Comparison of the Fermi-surface topologies of $\kappa$-(BEDT-TTF)$_2$Cu(NCS)$_2$ and its deuterated analogue

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We have measured details of the quasi one-dimensional Fermi-surface sections in the organic superconductor $\kappa$-(BEDT-TTF)$_2$Cu(NCS)$_2$ and its deuterated analogue using angle-dependent millimetre-wave techniques. There are significant differences in the corrugations of the Fermi surfaces in the deuterated and undeuterated salts. We suggest that this is important in understanding the inverse isotope effect, where the superconducting transition temperature rises on deuteration. The data support models for superconductivity which invoke electron-electron interactions depending on the topological properties of the Fermi surface.

PACS numbers: 71.18.+y, 71.27.+a, 72.80.Le, 74.70.-b, 78.70.Gq

The nature of the superconducting groundstate in quasi-two-dimensional charge-transfer salts such as $\kappa$-(BEDT-TTF)$_2$Cu(NCS)$_2$ has attracted much recent experimental and theoretical interest. The majority of the experimental data suggest that the superconductivity is not describable by a simple BCS-like, phonon-mediated approach (for a review, see Refs. [3,11] and refs. therein). Consequently, a number of the theoretical models invoke pairing mediated by electron-electron interactions and/or antiferromagnetic fluctuations [4,12]. In such a scenario, the "nestability" of the Fermi surface is an important consideration; it is expected that alterations of the Fermi-surface topology will affect the superconducting transition temperature.

In this context, the observation of a "negative isotope effect" in $\kappa$-(BEDT-TTF)$_2$Cu(NCS)$_2$ may be of great importance. On replacing the terminal hydorgen of the BEDT-TTF molecule in $\kappa$-(BEDT-TTF)$_2$Cu(NCS)$_2$ by deuterium, it was found that an increase ($\Delta T_c \approx 0.3 \text{ K}$) in the superconducting critical temperature $T_c$ occurred [13]. By contrast, isotopic substitutions of other atoms in the BEDT-TTF molecule or in the anion layer produce a very small, normal isotope effect or no significant isotope effect at all, respectively [14].

In this paper we describe millimetre-wave measurements which compare the Fermi surfaces of deuterated and conventional samples of $\kappa$-(BEDT-TTF)$_2$Cu(NCS)$_2$. The data suggest that it is primarily the changes in the topology of the Fermi surface brought about by deuteration that cause the observed isotope effect, supporting models for superconductivity involving pairing via electron-electron interactions [15].

The Fermi surface of $\kappa$-(BEDT-TTF)$_2$Cu(NCS)$_2$ comprises a quasi-two-dimensional (Q2D) pocket (the $\alpha$ pocket) and a pair of quasi-one-dimensional (Q1D) sheets [13]; it is similar to that used by Pippard to predict magnetic breakdown [15]. It is known that the $\alpha$ pocket is hardly affected by deuteration [16], and so our measurement concentrates on the Q1D sheets. Fermi-surface traversal resonances (FTRs) [15] (i.e. resonances in the high-frequency conductivity caused by magnetic-field-induced motion of quasiparticles across the Fermi sheets) are used to infer the corrugations of the sheets.

The experiments involved single crystals of $\kappa$-(BEDT-TTF)$_2$Cu(NCS)$_2$ ($\sim 0.7 \times 0.5 \times 0.1 \text{ mm}^3$; mosaic spread $\approx 0.1^\circ$), produced using electrocrystallization [11,12]. In some of the crystals, the terminal hydorgen of the BEDT-TTF molecules were isotopically substituted by deuterium; we refer to the deuterated samples as d8, and conventional hydrogenated samples as h8. A single sample is mounted at the centre (in a magnetic field antinode) of a rectangular cavity of inner dimensions $1.55 \times 3.10 \times 6.00 \text{ mm}^3$ resonating at 72 GHz in the $TE_{102}$ mode [17]: the oscillating H-field lies within the sample's Q2D ($b$, $c$) planes. In this configuration, the effective skin depth is very large, and the GHZ fields penetrate the bulk of the sample [18]. The cavity can be rotated with respect to the external quasistatic magnetic field $B$ so as to vary the angle $\theta$ between $B$ and the normal to the sample's Q2D planes [19]; the normal to the Q2D planes is the $a^*$ direction of the reciprocal lattice [17]. In addition, the sample can be turned about $a^*$ within the cavity, so as to vary the plane of rotation, defined by the azimuthal angle $\phi$ [20]. The angles $\theta$ and $\phi$ and their relationship to the Q1D sheet of the Fermi surface are given in the inset to Fig. 1.

Experiments were carried out on samples of d8 $\kappa$-(BEDT-TTF)$_2$Cu(NCS)$_2$ for angles $-70^\circ \leq \theta \leq 70^\circ$ for four different azimuthal angles $\phi$. Fig. 1 shows results for an azimuthal angle of $\phi = 5^\circ$ for $\theta$ values between $0^\circ$ and $70^\circ$ in $5^\circ$ steps, at a temperature of 1.5 K. At low fields (around 4 T at $\theta = 0^\circ$) one absorption can be seen (thick dashed line in Fig. 1). This is related to the superconducting to normal transition of the sample [21]; it follows the $\theta$ dependence of the upper critical field, $H_{c2}$, which varies approximately as $1/\cos \theta$ [22]. At high field magnetic quantum oscillations are observed, indicating that the sample is pure; the angular behaviour of the frequency $F$ of the oscillations ($F \propto 1/\cos \theta$) pro-
Each corrugation of the Q1D Fermi sheets is expected to give rise to a FTR with the ψ dependence \[ \frac{\omega}{B_{||}} = A \sin(\psi - \psi_0). \] (2)

Here, ω is the angular frequency of the millimetre-waves, A is a constant depending on details of the Fermi surface [15], and ψ₀ defines the axis of the corrugation R (Fig. 1, inset). As the millimetre-wave frequency is held constant, the FTRs should lie on sinusoidal “arches” when 1/B₀ is plotted as a function of ψ [14].

Fig. 2 shows the FTR positions plotted in terms of 1/B₀ and ψ. Apart from a region close to ψ = +100° where the feature associated with the superconducting-to-normal transition obscures the FTRs at some φ, making the exact position difficult to gauge, the data lie on two “arches”, shown as curves (Fig. 2); the curves were obtained by fitting the data to Eqn. 2. This indicates that the Q1D Fermi surface of d8 κ-(BEDT-TTF)₂Cu(NCS)₂ has two distinct corrugations, with their axes R₁ and R₂ at angles ψ₀ = 17.9° ± 2.0° and 39.8° ± 2.0° to a* respectively.

Equivalent experiments were carried out in h8 κ-(BEDT-TTF)₂Cu(NCS)₂. (Some representative data are shown in Ref. [15].) Again, the resonance positions lie on two “arches” shown in Fig. 2, implying that the Q1D Fermi surface of h8 κ-(BEDT-TTF)₂Cu(NCS)₂ has two distinct corrugations [15]. In this case, the corrugation axes R₁ and R₂ are at angles ψ₀ = 21.2° ± 2.0° and −20.8° ± 2.0° to a* respectively. The data for both samples are summarised in Table I.
\(\kappa-(\text{BEDT-TTF})_2\text{Cu(NCS)}_2\) has a monoclinic crystal structure with the crystallographic \(a\)-axis at an angle of 20.3° to the normal to the Q2D planes \((a^*)\) [17]; in this respect, the crystal structures of \(d8\) and \(h8\) \(\kappa-(\text{BEDT-TTF})_2\text{Cu(NCS)}_2\) appear identical [18]. Within tight-binding bandstructure, the corrugation axes of a Fermi surface usually relate to the primitive lattice translation vectors of the real-space lattice [24]. Rather than work in terms of the angle \(\psi_0\), which defines the directions of the corrugation axes with respect to \(a\), it is more useful to use the angle \(\Psi_0\), which relates to the real-space vector \(a\). Once this is done (Table I), it is plain that the corrugation axis \(R_1\) in both the \(d8\) and \(h8\) samples lies very close to the \(a\) (interlayer) direction. By contrast, the direction of \(R_2\), the second corrugation axis, differs; with reference to the primitive lattice translation vectors

\[
T_{mn} = ma + nc
\]  

where \(m\) and \(n\) are integers, we find that in \(d8\), \(R_2\) is very close in direction to \(T_{2-1}\), whereas in \(h8\), it is close in direction to \(T_{11}\) (Table I). The reasons for the dominance of these particular directions are unclear; however, interlayer coupling through the anion layer is presently poorly understood at a molecular-orbital level. It is possible that a variety of overlap-pathways may be operative and that the choice of dominant overlap-pathway through the anion layer depends very sensitively on the exact coordinates of the terminal end of the BEDT-TTF molecule. In this context, it will be very useful to have high-resolution structural experiments which address the detailed differences between \(h8\) and \(d8\) at low temperatures [18].

Finally, it is interesting to work out the relative amplitudes of the corrugations in \(h8\) and \(d8\) \(\kappa-(\text{BEDT-TTF})_2\text{Cu(NCS)}_2\). Models of FTR allow one to relate the intensity of the FTR to the amplitude of the Fermi-surface corrugation [21]. Measurements of the d.c. transport properties of \(d8\) and \(h8\) \(\kappa-(\text{BEDT-TTF})_2\text{Cu(NCS)}_2\) suggest that the transfer integral \(t_1\) in the \(a\) direction is very similar in the two materials \((t_1 \approx 0.04 \text{ meV})\) [13]. This implies that the Fermi-surface corrugations along \(R_1\) should be very similar in \(d8\) and \(h8\) [13]; in \(d8\) and \(h8\) samples with equal volume the corresponding FTRs should have the same intensity [21]. The average intensities \(I\) of each FTR for \(\theta = 0^\circ\) are shown in Table I. As the samples are of different size, the intensities of the FTRs have been normalised to that of the FTR corresponding to \(R_1\).

Using the periodicity in \(k\)-space [17], the relative intensities of the FTRs, and the orientations \(\Psi_0\) of \(R_1\) and \(R_2\), it is possible to make a comparison of the Q1D Fermi sheets for both materials. Fig. 3 shows these representations, with the corrugations, assumed sinusoidal, shown at the same scale. This scale is chosen so that the differences between \(d8\) and \(h8\) are clear; in reality, the small measured value of \(t_1\) [13] suggests that the corrugations will be on an extremely small scale.

Despite an intensive search, no features attributable to cyclotron resonance (CR) due to the Q2D Fermi-surface pocket were observed in either \(d8\) or \(h8\), in agreement with previous studies [13]. The response of our cavity system is dominated by the interlayer component of the sample's high-frequency conductivity [15, 16]. Fermi-surface sections with more complex corrugations in the interlayer direction will dominate the high-frequency interlayer conductivity [21]; recent resistivity measurements of \(\kappa-(\text{BEDT-TTF})_2\text{Cu(NCS)}_2\) suggest that the corrugations of the Q2D pocket are simpler and more regular than those of the Q1D sheets [13], perhaps explaining the absence of CR. Similarly, the in-plane corrugations of the Q1D sheet (Fig. 3(c)) will have little effect on the high-frequency interlayer conductivity [21], and therefore do not result in detectable FTRs.

It is obvious that there is a difference between the Q1D Fermi sheets of the two materials, with the corrugations
which define the directions of the corrugation axes \( \kappa \) with respect to \( \Psi \). Hence they are mediated by electron-electron interactions which depend on the "nestability" of the Fermi-surface; hence they are more amenable to nesting than that of \( h8 \).

In summary, we have measured details of the Fermi-surface topology of the deuterated organic superconductor \( \kappa-(BEDT-TTF)_2Cu(NCS)_2 \), and compared them with equivalent measurements of the undeuterated salt. We find that the quasi-one-dimensional Fermi-surface sheets are significantly more corrugated in the undeuterated salt, perhaps explaining the "inverse isotope effect" observed on deuteration. Our data support models for exotic d-wave superconductivity in the organics which invoke electron-electron interactions depending on the topological properties of the Fermi surface.

This work is supported by EPSRC (UK). NHMFL is supported by the US Department of Energy (DoE), the National Science Foundation and the State of Florida. Work at Argonne is sponsored by the DoE, Office of Basic Energy Sciences, Division of Materials Science under contract number W-31-109-ENG-38. We thank Stephen Hill and Stephen Blundell for constructive comments.

\[ \begin{array}{cccc}
A/\omega & d8 R_1 & d8 R_2 & h8 R_1 & h8 R_2 \\
0.198 \pm 0.004 & 0.131 \pm 0.002 & 0.204 \pm 0.004 & 0.168 \pm 0.004 \\
\psi_0 & 17.9 \pm 2.0^\circ & 39.8 \pm 2.0^\circ & 21.2 \pm 2.0^\circ & -20.8 \pm 2.0^\circ \\
\Psi_0 & -2.4 \pm 2.0^\circ & 19.5 \pm 2.0^\circ & 0.9 \pm 2.0^\circ & -41.1 \pm 2.0^\circ \\
T_{mn} & T_{10} & T_{2-1} & T_{10} & T_{11} \\
T & 1 & 0.34 \pm 0.06 & 1 & 6 \pm 2 \\
\end{array} \]

TABLE I: The values for \( A/\omega \) (see Eqn. 3), \( \psi_0 \) (angle of corrugation axis with respect to \( a^* \)) and \( \Psi_0 \) (angle of corrugation axis with respect to \( a \)) for each of the FTRs seen in \( d8 \) and \( h8 \) \( \kappa-(BEDT-TTF)_2Cu(NCS)_2 \). Also shown are the vectors \( T_{mn} \) which define the directions of the corrugation axes \( R_1 \) and \( R_2 \) (see Eqn. 3). \( I \) is the average intensity of the FTR at \( \theta = 0^\circ \) normalised as described in the text.

in \( h8 \) being stronger; the dominant corrugation has axis \( R_2 \), at -41.1° to \( a \). By contrast, the corrugations in \( d8 \) \( \kappa-(BEDT-TTF)_2Cu(NCS)_2 \) are weaker, and are dominated by that with axis \( R_1 \) lying along \( a \). This suggests that the Fermi surface of \( d8 \) \( \kappa-(BEDT-TTF)_2Cu(NCS)_2 \) would be more amenable to nesting than that of \( h8 \).

Our data support models for superconductivity such as those of Refs. [12, 13]. In these, the pairing of electrons is mediated by electron-electron interactions which depend on the "nestability" of the Fermi-surface; hence they predict a \( T_c \) which is sensitive to the details of the Fermi-surface topology. The difference between the Q1D Fermi sheets of \( d8 \) and \( h8 \) \( \kappa-(BEDT-TTF)_2Cu(NCS)_2 \) measured using FTR can thus explain the isotope effect; the Q1D Fermi sheets in the \( d8 \) samples are less corrugated (and therefore more nestable), leading to a higher \( T_c \).

In summary, we have measured details of the Fermi-surface topology of the deuterated organic superconductor \( \kappa-(BEDT-TTF)_2Cu(NCS)_2 \), and compared them with equivalent measurements of the undeuterated salt. We find that the quasi-one-dimensional Fermi-surface sheets are significantly more corrugated in the undeuterated salt, perhaps explaining the "inverse isotope effect" observed on deuteration. Our data support models for exotic d-wave superconductivity in the organics which invoke electron-electron interactions depending on the topological properties of the Fermi surface.

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