ELECTRONIC STRUCTURE OF URuGa₅ AND UIrGa₅

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As a part of a systematic study of the uranium isomorphic ternaries UTGa₅ with transition metals results of calculations of the electronic structure for T = Ru and Ir are presented. The compounds crystallize in the HoCoGa₅-type structure, space group P4/mmm, and (except for T = Ni, Pd and Pt) are not magnetically ordered [1]. The electronic structure was calculated by the full-potential LMTO method [2]. General features of the calculated densities of states show similarities to the ones for UCoGa₅ [3].

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