

## Hidden $T$ -Linear Scattering Rate in $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$ Revealed by Optical Spectroscopy

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The optical properties of  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$  have been determined in the normal state for a number of temperatures over a wide frequency range. Two Drude terms, representing two groups of carriers with different scattering rates ( $1/\tau$ ), well describe the real part of the optical conductivity  $\sigma_1(\omega)$ . A “broad” Drude component results in an incoherent background with a  $T$ -independent  $1/\tau_b$ , while a “narrow” Drude component reveals a  $T$ -linear  $1/\tau_n$  resulting in a resistivity  $\rho_n \equiv 1/\sigma_{1n}(\omega \rightarrow 0)$  also linear in temperature. An  $\arctan(T)$  low-frequency spectral weight is also strong evidence for a  $T$ -linear  $1/\tau$ . A comparison to other materials with similar behavior suggests that the  $T$ -linear  $1/\tau_n$  and  $\rho_n$  in  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$  originate from scattering from spin fluctuations and hence that an antiferromagnetic quantum critical point is likely to exist in the superconducting dome.

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Over the past several decades, it has been observed that the electrical resistivity  $\rho$  of some strongly correlated materials increases linearly with temperature ( $T$ -linear  $\rho$ ), particularly in the vicinity of an antiferromagnetic quantum critical point (QCP), a striking deviation from Landau’s Fermi-liquid description of metals. This anomalous  $T$ -linear  $\rho$ , extensively studied in the high- $T_c$  cuprate superconductors [1–4], organic Bechgaard salts [5,6], as well as heavy-fermion metals [7–10], may be intimately related to the emergence of superconductivity [4,5,11]. It is generally believed that in proximity to the antiferromagnetic QCP, spin fluctuations are so strong that the scattering process of quasiparticles is severely modified, inducing non-Fermi-liquid behavior. Such spin fluctuations may be responsible for the pairing of electrons in high- $T_c$  superconductors [12–14].

The presence of the  $T$ -linear  $\rho$  and a QCP in the newly discovered iron-based superconductors is highly desired since superconductivity arises in the vicinity of the spin-density-wave (SDW) instability [15,16]. Up to this point,  $T$ -linear  $\rho$  has been observed by transport, arousing considerable effort to describe it as evidence of possible quantum criticality in iron pnictides, especially in the “122” family [5,17–23]. However, unlike the high- $T_c$  cuprates, iron pnictides fall into the category of multiband materials [24,25]. Up to five Fe  $3d$  bands crossing the Fermi level contribute to the Fermi surface, leading to the presence of three holelike Fermi pockets at the  $\Gamma$  point and two electronlike pockets at the corners of the Brillouin zone. The scattering rate  $1/\tau$  and the response of quasiparticles to the external electrical field may vary considerably in different Fermi pockets [19,26–28]. As a result, the transport properties become extremely complicated in such a system and the question of whether a  $T$ -linear  $\rho$  in

iron pnictides originates from multiband effects or the presence of a QCP makes it inadequate to investigate the transport properties alone in search of a non-Fermi-liquid behavior and evidence for possible QCP [28,29].

Although many optical studies on iron pnictides have been reported [30–32], the above issue has never been touched due to the absence of detailed  $T$ -dependent optical data. In this Letter we address this issue by studying the detailed  $T$  dependence of the optical conductivity and low-frequency spectral weight in  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$ . The low-frequency optical conductivity is described by two Drude terms: a broad Drude with a large  $1/\tau_b$  that is basically  $T$  independent alongside a narrow one with a small  $T$ -linear  $1/\tau_n$  that reveals a  $\rho_n \equiv 1/\sigma_{1n}(\omega \rightarrow 0) \propto T$ . The low-frequency spectral weight increases with cooling, following an  $\arctan(T)$  dependence, which is demonstrated to be a clear signature of  $T$ -linear  $1/\tau$ . A comparison with similar behavior found in other materials attributes the  $T$ -linear  $1/\tau_n$  and  $\rho_n$  in  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$  to spin fluctuation scattering and the presence of a QCP in the superconducting dome.

High quality  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$  single crystals were grown using a self-flux method [19]. The inset of Fig. 1 shows the dc resistivity of  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$  as a function of temperature. The  $\rho(T)$  curve is characterized by a sharp superconducting transition at  $T_c \approx 39$  K and a tendency to saturation at room temperature. A visible change in slope occurs at about 175 K.

The near-normal incident reflectance  $R(\omega)$  has been measured for light polarized in the  $a$ - $b$  plane using FTIR spectrometers and an *in situ* evaporation technique [33]. Data from  $\approx 20$ – $12\,000$   $\text{cm}^{-1}$  were collected at 18 different temperatures from 5 to 300 K on a freshly cleaved surface. The visible and UV range ( $10\,000$ – $55\,000$   $\text{cm}^{-1}$ )

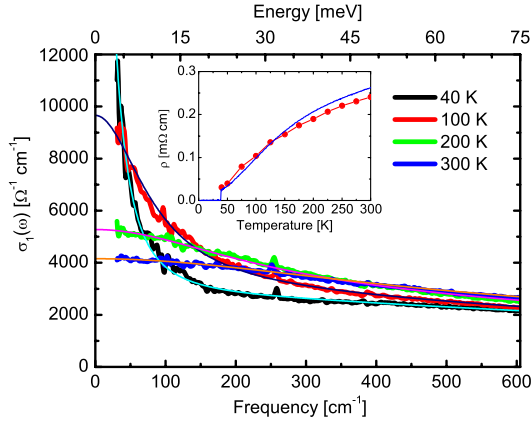


FIG. 1 (color online). Main panel: Thick solid lines are  $ab$  plane  $\sigma_1(\omega)$  of  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$  at different temperatures in the normal state. The smooth lines through the data are fits with the Drude-Lorentz model. Inset: dc resistivity as a function of temperature from transport measurement (solid curve) and the values derived from the zero frequency extrapolation of  $\sigma_1(\omega)$  (solid circles).

$R(\omega)$  was measured at room temperature with an AvaSpec-2048  $\times$  14 fiber optic spectrometer. The real part of the complex optical conductivity  $\sigma_1(\omega)$  is determined from  $R(\omega)$  via a Kramers-Kronig analysis. A Hagen-Rubens form ( $R = 1 - A\sqrt{\omega}$ ) is used for the low-frequency extrapolation. At high frequencies,  $R(\omega)$  is assumed to be constant to 40 eV, above which a free-electron response ( $\omega^{-4}$ ) is used.

The main panel of Fig. 1 shows  $\sigma_1(\omega)$  at four selected temperatures in the normal state (thick solid lines); all the spectra exhibit the well-known Drude-like metallic response. In order to quantitatively analyze the optical data, we fit  $\sigma_1(\omega)$  to the Drude-Lorentz model,

$$\sigma_1(\omega) = \frac{2\pi}{Z_0} \left[ \sum_k \frac{\Omega_{p,k}^2}{\tau_k(\omega^2 + \tau_k^{-2})} + \sum_j \frac{\gamma_j \omega^2 \Omega_j^2}{(\omega_j^2 - \omega^2)^2 + \gamma_j^2 \omega^2} \right], \quad (1)$$

where  $Z_0$  is the vacuum impedance. The first term describes a sum of free-carrier Drude responses, each characterized by a plasma frequency  $\Omega_p^2 = 4\pi n e^2 / m^*$ , where  $n$  is a carrier concentration and  $m^*$  is an effective mass, and a scattering rate  $1/\tau$ . The second term corresponds to a sum of Lorentz oscillators characterized by a resonance frequency ( $\omega_j$ ), a linewidth ( $\gamma_j$ ), and an oscillator strength ( $\Omega_j$ ). This Drude-Lorentz model is also used to determine the dc properties of the system [34].

As shown in Fig. 2,  $\sigma_1(\omega)$  at 150 K is described by a broad Drude with a large scattering rate  $1/\tau_b \approx 936 \text{ cm}^{-1}$ , and a narrow Drude with a small scattering rate  $1/\tau_n \approx 158 \text{ cm}^{-1}$  and an overdamped Lorentzian term. The linear superposition of these three components gives a very good description to  $\sigma_1(\omega)$  up to  $10\,000 \text{ cm}^{-1}$  at all the measured temperatures in the normal state. Fits for other temperatures are selectively

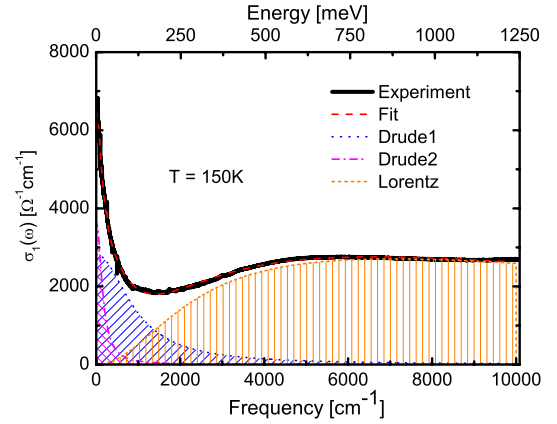


FIG. 2 (color online). The black solid curve is the measured  $\sigma_1(\omega)$  of  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$  at 150 K. The red dashed line through the data is the fit which is decomposed into a broad Drude (blue dotted line), a narrow Drude (pink dotted-dashed line), and a Lorentz (orange short-dashed line) term.

shown in the main panel of Fig. 1 as smooth thin lines through the corresponding data. The inset of Fig. 1 compares the optical estimate for the dc resistivity  $\rho \equiv 1/\sigma_1(\omega \rightarrow 0)$  (solid circles) to the transport measurements (solid line).

The two-Drude fit indicates the existence of two groups of carriers with different  $1/\tau$ 's in  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$ , which was first pointed out by Wu *et al.* in various iron pnictides [35]. The disparity of the  $1/\tau$ 's in different bands is also supported by both theoretical calculations [36] and measurements using other techniques on similar materials [19,26–28]. Tu *et al.* suggest that it is appropriate to describe the broad Drude term as bound excitations [37], because the mean free path  $l = v_F \tau$  ( $v_F$  is the Fermi velocity) associated with the broad Drude is less than the shortest interatomic spacing, violating the Mott-Ioffe-Regel condition [38]. In  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$ , the average Fermi velocities of the electron and hole pockets are estimated to be  $v_F^e \approx 0.40 \text{ eV \AA}$  and  $v_F^h \approx 0.36 \text{ eV \AA}$  [39], respectively. Furthermore, it is reported that in iron pnictides holes have a larger  $1/\tau$  than electrons [19,26–28]. If we associate the broad Drude component ( $1/\tau_b \approx 936 \text{ cm}^{-1}$ ) with the hole pockets, a mean free path of  $l_h \approx 3 \text{ \AA}$  is obtained. This value is close to the lattice parameter  $a \approx 4 \text{ \AA}$  of the 122 family compounds and probably too small for coherent transport. Since the broad Drude only produces an incoherent,  $T$ -independent, background contribution to the total  $\sigma_1(\omega)$ , the nature of the broad Drude component (whether or not bound excitations) does not affect our analysis on the temperature dependence of  $\sigma_1(\omega)$  and low-frequency spectral weight.

The temperature dependence of the two Drude components provides information about the nature of the two different types of carriers in this material. Figure 3 shows the  $T$  dependence of the Drude parameters from our fits. The subscripts  $n$  and  $b$  stand for the narrow and broad Drude terms, respectively. Figure 3(a) shows the  $T$  dependence of the plasma frequencies of the two Drude terms.

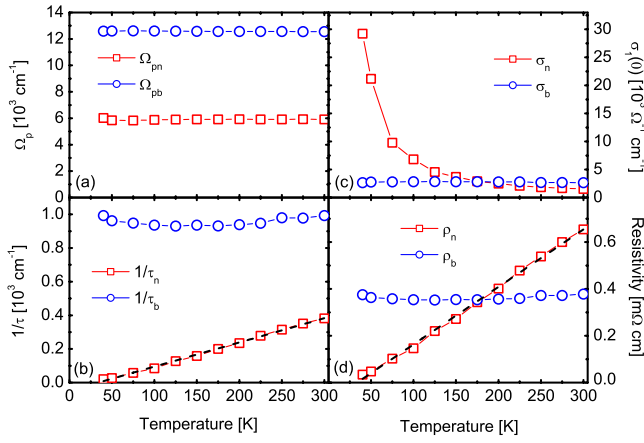


FIG. 3 (color online). The  $T$  dependence of (a) the plasma frequency  $\Omega_p$ , (b) the scattering rate  $1/\tau$ , (c) the contribution to dc conductivity  $\sigma_1(\omega \rightarrow 0)$ , and (d) the equivalent resistivity  $\rho$  for the narrow and broad Drude components, respectively. The dashed lines in (b) and (d) are linear fits to  $1/\tau_n$  and  $\rho_n$ , respectively.

Upon cooling,  $\Omega_p$  is roughly a constant for each of the two components, indicating that the band structure and  $n/m^*$  do not change with temperature, in agreement with a previous work [35]. Figure 3(b) portrays the  $T$  dependence of the scattering rate of the two Drude components, where  $1/\tau_b$  is basically  $T$  independent while  $1/\tau_n \propto T$ ; the black dashed line denotes a linear fit. Figure 3(c) displays the contribution of the two groups of carriers to the dc conductivity. As the broad Drude ( $\sigma_b$ ) is  $T$  independent, the temperature dependence of the total dc conductivity arises out of the narrow Drude band ( $\sigma_n$ ). The  $\rho(T)$  curve of  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$ , shown in the inset of Fig. 1, exhibits a tendency to saturation at room temperature, and a change of slope can be seen at about 175 K. This behavior can be explained by the different  $T$  dependence of the two Drude bands, which can be considered as a parallel circuit [40]:  $\sigma = \sigma_n + \sigma_b$ . A crossover region, where  $\sigma_n \approx \sigma_b$ , is found in Fig. 3(c) at  $\approx 175$  K. Below this temperature,  $\sigma_n > \sigma_b$ , so the total dc conductivity is dominated by  $\sigma_n$ , which exhibits strong temperature dependence. As a result, below 175 K,  $\rho(T)$  decreases quickly with decreasing temperature. Above 175 K,  $\sigma_n < \sigma_b$ , and the total dc conductivity is dominated by  $\sigma_b$ , which shows no temperature dependence. Hence, above 175 K, the growth of the dc resistivity slows with heating, resulting in the change of slope and the tendency to saturation in  $\rho(T)$ . Similar conclusions were obtained from investigations of the Hall effect [28,29] and theoretical calculations [41]. In the  $\omega \rightarrow 0$  limit, the inverse of  $\sigma_1(0)$  yields the resistivity from the two Drude components, as shown in Fig. 3(d). The resistivity of the broad Drude remains a constant at all measured temperatures, while a  $T$ -linear  $\rho_n$  is revealed for the narrow Drude component. This is in accord with transport measurements on hole-doped 122 compounds [18–20], where  $T$ -linear  $\rho$  was observed at

low temperatures in optimally doped materials. The  $T$ -linear  $\rho$  is found only at low temperatures as this is the region dominated by the narrow Drude component.

Further evidence for  $T$ -linear  $1/\tau$  lies in the temperature dependence of the low-frequency spectral weight. The spectral weight is defined as

$$W_0^{\omega_c} = \int_0^{\omega_c} \sigma_1(\omega) d\omega, \quad (2)$$

where  $\omega_c$  is a cutoff frequency. In a Drude metal the scattering rate decreases upon cooling, producing a narrowing of the Drude response and an increase of the dc conductivity, resulting in a transfer of spectral weight from high to low frequencies and an increase in  $W_0^{\omega_c}$ . To quantitatively analyze the  $T$  dependence of the low-frequency spectral weight, we adopt one Drude optical conductivity [see Eq. (1)] into Eq. (2), to obtain the spectral weight as a function of  $1/\tau$ ,

$$W_0^{\omega_c}(1/\tau) = \frac{2\pi\Omega_p^2}{Z_0} \arctan(\omega_c\tau). \quad (3)$$

In the case of  $1/\tau \propto T$ , Eq. (3) can be simplified as

$$W_0^{\omega_c}(T) = a_1 \arctan\left(\frac{a_2}{T}\right), \quad (4)$$

where  $a_1 = 2\pi\Omega_p^2/Z_0$ , and  $a_2 \propto \omega_c$ ; both are  $T$ -independent parameters. Considering the spectral weight arising from the incoherent part and interband transition (Lorentz), which are both  $T$  independent, we introduce the third  $T$ -independent parameter  $a_0$  into Eq. (4). Finally, the low-frequency spectral weight as a function of temperature for  $1/\tau \propto T$  is written as

$$W_0^{\omega_c}(T) = a_0 + a_1 \arctan\left(\frac{a_2}{T}\right). \quad (5)$$

$W_0^{\omega_c}$  can easily be determined by integrating the measured  $\sigma_1(\omega)$ . The open symbols in Fig. 4 denote  $W_0^{\omega_c}$  with different  $\omega_c$ 's: 150 (squares), 200 (triangles), and 250  $\text{cm}^{-1}$  (diamonds). As expected in a metallic system, for all the  $\omega_c$ 's,  $W_0^{\omega_c}$  increases with decreasing  $T$ . The solid curves through the data are the least-squares fits using Eq. (5). The excellent agreement indicates that a  $T$ -linear  $1/\tau$  dominates the low-frequency  $\sigma_1(\omega)$  over a very broad temperature range.

We now discuss the possible origin of the  $T$ -linear  $1/\tau_n$  and  $\rho_n$ . A  $T$ -linear  $\rho$  is usually expected in a system dominated by electron-phonon scattering in the regime  $T \gg \Theta_D$ , where  $\Theta_D$  is the Debye temperature; in the regime  $T < \Theta_D$ , the resistivity is approximated by a low-order polynomial that is neither linear nor quadratic in temperature. Transport measurements in the 122 compounds suggest  $\Theta_D \approx 250$  K [37,42]. However, the  $T$ -linear  $1/\tau_n$  and  $\rho_n$  span a very broad temperature range, from  $T_c$  up to 300 K, which is inconsistent with the electron-phonon dominated scattering process. Further clues on the  $T$ -linear  $1/\tau_n$  and  $\rho_n$  may be revealed by an examination of the phase diagram of the  $\text{Ba}_{1-x}\text{K}_x\text{Fe}_2\text{As}_2$  system and a comparison with similar behavior found in other materials.

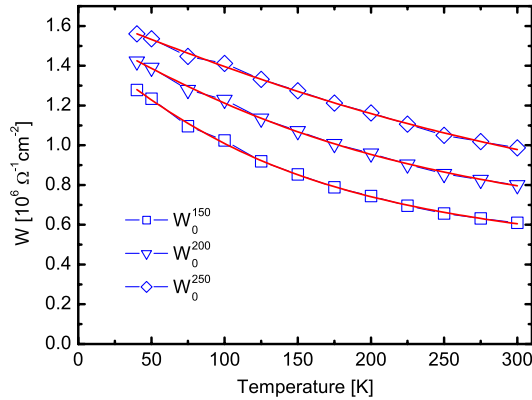


FIG. 4 (color online). Temperature dependence of the low-frequency spectral weight  $W_0^{\omega_c}$  with different  $\omega_c$ 's: 150 (squares), 200 (triangles), and 250  $\text{cm}^{-1}$  (diamonds). The solid lines through the data represent the expected temperature dependence for  $1/\tau \propto T$  at each  $\omega_c$ .

The parent compound  $\text{BaFe}_2\text{As}_2$  orders in a SDW state below  $T_{\text{SDW}} \approx 138$  K;  $T_{\text{SDW}}$  is suppressed by K doping allowing superconductivity to emerge. Here,  $1/\tau_n \propto T$  is observed at the doping where the SDW order is entirely suppressed, i.e.,  $T_{\text{SDW}} \rightarrow 0$ . At this point, spin fluctuations are expected to be very strong, which has been experimentally confirmed by NMR [43–45]. This brings us to models that attribute  $T$ -linear  $\rho$  to spin-fluctuation scattering [13,46]. The  $T$ -linear  $\rho$  (or  $1/\tau$ ) is also found in cuprates such as Nd-doped  $\text{La}_{2-x}\text{Sr}_x\text{CuO}_4$  [3] and electron-doped  $\text{La}_{2-x}\text{Ce}_x\text{CuO}_4$  [4], organic superconductors  $(\text{TMTSF})_2\text{PF}_6$  [5] and  $(\text{TMTSF})_2\text{ClO}_4$  [6], as well as a number of heavy fermions such as  $\text{CeCoIn}_5$  [7,8] and  $\text{YbRh}_2\text{Si}_2$  [9,10]. Studies on these materials have shown that  $T$ -linear  $\rho$ , arising on the border of antiferromagnetic order, is caused by spin-fluctuation scattering due to the proximity of an antiferromagnetic QCP. Studies on cuprates and Bechgaard salts [4,11] further show that the strength (or coefficient) of the  $T$ -linear  $\rho$  scales with  $T_c$  and disappears upon approaching the point where  $T_c \rightarrow 0$ , suggestive of an intimate relation between the  $T$ -linear  $\rho$  and superconductivity. Anomalous  $T$ -linear  $\rho$  (or  $1/\tau$ ) and pairing in unconventional superconductors may share a common origin: spin fluctuations. The hidden  $T$ -linear  $1/\tau_n$  and  $\rho_n$  in  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$  revealed by optical measurements may have the same origin as those found in the cuprates, organic superconductors, and heavy-fermion metals, because these materials share strikingly similar phase diagrams. Our observations may also imply a possible QCP in the superconducting dome. The existence of the QCP in iron pnictides is supported by transport properties [18–20], NMR studies [43,44], de Haas–van Alphen effect [47], penetration depth measurement [21], and first-principles calculations [17].

Spin-fluctuation induced  $T$ -linear  $\rho$  suggests an equivalent  $\omega$ -linear  $1/\tau(\omega)$  [13], obtainable through the extended Drude model provided that the interband contribution is negligible. Low energy interband transitions are important in iron pnictides, and their contribution has to be subtracted to

determine  $1/\tau(\omega)$  for mobile carriers [32,48]. We calculated  $1/\tau(\omega)$  via the extended Drude model with the interband contribution subtracted (see Supplemental Material [49]) and found that, by taking into account the interband transitions, a large fraction of the frequency dependence in  $1/\tau(\omega)$  is eliminated, which is consistent with the analysis of Charnukha *et al.* [32] on  $\text{Ba}_{0.68}\text{K}_{0.32}\text{Fe}_2\text{As}_2$  [32]. There is no confident evidence for the expected  $\omega$ -linear  $1/\tau(\omega)$ , since it could be masked by the multiband character of the iron pnictides.

In order to check if the  $T$ -linear  $1/\tau$  is unique in  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$  or general in iron pnictides, we applied the same analysis to  $\text{BaFe}_2(\text{As}_{0.7}\text{P}_{0.3})_2$  (see Supplemental Material [49]). Interestingly,  $T$ -linear  $1/\tau$  is also found for the narrow Drude. This suggests that  $T$ -linear  $1/\tau$  is not unique in  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$ , but most likely a general behavior in iron pnictides at the doping where SDW order is completely suppressed, i.e.,  $T_{\text{SDW}} \rightarrow 0$ .

In summary, the detailed  $T$  dependence of the normal state  $\sigma_1(\omega)$  and the low-frequency spectral weight in  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$  have been examined. Two Drude components with different  $1/\tau$ 's yield an excellent description of the low-frequency optical response, indicating the existence of two groups of carriers with different quasiparticle lifetimes. The broad Drude component produces an incoherent background conductivity with no temperature dependence, while the narrow Drude component reveals a  $T$ -linear  $1/\tau_n$  and  $\rho_n$ . This fact explains the  $T$ -linear  $\rho$  behavior at low temperatures and the tendency to saturation at room temperature observed in transport measurements in optimally hole-doped 122 compounds. An arctan( $T$ ) dependence of the low-frequency spectral weight is also strong evidence for a  $T$ -linear  $1/\tau$ . A comparison with similar behavior found in other materials suggests that the  $T$ -linear  $1/\tau_n$  and  $\rho_n$  in  $\text{Ba}_{0.6}\text{K}_{0.4}\text{Fe}_2\text{As}_2$  may arise out of spin-fluctuation scattering due to the possible existence of an antiferromagnetic QCP in the superconducting dome.

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