

# ELECTRONIC STRUCTURE AND MAGNETIC PROPERTIES OF $\text{Fe}_2\text{V}_{1-x}\text{Ti}_x\text{Al}$

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Recent transport, specific heat and magnetic susceptibility measurements indicated that the Heusler-type  $\text{Fe}_2\text{VAl}$  compound is a candidate for a 3d heavy-fermion system [1]. Experimental investigations have shown the large (compared to the normal metal) value of the low-temperature electronic specific heat coefficient  $\gamma$  and the semiconductor-like behavior of the resistivity in the paramagnetic state. Therefore,  $\text{Fe}_2\text{VAl}$  has been classified to the non-magnetic narrow-gap semiconductors (Kondo insulators) having similar properties as  $\text{FeSi}$  [2]. Another compound,  $\text{Fe}_2\text{TiAl}$  is weakly ferromagnetic [3],  $T_C \cong 120$  K in agreement to the Slater-Pauling behavior [4], however, its magnetic moment at  $T = 0$  is much smaller than expected.

In this presentation we discuss the results of our electronic structure investigations (XPS experimental results and LMTO calculations of the valence bands) for  $\text{FeV}_{1-x}\text{Ti}_x\text{Al}$  series of alloys. We also present the magnetic susceptibility measurements of  $\text{FeV}_{1-x}\text{Ti}_x\text{Al}$ . We try to understand the ground state properties of these alloys basing on the atomic disorder, which leads to the formation of the magnetic clusters.

[1] Y. Nishino et al., Phys. Rev. Lett. **79**, 1909 (1997).

[2] Z. Schlesinger et al., Phys. Rev. Lett. **71**, 1748 (1993).

[3] K.H.J. Buschow and P.G. Engen, J. Magn. Magn. Mater., **25**, 90 (1981).

[4] I. Galanakis et al., **66**, 174429 (2002).

9.7 cm

13.4 cm

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