# STRUCTURAL AND MAGNETIC PROPERTIES OF THE $\text{ErNi}_{1-x}\text{Cu}_x\text{Al SERIES}^*$

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The pseudo-ternary series  $\text{ErNi}_{1-x} \text{Cu}_x \text{Al}$  has been studied by means of X-rays and magnetization. Structure parameters exhibit a sudden "jump" between x = 0.5 and 0.6. The samples in the initial region with  $x \leq 0.2$  undergo two phase transitions and co-existence of F and AF ordering is possible. The samples with  $x \geq 0.4$  are already rather ferromagnets.

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## 1. Introduction

The RNiAl and RCuAl compounds belong to a large group of RTX compounds (R = rare earth, T = transition metal, X=p-metal). They crystallize in the ZrNiAl-type hexagonal structure, space group  $P\bar{6}2m$  (group No. 189), which consists of two types of layers. One of them contains all the rare-earth atoms and one third out of the transition-metal atoms. The p-metal and two out of the three transition metal atoms are included in the second, nonmagnetic layer separating the first one.

Interesting study was performed for  $\text{TbNi}_{1-x}\text{Cu}_x\text{Al series [1]}$ . In comparison to that case, the boundary stoichiometric concentration compounds in the  $\text{ErNi}_{1-x}\text{Cu}_x\text{Al series behave in the similar way with additional difference}$ in the directions of magnetic moments. ErNiAl is an antiferromagnet with

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magnetic moments arranged within the basal plane below 6.2 K [2], while ErCuAl is a simple ferromagnet with moments aligned in the *c*-axis direction with the ordering temperature of 6.8 K [3]. The change of magnetic-ordering type and also reorientation of the moments is expected.

## 2. Experimental

The polycrystalline samples have been prepared by arc-melting in monoarc furnace under protection of argon atmosphere. The initial materials consisted of pure elements with the purity of 99.9% for Er and Co, 99.995% for Ni and 99.999% for Cu and Al. All of the samples have been analyzed by X-ray powder diffraction at room temperature. Moreover, the quality of the selected samples has been proved by the microprobe experiment.

The DC-magnetization measurements have been performed using the PPMS instrument and on a SQUID magnetometer (both Quantum Design). The powder consisting of randomly oriented grains fixed by nonmagnetic glue in a small ampule has been used.

### 3. Results

The X-ray study of the  $\text{ErNi}_{1-x}\text{Cu}_x\text{Al}$  series at the room temperature showed that the lattice parameters radically "jump" between x = 0.5 and 0.6 while keeping the volume without any sudden change (Fig. 1).



Fig. 1. The concentration dependence of the lattice parameters in the  $\text{ErNi}_{1-x}\text{Cu}_x\text{Al}$  series, for several concentrations measured in low temperatures.

The lattice parameters measured on several samples (x=0.0; 0.4; 0.5; 1.0) at the temperatures down to 5 K show, that a and c follow the same type of concentration dependence at low temperatures as that at the room temperature.



Fig. 2. The low-temperature dependence of M/H in the  $\text{ErNi}_{1-x}\text{Cu}_x\text{Al}$  series. The full and open symbols represent zero-field cooled and field cooled regime, respectively. The values for different concentrations x are shifted for better view.



Fig. 3. Magnetization curves of the  $\text{ErNi}_{1-x}\text{Cu}_x\text{Al}$  series. In the right part the comparison with ErCuAl sample at 2 K — represented by empty symbols — is added.

By viewing magnetization data, one can think about two concentration regions. In the range of  $0.05 \le x \le 0.20$  there is possible co-existence of ferro- and antiferromagnetic alignment. The samples with x = 0.05, 0.1, 0.2exhibit two maxima in the temperature dependence of magnetization, what indicates two phase transitions (Fig. 2). These two points of phase transition have been confirmed by the heat capacity measurements. Also magnetization curves indicate the antiferromagnetic ordering. As it is seen from the double-curvature (Fig. 3), the ferromagnetic component is strengthening while increasing x.

The other region of the concentration of  $x \ge 0.4$  contains compounds exhibiting rather ferromagnetic behavior. They have a clear inflection point in low-temperature dependence of magnetization. Heat capacity measurements confirmed them to be phase-transition points. The magnetization curves contain already no double-curved behavior. And because of the similarity to the magnetization curve of the ErCuAl compound (under the same conditions) it is probable, that the  $\text{ErNi}_{1-x} \text{Cu}_x \text{Al compounds}$  with  $0.4 \le x \le 0.8$  are ferromagnets. The ferromagnetism of ErCuAl has been confirmed by neutron diffraction experiments [3].

### 4. Discussion

From the different alignment of magnetic moments in the ErNiAl and ErCuAl compounds, the change of magnetocrystalline anisotropy and the transition from antiferro- to ferromagnetic order within the  $\text{ErNi}_{1-x} \operatorname{Cu}_x \operatorname{Al}$  series is expected. The lattice "jump" can indicate the reorientation of the magnetic moments.

According to the evolution of the M vs. H dependencies within the concentration range, it seems that the ferromagnetic component becomes stronger while increasing the concentration parameter x. Also the temperatures of the second maximum in the low-temperature dependencies of M/H drives down with increasing x. So the change from antiferro- to ferromagnetic order happens consequently, without any sudden change.

If the hypothesis that compound with x = 0.40 is already ferromagnetic and if the sudden change in the lattice is connected with reorientation of the moments from the basal plane ( $x \le 0.5$ ) to the *c*-direction ( $x \ge 0.6$ ) it would mean that the concentrations  $0.5 \le x \le 0.6$  represent ferromagnetic compounds with magnetic moments within the basal plane. To confirm or negate this idea is the question for the neutron diffraction.

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