## The Solid Solution $CeAuIn_{1-x}Mg_x$ – Structure, Magnetic Properties and Specific Heat Data

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Z. Naturforsch. 61b, 495 – 502 (2006); received February 21, 2006

Four different samples of the solid solution  $\text{CeAuIn}_{1-x}\text{Mg}_x$  with x = 0.2, 0.4, 0.6, and 0.8 have been prepared from the elements in sealed tantalum tubes in an induction furnace. The samples were characterized through X-ray powder and single crystal data: ZrNiAl type,  $P\overline{6}2m$ , Z = 3, a =774.54(7), c = 420.32(10) pm, wR2 = 0.0203,  $395 F^2$  values, 15 variables for CeAuIn<sub>0.871</sub>Mg<sub>0.129</sub>, a = 775.25(7), c = 419.36(10) pm, wR2 = 0.0488, BASF = 0.10(1), 397  $F^2$  values, 16 variables for CeAuIn<sub>0.640</sub>Mg<sub>0.360</sub>, a = 774.62(7), c = 420.13(10) pm, wR2 = 0.0435, 376  $F^2$  values, 15 variables for CeAuIn<sub>0.445</sub>Mg<sub>0.555</sub>, a = 773.80(11), c = 420.82(8) pm, wR2 = 0.0415, 392  $F^2$  values. 15 variables for CeAuIn<sub>0.228</sub>Mg<sub>0.772</sub>. The lattice parameters show no pronounced changes within the solid solution. The largest shift occurs for the x parameter of the mixed occupied In/Mg positions. Due to the difference in size, the trigonal prisms around the Au1 atoms at the origin become smaller with an increasing content of magnesium. The In/Mg-In/Mg distances decrease from 334.5 (CeAuIn<sub>0.871</sub>Mg<sub>0.129</sub>) to 328.3 (CeAuIn<sub>0.228</sub>Mg<sub>0.772</sub>) pm, and consequently one observes also shorter bonds to the Au1 atoms with an increasing content of magnesium concentration. Susceptibility measurements reveal trivalent cerium for all CeAuIn<sub>1-x</sub>Mg<sub>x</sub> compounds, with no evidence sign of magnetic ordering down to 2 K. The disorder created by chemical substitution destroys the longrange magnetic ordering which can be attributed to the triggering of non Fermi-liquid (NFL) like behavior.

Key words: Cerium Compounds, Intermetallics, Crystal Chemistry, Magnetism