

# ELECTRONIC AND MAGNETIC PROPERTIES OF *UCoAs<sub>2</sub>* COMPOUND

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The *UCoAs<sub>2</sub>* compound crystallizes in the tetragonal *HfCuSi<sub>2</sub>* - type structure with space group *P4/nmm* [1]. The compound orders ferromagnetically at 150K with spontaneous magnetic moment of about  $1.8\mu_B$ . The magnetic behaviour in *UCoAs<sub>2</sub>* exhibits a giant anisotropy in both ordered and paramagnetic region, which was interpreted as being caused predominantly by strong *f* – *d* hybridization and pronounced crystal field effect [1]. In this paper we present results of *ab – initio* band structure calculations based on the Full-Potential Local-Orbital Minimum-Basis Scheme (FPLO) [2].

[1] D. Kaczorowski, H. Noël, M. Potel, J. Alloys Compd. **302** (2000) 1

[2] K. Koepernik and H. Eschrig, Phys. Rev. B **59** (1999) 1743

13.4 cm

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9.7 cm