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## LETTER TO THE EDITOR

## Remarks on the normal-state resistivity of high- $T_{\rm c}$ superconductors

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Abstract. The validity of semiclassical transport theory for two-dimensional Fermi and Bose conductors is analysed. It is pointed out that if the high- $T_c$  materials are Bose conductors, this theory breaks down. It is also argued that the regime of 'resistivity saturation' in this case occurs at low temperatures with a resistivity that increases linearly with temperature.

One of the characteristic features of the new Cu oxide superconductors is the unusual temperature dependence of the normal-state resistivity  $\rho$ . Measurements show that  $\rho$  parallel to the *ab*-plane is nearly linear in T for  $T > T_c [1-5]$ .

Several theoretical explanations have been offered for this behaviour [6–11]. Common to all these proposals is the use of semiclassical transport theory which leads, essentially, to the Drude formula for  $\rho$  [12]:

$$\rho = m/ne^2\tau \tag{1}$$

where *n* is the carrier density, *m* its mass and  $\tau$  the collision time. If one uses (1) to estimate  $\tau$ , taking  $\rho = \alpha T$ , with  $\alpha = 1 \,\mu\Omega \,\mathrm{cm}\,\mathrm{K}^{-1}$  [1],  $m \simeq (1-5)m_{\rm e}$  ( $m_{\rm e} =$  electron mass and  $n \simeq 3 \times 10^{21} \,\mathrm{cm}^{-3}$ ) one finds  $\hbar/\tau \ge \kappa T$ . Models that give  $\hbar/\tau \propto \kappa T$  are: two-dimensional Fermi conductors, in which case the linear *T*-dependence comes from electron-phonon scattering [6-9] and from Umklapp electron-electron scattering [7, 10], and two-dimensional Bose conductors in which the charge-carrying quasiparticles are bosons. In this case  $\hbar/\tau \propto \kappa T$  comes from boson-phonon scattering [11].

The conclusion that  $\hbar/\tau \ge \kappa T$  suggests that in these systems the quasiparticles are not well defined excitations, in the sense that the linewidth  $\Gamma \sim \hbar/\tau$  is greater than the excitation energy  $\sim \kappa T$ . One of the possible consequences of  $\Gamma \ge \kappa T$  is that (1) may be invalid [12].

This problem was studied a long time ago in connection with the theory of normal metals. In this case  $\Gamma \ge \kappa T$  occurs for  $T \ge \Theta_D$  ( $\Theta_D =$  Debye temperature) due to electron-phonon interactions. It was shown that (1) remains valid as long as  $\Gamma < \varepsilon_F$ ,  $\varepsilon_F$  being the Fermi energy [13]. Introducing the electron mean free path  $l = v_F \tau$ ,  $v_F = \hbar k_F/m$  being the Fermi velocity and  $k_F$  being the Fermi wavevector, this condition may be written as  $\Gamma \sim \hbar/\tau = \hbar v_F/l < v_F \sim \hbar v_F k_F$ , or  $l \ge k_F^{-1}$ . This result was first obtained by Ioffe [14] and by Ioffe and Regel [15]. They pointed out that semiclassical transport theory must break down when the characteristic linear dimension of the wavepacket

associated with the quasiparticle  $\sim \lambda$  ( $\lambda =$  typical quasiparticle wavelength) exceeds *l*. Thus (1) remains valid as long as  $l > \lambda$ . In ordinary metals  $\lambda \sim k_F^{-1} \sim a$ , *a* being the interatomic distance, and  $l > \lambda$  reads l > a.

The same condition,  $l > \lambda$ , must also dictate the region of applicability of (1) to Bose conductors. In the normal state,  $\lambda$  is the thermal De Broglie wavelength  $\lambda_T = \sqrt{2m\kappa T/\hbar^2}$  and the particle characteristic velocity is  $v_T = \sqrt{2\kappa T/m}$ . Thus (1) is valid if  $l = v_T \tau > \lambda_T$  or  $\hbar/\tau < \kappa T$ , a condition very different from that for ordinary conductors! If the Cu oxide superconductors are Bose conductors this restriction leads to a contradiction. As mentioned above, comparing  $\rho$  measured experimentally with (1) one finds  $\hbar/\tau \ge \kappa T$  which in turn violates  $l > \lambda$  rendering (1) invalid. This also suggests that scattering processes that give rise to  $\hbar/\tau \propto \kappa T$  do not explain  $\rho \propto T$ , since, because  $l > \lambda$  is violated,  $\rho$  is no longer proportional to  $\hbar/\tau$ .

The expression for  $\rho$  valid when  $l < \lambda$  is not known. What is generally accepted is the Ioffe-Regel criterion [16]. It states that  $l < \lambda$  is impossible. This means that as *l* decreases below  $\lambda$ ,  $\rho$  no longer increases with *l* but assumes the value  $\rho_{sat} \sim$  $mv/ne^2\lambda$ , obtained from (1) by setting  $l = v\tau \sim \lambda$  (v = quasiparticle characteristic velocity). Experimental evidence for this type of behaviour is found in several metals [17, 18]. What is observed is that  $\rho$  for these systems stops increasing with rising temperature. This phenomenon, called 'resistivity saturation' [18], is interpreted as evidence that  $l < \lambda$  in the temperature range where  $\rho$  no longer increases with T [17, 18].

This simple physical reasoning suggests that  $\rho$  can be written as

$$\rho = (m/ne^2\tau)f(\lambda/l) \tag{2}$$

where f(x) is a well behaved function of x. For  $\lambda/l \le 1, f \to 1$ , so (2) reduces to (1). For  $\lambda/l \ge 1$  the Ioffe-Regel criterion gives  $f \to \gamma l/\lambda$ ,  $\gamma$  being a constant ~1, so

$$\rho \to \rho_{\rm sat} = \gamma (mv/ne^2\lambda). \tag{3}$$

For Fermi systems  $v = v_F$ ,  $\lambda \sim k_F^{-1} \sim a \operatorname{so} \rho_{\operatorname{sat}} \sim \gamma m v_F/ne^2 a$ . In metals where resistivity saturation is observed equation (3) correctly predicts the order of magnitude of the maximum  $\rho$  [19].

Application of these ideas to Bose conductors gives  $\lambda/l \sim \lambda_T/v_T \tau \sim \hbar/\tau \kappa T$ . Thus, according to (2),

$$\rho^{\rm B} = (m/ne^2\tau)f(\hbar/\tau\kappa T). \tag{4}$$

For  $\hbar/\tau\kappa T < 1$ ,  $\rho^{B}$  reduces to (1). For  $\hbar/\tau\kappa T > 1$ ,  $\rho^{B} \rightarrow \rho^{B}_{sat}$ , where

$$\rho_{\rm sat}^{\rm B} = \gamma m v_{\rm T} / n e^2 \lambda_{\rm T} = \gamma (m \kappa T / n e^2 \hbar). \tag{5}$$

Thus, for Bose conductors in the regime  $l < \lambda$ , 'resistivity saturation' occurs not at a constant  $\rho$  but at  $\rho_{sat}^{B}$ , equation (5), which increases linearly with T.

If the Cu oxide superconductors are two-dimensional Bose conductors, then, due to boson-phonon scattering [11],  $\hbar/\tau = \Lambda \kappa T$  ( $\Lambda$  is a constant that depends on the strength of the boson-phonon interaction). Thus, for  $\Lambda > 1$ ,  $\rho$  is given by (5). Substituting in this equation the values appropriate for the oxide superconductors,  $n = 3 \times 10^{21} \text{ cm}^{-3}$ ,  $m \sim 5m_{\rm e}$ , and assuming  $\gamma \sim 1$  one finds  $\rho_{\rm sat}^{\rm B} = \alpha T$  with  $\alpha \sim 1 \,\mu\Omega \,{\rm cm} \,{\rm K}^{-1}$ . This  $\rho_{\rm sat}^{\rm B}$  is of the same order of magnitude as  $\rho$  in these materials and has the same temperature dependence.

In the RVB theory [20]  $\hbar/\tau$  has an additional contribution from holon-spinon scattering that varies as  $T^{3/2}$  [21]. Whether or not this is the most important contribution

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depends on the relative strengths of the holon-phonon and holon-spinon contributions. In any case, as long as  $\hbar/\tau > 1$ , the above conclusion holds true.

In conclusion then, the simple physical arguments given above suggest that if the normal state of Cu oxide superconductors is a two-dimensional Bose conductor, the linear temperature dependence of  $\rho$  arises from resistivity saturation, rather than from an intrinsic scattering mechanism.

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