Heat Transport in a Strongly Overdoped Cuprate: Fermi Liquid and Pure $d$-wave BCS Superconductor

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The transport of heat and charge in the overdoped cuprate superconductor Tl$_2$Ba$_2$CuO$_{6+\delta}$ was measured down to low temperature. In the normal state, obtained by applying a magnetic field greater than the upper critical field, the Wiedemann-Franz law is verified to hold perfectly. In the superconducting state, a large residual linear term is observed in the thermal conductivity, in quantitative agreement with BCS theory for a $d$-wave superconductor. This is compelling evidence that the electrons in overdoped cuprates form a Fermi liquid, with no indication of spin-charge separation.

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A fundamental question about the rich and baffling behavior of electrons in high-temperature superconductors is whether or not it can be understood in the framework of Fermi-liquid (FL) theory, the standard theory of electrons in solids. Several authors believe that when the concentration of electronic carriers in these cuprate materials is sufficiently low, as in the so-called underdoped region of the doping phase diagram, the basic excitations of the electron system are not the usual Landau quasiparticles characteristic of FL theory. In one class of proposals [1–3] for example, the electron is thought to fractionalize into a neutral spin-carrying excitation, called a “spinon”, and a spinless charge-carrying excitation, called a “holon” or “chargon”. However to this day, such “spin-charge separation” has not been confirmed experimentally. On the other hand, after 15 years of intensive research it is still not known with any certainty whether or not the ground state of cuprates is a Fermi-liquid in any region of the phase diagram. It is widely assumed that in the metallic-like overdoped regime at high carrier concentration FL theory does hold, but there is little solid evidence to support this lore.

In this Letter, we present the results of a study which show that strongly overdoped cuprates do not undergo spin-charge separation and their ground state is most likely a Fermi liquid. By measuring the transport of both heat and charge in the normal state at very low temperature, we were able to verify that one hole-doped cuprate in the overdoped regime obeys the Wiedemann-Franz (WF) law. This universal law is a robust signature of Fermi-liquid (FL) theory, the standard theory of electronic carriers concentration FL theory does hold, but there is still not known with any certainty whether or not the ground state of cuprates is a Fermi-liquid in any region of the phase diagram. It is widely assumed that in the metallic-like overdoped regime at high carrier concentration FL theory does hold, but there is little solid evidence to support this lore.

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tained previously [5], with $\rho_0 = 5.6 \, \mu\Omega \cdot \text{cm}$. Both heat and charge transport were measured using the same contacts, made by diffusing silver epoxy. A typical value for the contact resistance was $0.1 \, \Omega$ at 4 K. The thermal conductivity was measured down to below 100 mK with a standard one-heater two-thermometer technique in a dilution refrigerator. The magnetic field was applied along the $c$-axis. The geometric factor used to convert from resistance (electrical or thermal) to electrical resistivity $\rho$ or thermal conductivity $\kappa$ was set by requiring that $\rho(300 \, \text{K}) = 180 \, \mu\Omega \cdot \text{cm}$, the value obtained in previous studies of numerous crystals with the same doping level [5–8]. The uncertainty on this value is estimated at $\pm 10 \, \mu\Omega \cdot \text{cm}$.

The resistivity is shown in Fig. 1, for fields ranging from zero to above $H_{c2}(0)$. A slight positive magnetoresistance is observed, in agreement with previous work [7]. The resistivity below 30 K (and above $T_c$) is best fit by the function $\rho = \rho_0 + bT + cT^2$, with a substantial linear term (i.e. $bT > cT^2$ for $T < 15 \, \text{K}$). The fitting parameters are $\rho_0(H) = 5.84, 5.99$ and 6.15 $\mu\Omega \cdot \text{cm}$ at $H = 7, 10$ and 13 T, respectively, and $b = 0.064 \, \mu\Omega \cdot \text{cm} \cdot \text{K}^{-1}$, $c = 0.0054 \, \mu\Omega \cdot \text{cm} \cdot \text{K}^{-2}$ at 13 T. This unusual dependence was reported previously [8] and interpreted as “non-Fermi-liquid” behavior, in the sense that no linear term is expected in conventional FL theory. Deviations from the standard $T^2$ dependence have been observed in a number of heavy-fermion materials, for example, T$^{1.2}_1$ in CePd$_2$Si$_2$ below 20 K [9]. In these materials, this is associated with the proximity to a quantum critical point (QCP), where antiferromagnetic order sets in as a function of pressure or chemical composition. In the case of cuprates, the obvious QCP would be the onset of superconductivity at a critical concentration $p_c$ close to 0.3 hole/Cu atom, but a QCP has also been postulated to exist inside the superconducting region.

The thermal conductivity $\kappa$ is shown in Fig. 2. The data is plotted as $\kappa/T$ vs $T^2$ to separate the contribution of electrons from that of phonons, given that the asymptotic dependence of the former as $T \to 0$ is linear in $T$ while that of the latter is cubic. In other words, in Fig. 2, the electronic contribution is the residual linear term $\kappa_0/T$ given by the intercept of a linear fit with the $T = 0$ axis. The value of $\kappa_0/T$ obtained in this way is: 1.41, 2.76, 3.47, 3.75, 3.87, 3.90, 3.95, and 3.95 mW K$^{-2}$ cm$^{-1}$, at $H = 0, 1, 2, 5, 4, 5, 5, 7, 10, \text{and } 13 \, \text{T}$, respectively. As explained above, the uncertainty on the overall absolute value is approximately $\pm 5\%$. However, the relative uncertainty, e.g. between different fields, is much lower, around $1\%$. This high degree of reliability is due to the fact that in these samples electrons conduct much better than phonons, and hence the slope of $\kappa(T)/T$ in Fig. 2 is weak relative to the intercept. Note that at high fields, electrons scatter phonons very effectively and $\kappa(T)$ is entirely electronic below 1 K.

Fundamentally, the linear term in $\kappa$ at $T = 0$ reveals the presence of fermionic excitations in the electron system. We can then ask whether these excitations carry charge. This question can only be addressed in the absence of any superfluid that can also carry charge, which amounts to testing the WF law in the normal state. This law is one of the most fundamental properties of a Fermi liquid, reflecting the fact that the ability of a quasiparticle to transport energy is the same as its ability to transport charge, provided it cannot lose energy through collisions. It states that the heat conductivity $\kappa$ and the electrical conductivity $\sigma$ of a metal are related by a universal constant:

$$\frac{\kappa}{\sigma T} = \frac{\pi^2}{3} \left( \frac{k_B}{\epsilon} \right)^2 \equiv L_0$$

(1)

where $T$ is the absolute temperature, $k_B$ is Boltzmann’s constant and $L_0 = 2.44 \times 10^{-8} \, \text{W} \, \Omega \, \text{K}^{-2}$ is Sommerfeld’s value for the Lorenz ratio $L \equiv \kappa/\sigma T$. Theoretically, electrons are predicted to obey the WF law at $T \to 0$ in a wide range of environments: in both three or two dimensions (but not strictly in one dimension), for any strength of disorder and interaction [10], scattering and magnetic field [11]. Experimentally, the WF law does appear to be universal at $T \to 0$: until recently, no material had been reported to violate it. The first exception was found in optimally-doped Pr$_{2-x}$Ce$_x$CuO$_4$ (PCCO), an electron-doped cuprate [12].

It is in general difficult to test the WF law in cuprate superconductors because of their high upper critical
fields. In our crystals, the superconductivity has completely vanished by 13 T, at which field we find $\kappa_0/T = 3.95 \pm 0.04 \text{ mW K}^{-2} \text{ cm}^{-1}$ and $\rho_0 = 6.15 \pm 0.03 \mu\Omega \text{ cm}$, so that $L = \rho_0 \kappa_0/T = 0.99 \pm 0.01 L_0$, in perfect agreement with the WF law. Note that the Lorenz ratio does not suffer from the 5% uncertainty associated with the geometric factor, as both transport measurements are performed using the same sample with the same contacts. The error bars are therefore on the order of 1%.

In Fig. 2, the transport of heat and charge are compared directly by reproducing the charge conductivity at 13 T from Fig. 1. This is done by plotting $L_0/\rho(T)$ vs $T$ using the fit to the 13 T data for $\rho(T)$ (inset of Fig. 1). The charge conductivity $L_0\sigma(T)$ is seen to equal the heat conductivity $\kappa(T)/T$ at 13 T.

The basic implication of this result is that the fermions which carry heat also carry charge $e$ and are therefore indistinguishable from standard Landau quasiparticles. In particular, there is no evidence of any spin-charge separation. Indeed, if electrons were to fractionalize into neutral spin-carrying fermions (spinons) and charged bosons (chargons) [3], there would be no reason to expect the WF law to hold, as the heat-carrying fermions would not take part in the transport of charge. This result therefore imposes a constraint on theories of spin-charge separation (SCS): the critical hole concentration $p_{SCS}$ at which electron fractionalization starts to occur is not the QCP where superconductivity starts to occur (on the overdoped side of the phase diagram), but can only be lower. In other words, any hypothetical onset of SCS must obey $p_{SCS} < 0.26 < p_c$. It therefore appears that the mechanism for superconductivity in this overdoped region of the phase diagram is not the condensation of charge-$e$ bosons, but most likely Cooper pairing. Note that (barring any profound electron-hole asymmetry) this conventional picture is expected to break down with underdoping, as suggested by the violation of the WF law in PCCO near optimal doping [12].

Although the standard FL description fails, as revealed by the non-quadratic $T$ dependence of $\rho(T)$, the basic nature of the electronic excitations in the limit of zero energy is that of Landau FL quasiparticles. (A similar situation is seen in heavy-fermion materials [13].)

In the absence of a magnetic field, there is a large residual linear term in the thermal conductivity of TI-2201, namely $\kappa_0/T = 1.41 \text{ mW K}^{-2} \text{ cm}^{-1}$. A similar term has also been observed in other hole-doped cuprates, albeit at optimal doping, where it is much smaller: $\kappa_0/T = 0.14$, 0.15 and 0.11 mW K$^{-2}$ cm$^{-1}$, in Y-123 [14], Bi-2212 [15,16] and LSCO [17], respectively. Within BCS theory applied to a $d$-wave superconductor, this residual heat conduction is expected, arising from zero-energy quasiparticles induced by impurity scattering near the nodes in the $d_{x^2-y^2}$ gap function. In the clean limit, where the scattering rate $\Gamma \ll k_BT_c/h$, it is universal (in the sense that it is independent of $\Gamma$) and it depends only on the ratio of the two quasiparticle velocities ($v_F$ and $v_2$) which govern the Dirac-like spectrum of nodal quasiparticles, $E = \hbar \sqrt{v_F^2 \vec{k}_F^2 + v_2^2 \vec{k}_2^2}$ [18]:

$$\frac{\kappa_0}{T} = \frac{k_B^2 n}{3h} \left( \frac{v_F}{v_2} + \frac{v_2}{v_F} \right)$$

where $n$ is the number of CuO$_2$ planes per unit cell of height $c$ (along the c-axis), and $k_1$ and $k_2$ are unit vectors pointing in directions normal and tangential to the Fermi surface at the node, respectively. In other words, $v_F$ is the Fermi velocity in the nodal direction and $v_2$ is proportional to the slope of the gap at the node, $d\Delta/d\phi = \hbar k_F v_2$, with $k_F$ the Fermi wavevector.

Applying Eq. 2 to TI-2201, for which $n = 2$ and $c = 23.2 \text{ Å}$, we get $v_F/v_2 = 270$. A rough estimate using Fermi surface parameters typical of cuprates, namely $v_F = 2.5 \times 10^7 \text{ cm/s}$ and $k_F = 0.7 \text{ Å}^{-1}$ (the values measured in Bi-2212 [19]), and the simplest $d$-wave gap function, $\Delta = \Delta_0 \cos \phi$, with the weak-coupling relation for a $d$-wave superconductor, $\Delta_0 = 2.14 k_BT_c$, gives $v_F/v_2 = 210$. This shows that the magnitude of $\kappa_0/T$ is in good agreement with the simplest BCS analysis.

It should be recognized that even though the mean free path in these samples is rather long (in the range 500 – 1000 Å [8]), the scattering rate $\Gamma$ is not small compared to $T_c$. It may be estimated using the standard expression for the normal state conductivity: $\kappa_N/T = \frac{1}{3} \gamma_N v_F^2 \tau$, where $\gamma_N$ is the specific heat coefficient and...
\[ \tau = 1/(2\Gamma). \] With \( \gamma_N \approx 3 \text{ mJ K}^{-2} \text{ mole}^{-1} \) [6] and \( v_F = 2.5 \times 10^7 \text{ cm/s} \), one gets \( \hbar \gamma \approx 0.4 k_BT_c \). At finite \( \Gamma \), corrections to Eq. 2 give an increase in \( \kappa_0/T \) [20]. Assuming \( \Delta_0 = 2.14 k_BT_c \), the correction for \( \hbar \Gamma/k_BT_c \approx 0.4 \) is by a factor of approximately 1.5 [20]. Thus the correct value of \( v_F/v_2 \) is probably closer to 270/1.5 = 180 [21].

It will be interesting to investigate the doping dependence of \( \kappa_0/T \) as a way of measuring the dependence of the gap function on carrier concentration, via \( v_2 \). In the absence of further data on Ti-2201, we may compare with optimally-doped Y-123 (\( T_c = 93 \text{ K} \)) or Bi-2212 (\( T_c \approx 90 \text{ K} \)), for which \( v_F/v_2 = 14 \) and 19, respectively [15]. (The value of 19 for Bi-2212 agrees very well with the value of 20 obtained from ARPES measurements of \( v_F \) and \( v_2 \) separately [15,19].) Under the assumption of a doping independent \( v_F \), verified in both Bi-2212 [19] and LSCO [22], one immediately sees that \( v_2 \), or the magnitude of the gap (near the nodes), scales roughly with \( T_c \). This strongly suggests that the standard BCS relation between gap magnitude and transition temperature, \( \Delta_0 \propto T_c \), holds in the overdoped regime. This is in striking contrast with what is found in the underdoped region of the phase diagram. Indeed, our measurements on underdoped Y-123 and LSCO [23] reveal that \( v_F/v_2 \) decreases as \( T_c \) is reduced by underdoping (see also [24]).

Several authors have proposed the existence of a QCP within the superconducting dome in the phase diagram of cuprates, either as a theoretical prediction to explain the diagram itself or as suggested in various experiments. Its location is usually taken to be near optimal doping, in the neighbourhood of \( p = 0.2 \). If it is associated with a change in the symmetry of the superconducting order parameter, Vojta et al. have argued that the most likely scenario is a transition from a pure \( d_{x^2−y^2} \) state to a complex order parameter of the form \( d_{x^2−y^2} + ix \), where \( x \) can have either \( s \) or \( d_{xy} \) symmetry [25]. Dagan and Deutscher have recently reported a split zero-bias anomaly in their tunneling on Y-123 thin films as soon as the material is doped beyond optimal doping, a feature which they attribute to the appearance of a complex component to the order parameter in the bulk [26]. The presence of a subdominant component \( ix \) in the order parameter causes the nodes to be removed, as the gap can no longer go to zero in any direction. The observation of a residual linear term in the thermal conductivity, a direct consequence of nodes in the gap, therefore excludes the possibility of any such subdominant order parameter. (More precisely, since our measurement goes down to 100 mK, it puts an upper bound on the magnitude of \( |x| \) relative to \( |d_{x^2−y^2}| \) at about 0.5 \%). Moreover, there is no subdominant order parameter in Ti-2201 at optimal doping [4]. In other words, if there truly is a QCP between optimal doping at \( p \approx 0.16 \) and the critical point \( p_c \approx 0.3 \), it does not appear to be associated with the onset of a complex component in the order parameter.

In summary, the low-temperature transport properties of Ti-2201 with \( T_c = 15 \text{ K} \) show that spin-charge separation does not occur in strongly overdoped cuprates. The normal state at \( T \to 0 \) satisfies the Wiedemann-Franz law perfectly, demonstrating that the only electronic excitations carrying heat and charge are Landau quasiparticles. The superconducting state obeys BCS theory in that the residual heat conduction is of the expected magnitude for a \( d \)-wave gap and the dependence of the low-energy spectrum on doping strongly suggests that the gap scales with \( T_c \) in the conventional way. Finally, the possibility of a sub-dominant order parameter (\( ix \)) is ruled out.

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