



Pressure-induced recovery of the Fermi-liquid state in the non-Fermi liquid material U_2Pt_2In

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Abstract

In the study of non-Fermi-liquid (NFL) phenomena in correlated metals, U_2Pt_2In is of special interest as it is one of the rare stoichiometric (undoped) materials that show NFL behaviour at ambient pressure. Here we report on the stability of the NFL phase with respect to hydrostatic pressure ($p \leq 1.8$ GPa). Electrical resistivity data under pressure, taken on a single-crystalline sample for a current in the tetragonal plane, show that T_{FL} , i.e. the temperature below which the Fermi-liquid T^2 -term is observed, increases with pressure as $T_{FL} \sim (p - p_c)$, where $p_c \approx 0$ is a critical pressure. This provides evidence for the location of U_2Pt_2In at an antiferromagnetic quantum critical point. © 2002 Elsevier Science B.V. All rights reserved.

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Among the U_2T_2X compounds, where T is a transition metal and X is In or Sn, U_2Pt_2In takes a special place, because it is a non-ordering heavy-electron compound with a strongly renormalised quasiparticle mass ($c/T = 0.41$ J/molU-K² at $T = 1$ K) [1]. Moreover, it shows pronounced non-Fermi liquid (NFL) behaviour. The NFL properties are summarised by (i) the specific heat varies as $c(T) \sim -T \ln(T/T_0)$ over almost two decades of temperature ($T = 0.1$ –6 K) [2], (ii) the magnetic susceptibility [3] shows a weak maximum at $T_m = 8$ K for a magnetic field along the c -axis (tetragonal structure), while it increases as $T^{0.7}$ when $T \rightarrow 0$ for a field along the a -axis, and (iii) the electrical resistivity obeys a power law T^α with $\alpha = 1.25 \pm 0.05$ ($T < 1$ K) and 0.9 ± 0.1 ($T \rightarrow 0$), for the current along the a - and c -axis, respectively [3,4]. Notice that muon spin relaxation experiments [4] have demonstrated the absence of (weak) static magnetic order at least down

to 0.05 K. In order to investigate the stability of the NFL phase with respect to pressure, we have carried out a high-pressure transport study on single-crystalline U_2Pt_2In [4]. The electrical resistivity, $\rho(T)$, was measured for a current, I , along the a - and c -axis, up to pressures of 1.8 GPa. The pressure effect is strongly current direction dependent, indicating a significant anisotropy of the Fermi surface. For $I \parallel c$ $\rho(T)$ increases with pressure and develops a relative minimum at low temperatures ($T_{min} \sim 4.8$ K at 1.8 GPa) [5]. Here we concentrate on the low-temperature data obtained for $I \parallel a$.

A single-crystalline batch of U_2Pt_2In was prepared by a modified mineralisation technique. Single-crystalline U_2Pt_2In forms in the tetragonal Zr_3Al_2 -type of structure (space group $P4_2/mn$). The residual resistivity ρ_0 amounts to $\sim 115 \mu\Omega\text{cm}$ for $I \parallel a$ [3]. With $\rho_{RT} \equiv \rho(300 \text{ K}) = 220 \mu\Omega\text{cm}$, a low residual resistance ratio $\rho_{RT}/\rho_0 = 1.9$ results. The resistivity was measured in the temperature interval 0.3–300 K, using a standard low-frequency four-probe AC-technique with a typical excitation current of $\sim 100 \mu\text{A}$. Pressures were exerted using a copper–beryllium clamp cell [4].

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At ambient pressure the resistivity ($I\parallel a$) shows, upon cooling below 300 K, a weak maximum near 80 K, followed by a steady decrease at lower temperatures. Pressure leaves $\rho(300\text{ K})$ unchanged, but results in an overall reduction of $\rho(T)$ below 300 K. In Fig. 1 we show the low-temperature ($T < 3.2\text{ K}$) data in a plot of $\Delta\rho \equiv \rho(T) - \rho_0$ versus T^2 . The pressure data ($p \geq 0.2\text{ GPa}$) show a steady reduction of $\Delta\rho$ with increasing pressure. The zero-pressure data are not in-line with this trend, as they fall in between the 0.2 and 0.6 GPa curves. This is possibly due to the fact that the zero-pressure data have been measured in a separate experiment in a different experimental set-up, with other voltage contacts on the sample and thus with a different geometrical factor. The relative accuracy in the resistivity values for the different experiments amounts to $\sim 10\%$, because of the uncertainty in the determination of the geometrical factor.

A most important conclusion that can be drawn from the data presented in Fig. 1, is the recovery of the Fermi-liquid (FL) $\Delta\rho \sim T^2$ law at moderate pressures. We have extracted T_{FL} , i.e. the upper temperature limit for the $\Delta\rho \sim T^2$ behaviour, by a least-squares fitting procedure. The pressure variation of T_{FL} is shown in Fig. 2. Within the error bars, the data are consistent with T_{FL} being a linear function of pressure. Such a linear pressure dependence has been proposed by Rosch [6] for itinerant antiferromagnets in the paramagnetic regime close to a magnetic quantum critical point (QCP). Within this magnetotransport theory T_{FL} is calculated as a function of the distance (measured by the pressure) to the QCP. T_{FL} varies initially as $T_{\text{FL}} = a_1(p - p_c)$ with a cross-over to $T_{\text{FL}} = a_2(p - p_c)^{1/2}$ at higher distances, where p_c is the pressure at the QCP. The pressure intervals in which the different laws are observed depend on the amount of disorder x in the system ($x \approx 1/\text{RRR} = \rho_0/\rho(300\text{ K})$). For $I\parallel a$, $x \sim 0.6$, which indicates that our sample is in a regime of intermediate disorder. In this regime, theory [6] predicts $T_{\text{FL}} = a_1(p - p_c)$, while at the QCP $\Delta\rho \sim T^\alpha$ with the NFL exponent $\alpha = 1.5$. The measured NFL exponent $\alpha = 1.25 \pm 0.05$ at $p_c = 0$ [4], is rather close to the predicted value. Correspondingly, the coefficient A of the FL T^2 term diverges upon approaching p_c . A strong increase of A can be deduced from the data in Fig. 1. A more detailed account of our high-pressure study can be found elsewhere [7].

The large residual resistivity value of our samples brings about the question whether the NFL behaviour in $\text{U}_2\text{Pt}_2\text{In}$ is due to Kondo disorder. However, the rapid recovery of the FL behaviour under pressure as probed by the resistivity data for $I\parallel a$ does not support this scenario. Since the compressibility is isotropic [4], pressure is expected to result in the further broadening of the distribution of Kondo-temperatures and thus the concurrent NFL behaviour would prevail.

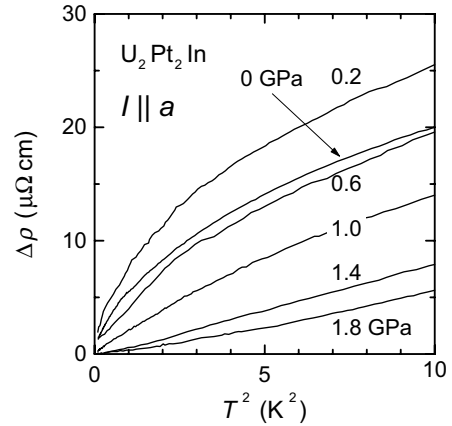


Fig. 1. Resistivity under pressure measured for a current along the a -axis of single-crystalline $\text{U}_2\text{Pt}_2\text{In}$ in a plot of $\Delta\rho$ versus T^2 .

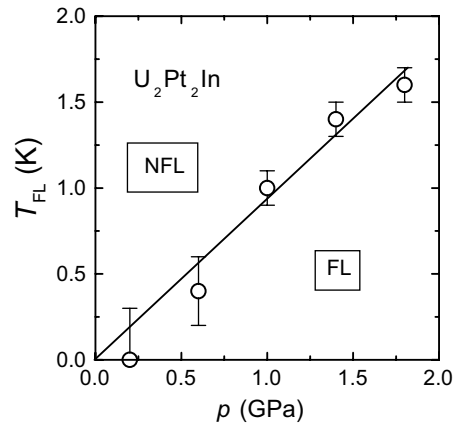


Fig. 2. T_{FL} versus pressure for single-crystalline $\text{U}_2\text{Pt}_2\text{In}$ determined from resistivity data for $I\parallel a$. The solid line separates the NFL regime from the FL regime. The linear behaviour is consistent with the presence of an antiferromagnetic quantum critical point at zero pressure.

In summary, we have measured the resistivity of the NFL material $\text{U}_2\text{Pt}_2\text{In}$ under pressure. For $I\parallel a$, the FL state is rapidly recovered. T_{FL} is a linear function of the pressure, $T_{\text{FL}} \sim (p - p_c)$, with $p_c \approx 0$. This provides evidence for the location of $\text{U}_2\text{Pt}_2\text{In}$ at an antiferromagnetic quantum critical point.

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