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# Electronic transport in $Ce_3Bi_4Pt_3$ : evidence for a temperature-dependent hybridization gap

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#### Abstract

We present an analysis of transport data from  $Ce_3Bi_4Pt_3$  to determine the energy gap temperature dependence  $E_g(T)$ .  $E_g(T)$  is a decreasing function of temperature. This result is consistent with the gap being a product of many-body correlations in which  $E_g(T)$  is scaled by the Kondo energy scale  $T_K$ .

Ce<sub>3</sub>Bi<sub>4</sub>Pt<sub>3</sub> is a narrow-gap Kondo insulator [1] in which the energy gap results from band hybridization driven by many-body effects. Mean-field treatments [2,3] of the Anderson lattice Hamiltonian suggest that the gap should be a decreasing function of temperature scaled by  $T_{\rm K}$ . In this note we present the results of an analysis of the energy gap  $E_{\rm g}(T)$  as determined from the resistivity  $\rho$  and thermoelectric power S of Ce<sub>3</sub>Bi<sub>4</sub>Pt<sub>3</sub>. The analysis indicates that  $E_{\rm g}$  is strongly temperaturedependent above 50 K, dropping close to zero at 300 K. As the gap shrinks above 100 K, single-impurity Kondo (SIK) scattering becomes evident in the resistivity. Upon comparing  $E_{\rm g}(T)$  with bulk modulus B and pressuredependent  $E_{\rm g}(P)$  data, a single energy scale ( $T_{\rm K}$ ) appears to govern both  $E_{\rm g}(T)$  and the thermal expansion  $\beta$ .

The  $\rho$  and S of Ce<sub>3</sub>Bi<sub>4</sub>Pt<sub>3</sub> are plotted in Fig. 1. With a temperature-independent gap and electron-phonon (EP) scattering, both data sets would appear as straight lines in Fig. 1; in both cases, the slope, and therefore  $E_g$ , drops with increasing temperature. Some curvature is also expected because this material is an extremely degenerate semiconductor (i.e.,  $E_{g} \sim k_{B}T$ ). Therefore, a careful quantitative analysis employing degenerate semiconductor statistics [4] is required to determine  $E_{\alpha}(T)$ . We model the system as involving two wide parabolic bands separated by  $E_{e}(T)$ ; this is a reasonable assumption for the lower and upper hybridized bands away from the Brillouin zone edge. In addition, we assume that EP scattering predominates. The results of a degenerate semiconductor analysis with these assumptions are presented in Fig. 2, along with a mean-field prediction [3].  $E_{\rm e}(T)$  is a decreasing function of temperature, as expected for a gap produced by many-body effects. Further, the gaps determined from  $\rho$  and S are equivalent; this is an indication that the assumption concerning scattering is valid since  $\rho$  and S are affected by EP scattering in fundamentally different ways. Above 100 K the uncertainty in the gap determined from S increases, but it still consistent with a shrinking gap. Despite considerable scatter in  $E_g$  above 100 K as determined from  $\rho$ , the gap appears to saturate to  $E_{g} = 70$  K. This is an artifact of the assumption that only EP scattering is present. It suggests that SIK scattering develops as the gap shrinks with increasing temperature.

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Fig. 1.  $\text{Log}_{10} \rho$  and S plotted as a function of 1/T. In both cases the slope drops with increasing temperature, indicating that  $E_g$  is a decreasing function of temperature.



Fig. 2.  $E_{g}(T)$  as determined from analyses of  $\rho$  (filled squares) and S data (open squares). The solid line is a mean-field prediction [3] based on the Anderson lattice Hamiltonian.

SIK scattering involves electrons in an Abrikosov-Suhl resonance of width  $T_{\rm K}$  near  $E_{\rm F}$ . With a wide gap centered at  $E_{\rm F}$  no scattering occurs. As the temperature is raised and  $E_{\rm g}$  falls well below  $T_{\rm K}$ , a partially gapped resonance will develop; therefore, SIK scattering should become evident as the temperature is raised. Preliminary high-temperature  $\rho$  measurements suggest that the gap disappears above 350 K and Ce<sub>3</sub>Bi<sub>4</sub>Pt<sub>3</sub> acts as an SIK metal.

The volume thermal expansion  $\beta$  of Ce<sub>3</sub>Bi<sub>4</sub>Pt<sub>3</sub> also displays a temperature dependence dominated by manybody effects [5]. If the temperature dependence of  $\beta$  and  $E_{g}$  result from hybridization driven by correlations they must be quantitatively interconnected. We can ensure that the gap scales with the thermal expansion by comparing  $E_{\bullet}(T)$  with bulk modulus and pressure-dependent  $g_{a_{P}} E_{c}(P)$  measurements. With a bulk modulus  $B \approx 900$  kbar, the thermal contraction at 4 K relative to 300 K is equivalent to 14.4 kbar pressure [5]. With  $\partial E_{o}/\partial P \approx +11 \pm 6 \,\mathrm{K/kbar}$  [1], the expected increase upon cooling from 300 K to 4 K is  $\Delta E_g = +160 \pm 90$  K; despite the crude nature of this calculation it is in good qualitative agreement with the data in Fig. 2. This quantitative consistence is a manifestation of the fact that all physical quantities controlled by many-body interactions in Ce<sub>3</sub>Bi<sub>4</sub>Pt<sub>3</sub> are fundamentally interrelated and their respective temperature dependencies are scaled by the Kondo energy scale  $T_{\rm K}$ .

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