ni 3d and Gd/Tb states.

**←** 13.4 cm −

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 $9.7~\mathrm{cm}$