Low-temperature electrical resistivity of $Yb_{1-x}La_xRh_2Si_2$

F. Weickert*, P. Gegenwart, J. Ferstl, C. Geibel, F. Steglich

Max-Planck Institute for Chemical Physics of Solids, D-01187 Dresden, Germany

Abstract

Substitution of Yb-atoms with La in Yb_{1-x}La_xRh₂Si₂ leads to an increase of the Kondo temperature and suppression of antiferromagnetic order. Here, we present low-temperature ($T \ge 20 \text{ mK}$) electrical resistivity measurements on x = 0.05 and x = 0.1 single crystals at various magnetic fields. For x = 0.05, a linear temperature dependence is observed for $0 \le B \le 0.04 \text{ T}$ which turns into Landau Fermi liquid (LFL) behavior $\Delta \rho = AT^2$ at larger fields. For x = 0.1, LFL behavior is found already at 0.02 T. The evolution of A(B) suggests x = 0.05 and x = 0.1 to be located on the magnetically ordered and LFL side of the quantum critical point, respectively.

Keywords: YbRh2Si2; Quantum critical point

YbRh₂Si₂ is a heavy fermion system situated very close to a magnetic quantum critical point (QCP) [1]. A very weak antiferromagnetic (AF) ordering below $T_N = 70 \,\mathrm{mK}$ is suppressed by a tiny critical magnetic field of 0.06 T applied in the easy-plane perpendicular to the tetragonal caxis [2]. Pronounced non-Fermi liquid (NFL) effects like a stronger than logarithmic mass divergence and temperature over magnetic field scaling are incompatible with the itinerant spin-density-wave scenario for an AF QCP [3]. Furthermore, Hall-effect measurements suggest a dramatic change of the Fermi surface at the magnetic field tuned OCP [4]. The quantum critical fluctuations at magnetic fields well beyond 0.06 T have a dominating ferromagnetic component [5,6]. Therefore, it would be very interesting to compare the properties of such a magnetic field-driven OCP with the one at B = 0 in which the magnetic order is suppressed by an increase of the 4f-conduction electron hybridization q.

Hydrostatic pressure experiments on YbRh₂Si₂ have revealed an increase of T_N , as expected for Yb-systems, and allow to extrapolate to a negative critical pressure of -0.3(1) GPa corresponding to a tiny 0.2% volume expansion [7]. A volume expansion could either be achieved by the partial substitution of Si-atoms with the larger Ge or by La-substitution on the Yb-site in $Yb_{1-x}La_xRh_2Si_2$. The attempt to grow Ge-doped single crystals in which the AF ordering is completely suppressed has failed due to problems in sample preparation [8]. On the other hand, $Yb_{1-x}La_xRh_2Si_2$ single crystals have been successfully prepared for $x \leq 0.3$ and their characterization down to 0.4 K has revealed a systematic evolution of the characteristic maximum in the electrical resistivity suggesting an increase of q with increasing La-concentration [8]. Indeed the specific heat coefficient has been found to saturate below 1 K for x = 0.2 at $\gamma = 0.38 \text{ J/K}^2 \text{ mol}$, indicating heavy Landau Fermi liquid (LFL) behavior [8]. Here, we report electrical resistivity measurements on x = 0.05(residual resistivity ratio RRR = 12) and x = 0.1(RRR = 8) single crystals down to mK temperatures and at magnetic fields applied in the easy plane perpendicular to the *c*-direction.

Fig. 1(a) shows the low-temperature resistivity data of $Yb_{0.95}La_{0.05}Rh_2Si_2$. For $B \le 0.04 T$ and below 0.4 K a linear temperature dependence is observed, characteristic for NFL behavior. At larger fields, the low-*T* resistivity turns into a AT^2 dependence, as expected for a LFL. For $Yb_{0.9}La_{0.1}Rh_2Si_2$ (Fig. 1(b)) the zero-field data show power-law behavior $\Delta \rho \propto T^{\varepsilon}$ with $\varepsilon \approx 1.05$. This is in

^{*}Corresponding author. Tel.: +4935146462323;

fax: +49 351 4646 2360.

E-mail address: weickert@cpfs.mpg.de (F. Weickert).



Fig. 1. Low-temperature electrical resistivity of Yb_{1-x}La_xRh₂Si₂ for x = 0.05 (a) and x = 0.1 (b) at varying magnetic fields applied perpendicular to the *c*-direction. For clarity, the different curves at B>0 are shifted subsequently by 0.1 µΩcm. Solid, dashed and dotted lines represent $\Delta\rho \propto T^{\varepsilon}$ with $\varepsilon = 1$, $\varepsilon = 1.07$ and 2, respectively. Arrows indicate upper limit of T^2 behavior.

accordance with the trend observed at $T \ge 0.4$ K, that the resistivity exponent increases with increasing La-substitution [8]. T^2 behavior is observed at $B \ge 0.02$ T.

To get further information on the ground state behavior in the two systems, we analyze the magnetic field dependence of the coefficient A(B) which is a measure of the quasiparticle-quasiparticle scattering cross-section in the LFL state. In Fig. 2, A(B) values extracted from the data shown in Fig. 1 are compared with corresponding results for undoped YbRh₂Si₂ [2]. We use a double-log representation for A(B) since this allows to separate between three different cases: (i) for a divergence at a finite critical field B_c , like 0.06 T for the undoped system, an upwards curvature in A(B) is observed upon reducing the field towards the critical field, (ii) for a zero-field OCP, a straight line would indicate a power-law divergence towards $B_{\rm c} = 0$, whereas (iii) for a system located on the LFL side beyond the QCP, a saturation in A(B) is expected at small fields. As shown in Fig. 2, the A(B) data for the x = 0.05 system belong to case (i) and suggest a critical field slightly below 0.04 T, whereas the saturation of A(B)observed for x = 0.1 indicates a LFL ground state in the latter system. The strong reduction of A for fields beyond 10 T, observed for all three different systems, is related to the field-induced suppression of the Kondo effect [9].

A finite critical field for $Yb_{0.95}La_{0.05}Rh_2Si_2$ is also consistent with the temperature-field diagram determined from the field-dependence of the upper limit of T^2 behavior shown in the inset of Fig. 2. The linear extrapolation of this cross-over towards zero temperature indicates a critical



Fig. 2. Coefficient $A = \Delta \rho / T^2$ vs. field $B \perp c$ (on a double-log scale) for various concentrations x in Yb_{1-x}La_xRh₂Si₂. Inset: Temperature vs. field diagram with upper limit temperatures of $\Delta \rho = AT^2$ behavior (cf. arrows in Fig. 1), separating the non-Fermi liquid (NFL) regime from the LFL state. Dashed and dotted lines indicate linear and quadratic dependences, respectively.

field $B_c \approx 0.03 \text{ T}$ for this system. In Yb_{0.9}La_{0.1}Rh₂Si₂ the boundary of the NFL-LFL cross-over extrapolates towards a LFL state at B = 0.

The fact that an AF transition in the x = 0.05 system cannot be resolved in the temperature dependence of the resistivity might hint at a transition temperature below 20 mK. Since the signature of the AF transition is very sensitive to disorder and has neither been observed for the first generation of YbRh₂Si₂ single crystals ($\rho_0 = 3 \mu\Omega cm$) [1], nor in YbRh₂(Si_{0.95}Ge_{0.05})₂ ($\rho_0 = 5 \mu\Omega cm$) [3] it is also very likely that the disorder introduced by La-substitution ($\rho_0 = 5.6 \mu\Omega cm$ for x = 0.05) prevents the detection of the transition in the electrical resistivity.

To summarize, the analysis of the low-temperature electrical resistivity suggests a QCP in $Yb_{1-x}La_xRh_2Si_2$ for 0.05 < x < 0.1. Further low-*T* experiments are needed to determine the nature of the zero-field QCP in this system.

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