

Journal of Physics Condensed Matter 2008 vol.20 N45

Magnetic susceptibility of YbRh₂Si₂ and YbIr₂Si₂ on the basis of a localized 4f electron approach

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Abstract

We consider the local properties of the Yb³⁺ ion in the crystal electric field in the Kondo lattice compounds YbRh₂Si₂ and YbIr₂Si₂. On this basis we have calculated the magnetic susceptibility, taking into account the Kondo interaction in the simplest molecular field approximation. The resulting Curie-Weiss law and Van Vleck susceptibilities could be excellently fitted to experimental results over a wide temperature interval where thermodynamic and transport properties show non-Fermi-liquid behavior for these materials. © 2008 IOP Publishing Ltd.

<http://dx.doi.org/10.1088/0953-8984/20/45/455208>
