

Letter

# Crystal structure and magnetism of $\text{YbT}_x\text{Ga}_{4-x}$ , $T = \text{Zn, Cd}$ with the $\text{BaAl}_4$ -type

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## Abstract

Intermetallic compounds of the composition  $\text{YbT}_x\text{Ga}_{4-x}$  ( $T = \text{Zn, Cd}$ ) were synthesized and their crystal structure and magnetic properties were studied. These compounds crystallize in the tetragonal  $\text{BaAl}_4$  type structure with a random occupation of the 4e site by Ga, Zn(Cd). The magnetic ground state of Yb in these compounds is nonmagnetic.

*Keywords:* Crystal structure; Magnetism; Intermetallic compounds

## 1. Introduction

Recent results on the ternary  $\text{BaAl}_4$ -type gallides with ytterbium and transition metals such as Cu, Ag, Au, Pd, and Pt showed magnetically dipositive ytterbium for Cu, Ag, Au and Pd with a tendency towards tripositive behavior for  $\text{YbPt}_{0.5}\text{Ga}_{3.5}$  [1]. An extension of this research was designed to cover the metals with a filled 3d or 4d shell, zinc and cadmium, and became the subject of the present paper.

## 2. Experimental

Alloys with the nominal compositions  $\text{YbT}_x\text{Ga}_{4-x}$ ,  $x$  at 0.10, 0.25, 0.5, 0.75 and 1.0 for  $T = \text{Zn}$  and  $x = 0.5$  for  $T = \text{Cd}$ , each with a total weight of 2 g, were prepared by high frequency melting in tantalum crucibles under argon. Further details of sample preparation, heat treatment (400 °C for 500 h), X-ray powder diffraction techniques and methods of magnetic measurements in the range from 4.2 K to 300 K can be found from our earlier reports on a similar subject [1,2].

## 3. Results and discussion

### 3.1. Formation, homogeneity region and crystal structure for $\text{YbT}_x\text{Ga}_{4-x}$

Evaluation of the X-ray powder diagrams of the two alloy systems  $\text{YbZn}_x\text{Ga}_{4-x}$  and  $\text{YbCd}_x\text{Ga}_{4-x}$  confirmed the isotypism with the crystal structure of the  $\text{BaAl}_4$ -type. Whereas for  $\text{YbZn}_x\text{Ga}_{4-x}$  an extended homogeneity region for  $0.4 \leq x \leq 1.0$ , was observed on the alloys annealed at 400 °C, only a very small field of existence was encountered at 400 °C for  $\text{YbCd}_x\text{Ga}_{4-x}$ , essentially centered around  $x = 0.5$ . The crystallographic data for the two alloy systems are summarized in Table 1. Adopting the  $z$ -parameter for atoms in 4e,  $z = 0.390$ , as derived from a Rietveld refinement of  $\text{YbAu}_{0.25}\text{Ga}_{3.75}$  [1], satisfactory agreement is found between observed and calculated X-ray powder intensities, assuming random Ga, Zn (Cd) atom distribution in the 4e position. Therefore, both phases belong to the  $\text{BaAl}_4$  type of structure.

The unit cell volume of  $\text{YbZn}_x\text{Ga}_{4-x}$  vs. increasing values of  $x$  shows a monotonic decrease for  $0.25 \leq x \leq 1.0$ . The extrapolated value for  $x = 0$  correlates nicely with the value for binary  $\text{YbGa}_4$  ( $V = 200.0 \text{ \AA}^3$  [1]). The sharp drop of the  $c$ -parameter from  $x = 0.25$  to  $x = 0.5$  may be seen as an increase in the covalent interatomic 4e-4e interaction, which starts to reduce again after the  $c$ -parameter passes through a minimum

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Table 1  
Crystallographic and magnetic data of  $\text{YbZn}_x\text{Ga}_{4-x}$  and  $\text{YbCd}_x\text{Ga}_{4-x}$  alloys (structure type  $\text{BaAl}_4$ , space group  $I4/mmm$ )

Composition	$a, \text{Å}$	$c, \text{Å}$	$V, \text{Å}^3$	$\mu_{\text{eff}}/\text{Yb}$ $\mu_{\text{B}}$	$\chi_{(5\text{K})} \cdot 10^6$ $\text{cm}^3 \text{g}^{-1}$	$\chi_{(300\text{K})} \cdot 10^7$ $\text{cm}^3 \text{g}^{-1}$
$\text{YbZn}_{0.25}\text{Ga}_{3.75}$	4.2428(5)	10.968(2)	197.4(1)	0.5	7.8	3.6
$\text{YbZn}_{0.5}\text{Ga}_{3.5}$	4.2290(7)	10.944(3)	195.7(1)	0.7	9.7	5.1
$\text{YbZn}_{0.75}\text{Ga}_{3.25}$	4.2141(6)	10.951(2)	194.5(1)	0.6	7.4	3.8
$\text{YbZnGa}_3$	4.2021(3)	10.955(1)	193.4(1)	0.7	9.5	5.5
$\text{YbCd}_{0.5}\text{Ga}_{3.5}$	4.2665(7)	10.965(2)	199.6(1)	0.5	2.5	7.8

(see Table 1,  $0.5 < x \leq 1.0$ ). The importance of this interaction phenomenon has been disclosed from band structure calculations of  $\text{BaAl}_4$  [3].

### 3.2. Magnetism

Magnetic susceptibilities of the compounds  $\text{YbT}_x\text{Ga}_{4-x}$  with  $T = \text{Zn, Cd}$  have been measured in the temperature range between 4.2 to 300 K. In general, the magnetic behavior is characterized by a nonmagnetic groundstate  $^1S_0$  of the ytterbium atom. However at low temperatures an upturn of the susceptibility curves (Curie-tail) is observed arising from small amounts of secondary phases. The calculated magnetic moment, according to a Curie–Weiss behaviour is smaller than  $0.7 \mu_{\text{B}}$  in all samples investigated. Hence less than 3% of impurity, where Yb adopts the magnetic  $^2F_{7/2}$  groundstate can cause the observed weak temperature dependency of  $\chi$  below 200 K.

### 4. Acknowledgement

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