Unconventional superconductivity in Cu\textsubscript{x}Bi\textsubscript{2}Se\textsubscript{3} from magnetic susceptibility and electrical transport

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Abstract

Although the Cu doped Bi\textsubscript{2}Se\textsubscript{3} topological insulator was discovered and intensively studied for almost a decade, its electrical and magnetic properties in normal state, and the mechanism of 'high-'\textsubscript{T}\textsubscript{c} superconductivity regarding the relatively low-carrier density are still not addressed yet. In this work, we report a systematic investigation of magnetic susceptibility, critical fields, and electrical transport on the nominal Cu\textsubscript{0.09}Bi\textsubscript{2}Se\textsubscript{3} single crystals with \textit{T}\textsubscript{onset} = 4.18 K, the highest so far. The composition analysis yields the Cu stoichiometry of \(x = 0.09\). The magnetic susceptibility shows considerable anisotropy and an obvious kink at around 96 K was observed in the magnetic susceptibility for \(H \parallel c\), which indicates a charge density anomaly. The electrical transport measurements indicate the two-dimensional (2D) Fermi liquid behavior at low temperatures with a high Kadowaki–Woods ratio, \(\lambda/\gamma^2 = 30.3a_0\). The lower critical field at 0 K limit was extracted to be 6.0 Oe for \(H \parallel ab\). In the clean limit, the ratio of energy gap to \(T_c\) was determined to be \(\Delta_0/k_B T_c = 2.029 \pm 0.124\), exceeding the standard BCS value 1.764, suggesting Cu\textsubscript{0.09}Bi\textsubscript{2}Se\textsubscript{3} is a strong-coupling superconductor. The in-plane penetration depth at 0 K was calculated to be 1541.57 nm, resulting in an unprecedented high ratio of \(\lambda/\lambda_0(0) = 9.86\). Moreover, the ratio of \(T_c\) to Fermi temperature is estimated to be \(T_c/T_F^{2D} = 0.034\). Both ratios fall into the region of unconventional superconductivity according to Uemura’s regime, supporting the unconventional superconducting mechanism in Cu\textsubscript{x}Bi\textsubscript{2}Se\textsubscript{3}. Finally, the enhanced \(T_c\) value higher than 4 K is proposed to arise from the increased density of states at Fermi energy and strong electron–phonon interaction induced by the charge density instability.

1. Introduction

The superconducting critical temperature (\(T_c\)) up to 3.8 K in doped superconducting topological insulators such as Cu\textsubscript{x}Bi\textsubscript{2}Se\textsubscript{3} are unexpectedly 'high' for a low carrier density semiconductor [1]. According to the quantum oscillations results [2], the Fermi wave number \(k_F\) in Cu\textsubscript{x}Bi\textsubscript{2}Se\textsubscript{3} is increased from 0.69 nm\textsuperscript{-1} to 0.97 nm\textsuperscript{-1} by Cu doping. Such value lies within the range of 0.1–1 nm\textsuperscript{-1} for typical densities in a classical two-band single valley semiconductor [3], in which the density of states (DOS) is only about one-tenth of that of a metal. Actually, the superconductivity was also observed in some semiconductors with \(T_c\) far below 1 K, such as SrTiO\textsubscript{3} (\(T_c \sim 0.05-0.29\) K, \(n_e = 0.65\) to 42.4 \texttimes 10\textsuperscript{20} cm\textsuperscript{-3}) [3, 4], GeTe (\(T_c \sim 0.07-0.3\) K, \(n_e = 8.6\) to 15.2 \texttimes 10\textsuperscript{20} cm\textsuperscript{-3}) [5], and SnTe (\(T_c \sim 0.034-0.214\) K, \(n_p = 7.5\) to 20.0 \texttimes 10\textsuperscript{20} cm\textsuperscript{-3}) [6]. To compare with these systems, the \(T_c\) value of Cu\textsubscript{x}Bi\textsubscript{2}Se\textsubscript{3} is relatively 'high' given it is a highly doped narrow semiconductor with lower carrier density. However, the mechanism of such anomalous enhanced \(T_c\) phenomenon remains unclear despite nearly a decade of extensive research.
The emergence and enhancement of superconductivity may occur in various cases by tailoring the ground states of quantum materials. For example, suppression of charge density wave (CDW) by chemical doping or pressure is observed in many systems [7–10], including low dimensional layered Cu4TiSe2 and three dimensional (3D) alloys Lu(Pr1−xPdx)3In and (Sr,Ca)3Ir5Sn13, 3D oxides Ba1−xKxBiO3. Similarly, the magnetic ordering or spin density wave (SDW) have also been suppressed in strongly correlated heavy fermions materials, high-Tc cuprates and iron-based superconductors [11–13]. Lastly, the value of Tc can be significantly suppressed by introducing disorders [14], leading to metal–insulator transition. Note that such mechanism may be also relevant to superconducting semiconductors, since a large number of disorders are expected inside.

In this work, we revisit the superconducting Cu0.20Bi2Se3 system, which is widely studied in the past but rarely emphasizing on the magnetic susceptibility and electrical transport properties in the normal state. Interestingly, we observed Tc = 4.18 K in Cu0.09Bi2Se3 for our sample, the highest value to date. In addition, for the first time, we reveal a charge density anomaly at around 96 K in the Cu0.09Bi2Se3 crystal from the magnetic susceptibility in the normal state, which is reminiscent of unconventional superconductivity proximity to CDW. We show the Cu0.09Bi2Se3 displays the 2D Fermi liquid behavior at low temperature with a high Kadowaki–Woods ratio, A/γ2 = 30.3a0, as may be enhanced by proximity to charge density instability and/or electronic anisotropy. Furthermore, we obtain the upper and lower critical fields from the measurement of temperature dependent magnetic susceptibility under different fields. We find the Cu0.09Bi2Se3 is a strong-coupling superconductor with ∆0/kBTc = 2.029 ± 0.124 exceeding the standard BCS value 1.764. Finally, we discuss the possible underlying mechanisms for the enhanced Tc higher than 4 K and the unconventional superconducting features. Importantly, we demonstrate the Cu0.09Bi2Se3 belongs to one of unconventional superconducting families according to Uemura’s plot, with unprecedented high ratios of Tc/λ2(0) ≈ 9.86 and Tc/Δ0D = 0.034 among superconducting semiconductors.

2. Experimental details

The samples were prepared using high purity (99.999%) elements in the nominal composition Cu0.20Bi2Se3 (x = 0.20), as reported elsewhere [15, 16]. We find some crystals cleaved from the large batch are superconducting at 4.18 K (denoted as CBS418), which is higher than previous report [1]. The phase purity of obtained single crystals was examined by x-ray diffraction (XRD) measurements with Cu Kα radiation. To check the homogeneity, we performed the measurements of the energy-dispersive spectrum on the sample with Tc,onset = 4.18 K. In addition, we also present the measurements of pure Bi2Se3 (denoted as BS) and nominal Cu0.20Bi2Se3 with Tc,onset = 3.79 K (denoted as CBS379). The in-plane resistivity measurements were carried out on a physical property measurement system (PPMS-9, quantum design) using the standard four probes with silver paste as the contacts. The DC magnetic susceptibility and initial magnetizations curves were measured with a SQUID-VSM magnetometer (quantum design) at different magnetic fields and temperatures.

3. Results and discussion

3.1. Crystal structure and composition

Figures 1(a) and (b) show the crystal structure and XRD patterns as collected on Cu0.20Bi2Se3 single crystal. The parent Bi2Se3 compound crystallizes to the rhombohedral structure with space group R3m (no. 166) [17]. As suggested by the previous report [1], the intercalated Cu atoms were in the van der Waals gap between Bi2Se3 layers, which partially occupy the octahedrally coordinated 3b (0, 0, 1/2) sites, as indicated in figure 1(a). Only a series of (00l) reflections were observed and the background is clean, indicating the phase purity. This is consistent with those observed from the rhombohedral space group of Bi2Se3 [17].

To investigate the sample homogeneity, we performed composition analysis on the sample with the onset superconducting transition temperature Tc,onset = 4.18 K by energy dispersive x-ray spectroscopy (EDS). A total of 30 random EDS spectra locating at different regions were measured, as indicated by yellow frames in the inset of figure 2(a). The atomic ratios were plotted in figure 2(b). We observed that there are three measured regions with higher Cu content than others, where the Cu content is between 0.15 and 0.25. In these regions, all the element ratios of Bi and Se are very close to their stoichiometric composition, indicating the Bi and Se elements are generally homogeneous on a large scale. The remaining measured EDS spectrum yields an averaged composition Cu0.09(1)Bi0.02(2)Se2.98(2) (referred as Cu0.09Bi2Se3), in which the Cu content ratio still fluctuates within 23% compared to the averaged value. This demonstrates the melt-grown crystal is quite inhomogeneous on a micrometer scale. It is reported that the distribution of Cu
content in the electrochemical synthesized samples shows a ∼19% variation [18], which is slightly smaller than that in our melt-grown crystals. According to phase diagram reported by Kriener et al [18], the higher $T_c$ phase is deemed to have a lower Cu doping level.

3.2. Magnetic susceptibility

Figure 3 shows the temperature dependence of volume susceptibility $\chi(T)$ measured under $H = 3$ Oe on one crystal with dimensions of $\sim 5.83 \times 4.67 \times 0.59$ mm$^3$. We observe $T_c^{\text{onset}}$ occurs at 4.18 K, which is the highest in doped Bi$_2$Se$_3$ to our knowledge. The apparent difference between zero-field cooling (ZFC) and field cooling (FC) is caused by the vortex pinning, which has been investigated in detail for Cu$_x$Bi$_2$Se$_3$ [15, 16]. Interestingly, the peak effect, a feature of anomalous vortex pinning phenomenon, is observed in the high quality clean Cu$_x$Bi$_2$Se$_3$ crystals [15, 16]. To estimate the volume shielding fraction, we take the demagnetization factor $N$ into account, which can be calculated by [19]

$$N = \left[ 1 + \frac{3}{4} \left( 1 + \frac{a}{b} \right) \right]^{-1}, \quad (1)$$

where $a$, $b$, $c$ are dimensions ($2a \times 2b \times 2c$) of the experimental specimen with rectangular cuboid shape. According to equation (1), we obtain $N_{ab} = 0.107$ for $H \parallel ab$ and $N_c = 0.854$ for $H \parallel c$. The intrinsic
volume susceptibility $\chi_{\text{int}}$ is given by

$$\chi_{\text{int}} = \chi_{\text{exp}}/(1 - N\chi_{\text{exp}}),$$

where $\chi_{\text{exp}}$ is the experimental measured susceptibility. As seen in figure 3, a 17% superconducting shielding fraction was estimated, which is consistent with other reports [1, 20–22].

To investigate the magnetic properties in normal state, we measured $\chi(T)$ under 0.3 T, as is plotted in figure 4. The magnetic susceptibility $\chi \equiv M/H$ for both $H \parallel c$ and $H \parallel ab$ shows a small negative background, essentially presenting a weak temperature dependence at high temperatures ($T > 96$ K). This indicates that the Larmor or Langevin diamagnetic and Landau diamagnetic contributions are larger than the sum of the spin Pauli paramagnetic susceptibility. With decreasing the temperature, the magnitude of $\chi$ firstly shows a quasilinear increase up to 96 K, below which a kink occurs. The drop of $\chi$ across the kink

Figure 3. Volume susceptibility versus temperature $\chi(T)$ under 3 Oe for $H \parallel c$ in ZFC and FC modes. Inset shows the detailed region at the onset superconducting transition.

Figure 4. Temperature dependence of magnetic susceptibility under 0.3 T for (a) $H \parallel c$, (b) $H \parallel ab$. The insets is the Curie–Weiss fit.
temperature \((T_k)\) implies a decrease of electronic DOS at Fermi energy \((E_F)\). The quasilinear behavior of \(\chi\) above the \(T_k\) and its weak temperature dependence have been widely observed in some CDW materials like \(\text{Cu}_x\text{TiSe}_2\) \([7]\), \(\text{Lu}(\text{Pt}_{1-x}\text{Pd}_x)\text{In}\) \([8]\), and \((\text{Sr}, \text{Ca})_3\text{Ir}_4\text{Sn}_{13}\) \([9]\). In \(\text{Cu}_{0.09}\text{Bi}_2\text{Se}_3\), the superconductivity is induced only by Cu intercalation, which is in favor of \(\text{Cu}^{+}\)–non-magnetic state (as is in the case of interstitial site, see reference \([23]\)). While the Cu has a possibility to substitute the Bi \([23]\), which can create two holes and act as carrier acceptors to depress the carrier density. Therefore, almost all the ions in \(\text{Cu}_x\text{Bi}_2\text{Se}_3\) are fully shelled, so a magnetic ordering seems to be highly impossible. The observed kink in \(\chi(T)\) is likely to be attributed to the charge density anomaly, or possibly the CDW transition, below which the Fermi surface (FS) is partially or fully gapped resulting in the reduction of DOS at \(E_F\). For this reason, the resistivity will sometimes display an upturn feature when the temperature crosses the \(T_k\) \([7–9]\).

However, other factors, like changes in the effective carrier mass and the scattering rate \([9]\), shall also be taken into account to interpret this upturn. Moreover, the weaker kink feature in \(\chi_{ab}(T)\) implies that the measured sample has a prominent anisotropic bulk FS since it holds a layered crystal structure and the Bi–Se quintuple layers are weakly bonded by van der Waals force.

Using the electron diffraction measurements at room temperature, the existence of CDW was reported in other Cu intercalated \(\text{Bi}_2\text{Se}_3\)-based topological insulators \([24, 25]\). However, the evidence of forming superlattice spots or diffusive order was given only in higher Cu doping content, e.g., >10 atom\% for zero-valent intercalated \(\text{Cu}_x\text{Bi}_2\text{Se}_3\) \([24]\) and 5.6 atom\% for \(\text{Cu}^{+}\) intercalated \(\text{Bi}_2\text{Te}_2\text{Se}\) \([25]\). Thus, the observed charge density anomaly in our crystal (~2 atom\% for \(\text{Cu}^{+}\) intercalation) is surprising, the content of which is close to the optimal level \([1]\). According to theoretical calculations \([26, 27]\) and angle-resolved photoemission electron spectroscopy (ARPES) experiments \([28, 29]\), the conduction bands are dominated by Bi-p orbital, which is in favor of Cu-2p orbital and would be accumulated by the electron carriers through Cu doping, leading to the lift of Fermi level. Further, since the electron doping \((x \sim 0.12)\) can induce large phonon singular electron–phonon interaction behavior and strong FS nesting in the Brillouin zone \([27]\), it is likely that the charge density anomaly is related to the \(\text{Bi}_2\text{Se}_3\) quintuple layers for nearly optimal electron-doped \(\text{Cu}_x\text{Bi}_2\text{Se}_3\).

As seen in figure 4, the \(\chi(T)\) shows an upward behavior below 50 K, which can be attributed to a tiny amount of paramagnetic defects in non-magnetic or weakly-correlated magnetic systems \([8, 30]\). Such upturn in \(\chi(T)\) is also observed in other non-magnetic materials, for instance, \(\text{Lu}(\text{Pt}_{1-x}\text{Pd}_x)\text{In}\) \([8]\) and \(\text{LaPtGe}_3\) \((T_c = 1.2\) K\) superconductors \([31]\). Furthermore, we find the \(\chi(T)\) curve between 5 K and 50 K could be well fitted by a modified Curie–Weiss law \([32–34]\), reading

\[
\chi = \chi_0 + \frac{C}{T - \Theta_{\text{cw}}},
\]

where \(C\) is Curie constant, \(\Theta_{\text{cw}}\) is Weiss temperature, \(\chi_0\) is an isotropic temperature-independent term given by \(\chi_0 = \chi_{\text{dia}} + \chi_{\text{para}} + \chi_{\text{Landau}} + \chi_{\text{vanVleck}} + \chi_{\text{Pauli}}\) in which \(\chi_{\text{core}}\) is the diamagnetic (orbital) contribution from the atomic core electrons, \(\chi_{\text{Landau}}\) is the diamagnetic (orbital) Landau susceptibility of the conduction electrons, \(\chi_{\text{vanVleck}}\) is the paramagnetic (orbital) van Vleck susceptibility, and \(\chi_{\text{Pauli}}\) is the paramagnetic (spin) Pauli susceptibility of conduction electrons. The fitted values for \(H \parallel c\) and \(H \parallel ab\) are \(\chi_0(\chi_{\text{dia}}) = -89.3 \times 10^{-6} (-159 \times 10^{-6}) \text{ cm}^3 \text{ mol}^{-1}\), \(C(C^{ab}) = -2.42 \times 10^{-4} (-1.80 \times 10^{-4}) \text{ cm}^3 \text{ K mol}^{-1}\), and \(\Theta_{\text{cw}}(\Theta_{\text{Pauli}}) = -1.25 (-0.73)\) K. The Curie constant per mole of spins is given by \([35]\)

\[
C = \frac{N_Ag^2S(S+1)\mu_B^2}{3k_B} = \frac{N_A\mu_{\text{eff}}^2\mu_B^2}{3k_B},
\]

where \(N_A\) is Avogadro’s number, \(g\) is the Landé g-factor, \(\mu_B\) is the Bohr magneton, \(k_B\) is Boltzmann’s constant, and \(\mu_{\text{eff}} = g\sqrt{S(S+1)} = \sqrt{\frac{2\mu_B}{N_Ak_B}}\) is the effective moment of a spin in unit of \(\mu_B\) (the orbital quantum number \(L = 0, J = S\)). Taking the Gaussian–CGS values of the fundamental constants, \(N_A = 6.022 \times 10^{23} \text{ mol}^{-1}, k_B = 1.381 \times 10^{-16} \text{ erg K}^{-1}\), and \(\mu_B = 9.274 \times 10^{-16} \text{ erg K}^{-1}\), \(\mu_{\text{eff}}\) is given as \(\mu_{\text{eff}} \approx \sqrt{7.99684C}\). The fitted Curie constants yield an effective moment of \(\mu_{\text{eff}} = 0.044 \mu_B\) and \(\mu_{\text{eff}} = 0.038 \mu_B\) for \(H \parallel c\) and \(H \parallel ab\), respectively. On one hand, assuming the g-factor \(g = 2.25\) typically found for \(\text{Cu}^{2+}\) ions, the net moment of \(\text{Cu}^{2+}\) ion is estimated to be 1.95 \(\mu_B\), which is much larger than experimental value of the effective moment. On the other hand, according to the Cu–Se binary diagram \([37]\), the possible precipitated CuSe phase is more favorable than the CuSe, so the paramagnetism arising from \(\text{Cu}^{2+}\) is also not promising from this point. For these reasons, we ascribe the observed anomaly at 96 K to charge density anomaly rather than the onset of magnetic ordering. The negative values of \(\Theta_{\text{Pauli}}\) and \(\Theta_{\text{Pauli}}\) indicate there exists weak antiferromagnetic exchange interactions, likely originating from the defects since the \(\text{Cu}^{2+}\) is non-magnetic with \(S = 0\).
Furthermore, one can estimate the $\chi_{\text{Pauli}}$ at $T = 0$ K using the relation [38]

$$\chi_{\text{Pauli}} \equiv \frac{g^2}{4} \mu_B^2 N(E_F), \quad (5)$$

where $N(E_F)$ is the DOS at the $E_F$. Inserting $g = 2$ into equation (5) gives $\chi_{\text{Pauli}} = (3.233 \times 10^{-5}) N(E_F)$, with $\chi_{\text{Pauli}}$ is in unit of cm$^3$ mol$^{-1}$ and $N(E_F)$ is in unit of states/eV f.u. for both spins directions (f.u. means formula unit) with $N(E_F) = N_{\text{band}}(E_F) (1 + \lambda_{\text{ep}})$. The band DOS $N_{\text{band}}(E_F)$ is given by an exact band-structure calculation. In the absence of electron–phonon (e–p) coupling, the bare Sommerfeld coefficient $\gamma_0$ is given by

$$\gamma_0 = \frac{\pi^2 k_B^2}{3} N_{\text{band}}(E_F) = 2.359 N_{\text{band}}(E_F). \quad (6)$$

Here, $\gamma_0$ is in unit of mJ mol$^{-1}$ K$^{-2}$. Taking the e–p coupling into account, the Sommerfeld coefficient is given by $\gamma = \gamma_0 (1 + \lambda_{\text{ep}})$ with $\lambda_{\text{ep}}$ is the e–p coupling constant. Combining equations (5) and (6), the $\chi_{\text{Pauli}}$ can be rewritten as

$$\chi_{\text{Pauli}} = \frac{3 \mu_B^2 \gamma_0}{\pi^2 k_B^2} = (1.370 \times 10^{-5}) \gamma. \quad (7)$$

Using the estimated value $\gamma_s = 3.46$ mJ mol$^{-1}$ K$^{-2}$ (see the discussion section for details), we obtain $\chi_{\text{Pauli}}$ is 47.4 $\times$ 10$^{-6}$ cm$^3$ mol$^{-1}$. In a normal metal or a highly doped extinct semiconductor [39], the $\chi_{\text{Landau}}$ is approximately given by

$$\chi_{\text{Landau}} = -\frac{1}{3} \left( \frac{m_r}{m^*} \right)^2 \chi_{\text{Pauli}}. \quad (8)$$

where $m_r$ and $m^*$ are the electron mass and the effective electron mass. Adopting $m^* = 0.194 m_e$ [2], we estimate $\chi_{\text{Landau}} = -419.8 \times 10^{-6}$ cm$^3$ mol$^{-1}$, showing a considerable enhancement compared to the counterpart of a normal metal. Moreover, similar to parent Bi$_2$Se$_3$ [26, 40], the Cu$_{0.09}$Bi$_2$Se$_3$ is assumed to be strongly covalent bonding in quintuple-layers with only a small ionic contribution. Since the Cu atoms are intercalated into van der Waals gap between quintuple layers, the ionic bonding is supposed to be dominant. To calculate the total core contribution, the $\chi_{\text{core}}$ values of Cu$^{1+}$, Bi$^{3+}$, Se$^{2-}$ ions are taken from the tabulated data in reference [41], $\chi_{\text{core}}$ (Cu$^{1+}$) = $-12 \times 10^{-6}$ cm$^3$ mol$^{-1}$, $\chi_{\text{core}}$ (Bi$^{3+}$) = $-25 \times 10^{-6}$ cm$^3$ mol$^{-1}$, $\chi_{\text{core}}$ (Se$^{2-}$) = $-48 \times 10^{-6}$ cm$^3$ mol$^{-1}$, yielding the core susceptibility per mole of Cu$_{0.09}$Bi$_2$Se$_3$ as $\chi_{\text{ionic}}^{\text{atomic}} = -195.1 \times 10^{-6}$ cm$^3$ mol$^{-1}$. If making use of the values of atomic core diamagnetic susceptibility calculated by the relativistic Hartree–Fock method [42] (see table 2.1 in reference [42]), $\chi_{\text{atomic}}$(Bi) = $-58.9 \times 10^{-6}$ cm$^3$ mol$^{-1}$, $\chi_{\text{atomic}}$(Se) = $-32.4 \times 10^{-6}$ cm$^3$ mol$^{-1}$, we obtain $\chi_{\text{atomic}}$ = $-215.1 \times 10^{-6}$ cm$^3$ mol$^{-1}$. One finds that there is no significant difference between $\chi_{\text{ionic}}$ and $\chi_{\text{atomic}}$. The van Vleck susceptibility is ignored due to absence of any ion with a shell that is one electron short of being half filled ($J = |S + L|$ with $J = 0$, the ground state is non-degenerate). Finally, one gets the sum $\chi_{\text{atomic}}^{\text{exp}}$ = $-58.8 \times 10^{-5}$ cm$^3$ mol$^{-1}$, which is almost four times smaller than the experimental observed value $\chi_{\text{atomic}}^{\text{exp}}$ = $15.6 \times 10^{-6}$ cm$^3$ mol$^{-1}$ at 300 K. This may be mainly caused by two reasons. One possibility is that the absolute value of $\chi_{\text{Landau}}$ is calculated based on the reported value of $m_r/m^*$ in higher Cu doping sample [2], which somehow results in uncertainty. Especially, the heat capacity measurement shows the $m^*$ in Cu$_{0.25}$Bi$_2$Se$_3$ was as large as $m^* = 2.6 m_e$ [43], strongly reducing the diamagnetic Landau contribution if this is the case. However, according to the quantum oscillation [2, 29, 44] and

| Properties | Units | Values |
|------------|-------|--------|
| $\chi^{\text{atomic}}_0$ | cm$^3$ mol$^{-1}$ | $-89.3 \times 10^{-6}$ |
| $\chi^{\text{ionic}}_0$ | cm$^3$ mol$^{-1}$ | $-159 \times 10^{-6}$ |
| $\chi_{\text{atomic}}$ | cm$^3$ K mol$^{-1}$ | $-2.42 \times 10^{-4}$ |
| $\chi_{\text{ionic}}$ | cm$^3$ K mol$^{-1}$ | $-1.80 \times 10^{-4}$ |
| $\Theta_{\text{cv}}$ | K | $-1.25$ |
| $\Theta_{\text{cv}}^{\text{Cu}}$ | K | $-0.73$ |
| $\mu_{\text{eff}}^\text{atomic}$ | $\mu_B$ | 0.044 |
| $\mu_{\text{eff}}^\text{ionic}$ | $\mu_B$ | 0.038 |

Table 1. Magnetic parameters obtained by fitting and calculation for CBS418 crystal.
optical spectroscopic studies [44, 45], the large electronic mass enhancement \(m'/m_e = 0.14–0.3\) in 
\(\text{Cu}_x\text{Bi}_2\text{Se}_3\) is not observed and thus the discrepancy from the Landau diamagnetic term seems unfavorable.

The other possibility is there might be some contributions from other kinds of paramagnetic defects [30], like the Se vacancies, as proposed in \(\text{CuInSe}_2\) compound. An exact dominant factor to explain the discrepancy between the experimental observed value and the calculated value of \(\mu_0\chi_0\) needs further explorations. Regardless of the origin of the discrepancy, the total paramagnetic susceptibility could be calculated in terms of \(\chi_p = \chi_{\text{Pauli}} + (\chi_0^\text{exp} - \chi_0^\text{atomic}) = 47.9 \times 10^{-5} \text{ cm}^3 \text{ mol}^{-1}\). The magnetic parameters are summarized in table 1.

3.3. Upper critical field
To investigate the effect of applied fields on the \(T_c\), we measured the \(M(T)\) curves under various magnetic fields. Figures 5(a) and (b) show a series of normalized \(M(T)\) curves, \(M_{\text{norm}}\), with a parallel shift of the curves for clarity. The \(T_c\) was determined by the intersection point between the extrapolation lines below and above the superconducting transition, as indicated in the inset of figure 5(a). We plot the upper critical field \(\mu_0H_{c2}^0(0)\) and \(\mu_0H_{c2}^{ab}(0)\) for \(H \parallel c\) and \(H \parallel ab\) in figure 6. The early reports on \(\text{Bi}_2\text{Se}_3\) and \(\text{Cu}_x\text{Bi}_2\text{Se}_3\) \((T_c \approx 3.5 \text{ K})\) under high pressure indicates the absence of Pauli limiting effect and an enhancement of \(\mu_0H_{c2}^{ab}(T)\) at zero-temperature limit [46, 47]. The estimated electron mean free path \(l\) is larger than the coherence length \(\xi\), suggesting the superconductivity in \(\text{Cu}_x\text{Bi}_2\text{Se}_3\) crystals grown by melt method are in the clean limit [15, 47]. We further fit the experimental quantities \(\mu_0H_{c2}^0(T)\) and \(\mu_0H_{c2}^{ab}(T)\) by the orbital-limited formula of Werthamer–Helfand–Hohenberg (WHH) theory [48] and the data above 1.8 K agree well with the WHH prediction, as is shown in figure 6. At 0 K, it yields \(\mu_0H_{c2}^0(0) = 2.22 \text{ T}\) and \(\mu_0H_{c2}^{ab}(0) = 3.60 \text{ T}\). Moreover, we have plotted the normalized upper critical field, \(h^*(t) = (\mu_0H_{c2}^0/T_c)/|d\mu_0H_{c2}^0/dT|_{T_c}\), both for \(H \parallel c\) and \(H \parallel ab\). The two \(\mu_0H_{c2}(T)\) curves collapse onto a single universal function \(h^*(t)\), as is similar to that reported by Bay et al [47]. Using the Ginzburg–Landau (GL) relations, we extract the \(\xi_{ab}(0) = 12.18 \text{ nm}\) and \(\xi_c(0) = 7.51 \text{ nm}\) at 0 K. This manifests the anisotropy ratio \(\Gamma = \xi_{ab}/\xi_c = 1.62\), which is smaller than the value determined by magnetotransport measurements [15, 43]. The anisotropy may result from the in-plane nematic superconducting state (about \(\pm 0.5\) difference in \(\Gamma\) from \(\mu_0H_{c2}^{ab}(0)\) along \(x\) and \(y\) axis) evidenced from the high-resolution heat capacity measurements of \(\text{Cu}_x\text{Bi}_2\text{Se}_3\) \((T_c \approx 3.2 \text{ K})\) [49].

3.4. Lower critical field
To determine the lower critical field \(\mu_0H_{c1}\), the demagnetization effect due to the sample shape shall be taken into account. To this end, the \(\mu_0H_{c1}\) is obtained by [50]

\[
\mu_0H_{c1}(T) = \frac{\mu_0H_p(T)}{(1 + N \cdot \chi_{\text{int})}},
\]

(9)

where \(\mu_0H_p(T)\) is the first penetration field as can be determined by the initial magnetization versus the magnetic field \(M(H)\), \(\chi_{\text{int}} = -1\) for a homogeneous bulk superconductor at most temperatures below \(T_c\).
Figure 6. Upper critical field versus temperature for (a) $H \parallel c$, (b) $H \parallel ab$. Solid lines are fitting curves by WHH theory. Inset show the normalized upper critical field versus $t$, $h^*(t) = (\mu_0 H_{c2}/T_c)/|d\mu_0 H_{c2}/dT|_{t_c}$.

Figure 7. Initial magnetization versus temperature for $H \parallel ab$ and $H \parallel c$ directions. The temperature interval $\Delta(T)$ is 0.1 K. The dash line is a linear fit.

Given the non-bulk superconductivity in Cu$_{0.20}$Bi$_2$Se$_3$, we mainly focus on the $\mu_0 H_{c1}$ for $H \parallel ab$ due to the small effect of demagnetization. Figures 7(a) and (b) show the initial $M(H)$ curves for $H \parallel ab$ and $H \parallel c$. The extending tails of diamagnetic signals above the minimum indicates a typical feature of the type-II superconductors.

To extract the $\mu_0 H_{c1}$, one common used criterion is assigning the deviation from the linear behavior of initial $M(H)$ as the $\mu_0 H_{c1}^{I}$, designated as criterion I. Figures 8(a) and (b) show the plots of $\Delta M = M_{\text{exp}} - M_{\text{fit}}$ versus field, where the $M_{\text{fit}} = s \cdot \mu_0 H$ with $s$ being the slope of the initial magnetization. Due to the small $\mu_0 H_{c1}$, the boundary of deviation from linear behavior is not well defined, especially for $H \parallel c$. This would bring much noise in the directly extracted $\mu_0 H_{c1}$. Here, we also use an alternative criterion, designated as criterion II, i.e., the intersection point as $\mu_0 H_{c1}^{II}$ from the linear fitting of $\Delta M$ slightly above the field where $\Delta M \approx 0$, as seen in figure 8(c). From these two criterions, we obtain the $\mu_0 H_{c1}$ as shown in figures 9(a) and (b). The experimental data points follow on a parabolic line, which is fitted by a GL phenomenological model as given by

$$\mu_0 H_{c1} (T) = \mu_0 H_{c1} (0) \left[ 1 - \left( T/T_c \right)^2 \right]. \tag{10}$$

We thus obtain the $\mu_0 H_{c1}^{ab} (0) = 10.6(1)$ Oe and $\mu_0 H_{c1}^{c} (0) = 6.82(6)$ Oe. The smaller value of $\mu_0 H_{c1}^{c}$ than $\mu_0 H_{c1}^{ab}$ is attributed to the non-bulk superconductivity. If one assumes $\chi_{\text{int}} = -1$, this will yield $\mu_0 H_{c1}^{ab} (0) = 11.86$ Oe and $\mu_0 H_{c1}^{c} (0) = 41.50$ Oe, respectively. To calculate the GL parameter $\kappa$, we adopt a relation by
Figure 8. Plots of selected curves of $\Delta M = M_{\text{exp}} - M_{\text{fit}}$ versus field for (a) $H \parallel ab$, (b) $H \parallel c$. (c) Zoom view of the lower field region and the criterion to determine the $\mu_0 H_{Ic}^\text{I}$ and $\mu_0 H_{Ic}^\text{II}$.

Figure 9. Lower critical field versus temperature for CBS418 determined by two criterions of (a) criterion II, (b) criterion I. The solid lines represent fitting curves by equation (10).

Taking into account the vortex core energy given by [51, 52]

$$\mu_0 H_{c1}^{\mu}(T) = \left(\Phi_0/4\pi \lambda_{ab}\lambda_c\right) (\ln \kappa_{ab} + 0.5),$$

$$\mu_0 H_{c1}^{\mu}(T)/\mu_0 H_{c1}^{\mu}(0) = 2\kappa_{ab}^2/(\ln \kappa_{ab} + 0.5).$$

The $\kappa$ is only weakly dependent on the temperature and the doping [52], as is the same case of $\Gamma$ ($\Gamma = \lambda_c/\lambda_{ab}$) in Cu$_{0.3}$Bi$_2$Se$_3$ [15, 53]. Thus the relation $\mu_0 H_{c1}^{\mu}(T) \propto 1/\lambda_{ab}\lambda_c \propto 1/\lambda_{c1}^2$ is a good approximation. For anisotropic superconductors, the $\kappa$ is defined as $\kappa_{ab} = \sqrt{\lambda_{ab}\lambda_c/\xi_{ab}\xi_c} = \lambda_{ab}/\xi_c$ and $\kappa_{c} = \lambda_c/\xi_{ab}$. Using equations (11) and (12) with $\mu_0 H_{c1}^{\mu}(0) = 6.0(3)$ Oe and $\mu_0 H_{c1}^{\mu}(0) = 9.7(3)$ Oe obtained in the clean limit (see next section), we obtain $\kappa_{ab}^I = 126.58$ and $\kappa_{ab}^II = 97.05$, respectively. The value of $\kappa_{ab}^I$ is very close to the early report ($\sim 128$) [43]. Hereafter, we mainly use this value to derive the related physical parameters as following. The penetration depth at 0 K are $\lambda_{ab}^I = 1541.57$ nm and $\lambda_{c}^I = 2499.84$ nm, yielding the averaged value $\lambda_{av}^I = (\lambda_{ab}^I\lambda_{c}^I)^{1/2} = 1811.12$ nm. Matano et al reported $\lambda_{ab} = 1038$ nm and $\lambda_{c} = 1612$ nm at 0.5 K in Cu$_{0.3}$Bi$_2$Se$_3$ [54]. Using muon spin rotation ($\mu$SR) spectroscopy, Krieger et al obtained the effective penetration depth $\lambda_{\text{eff}} \sim 1600$ nm [55]. The penetration depth at 0 K we gain is in good line with the reported results. The thermodynamic critical field is further determined by
Temperature dependence of the superfluid density, \( n_s = \lambda^{-2}(T)/\lambda^{-2}(0) \) for (a) in dirty limit, (b) in clean limit. Inset shows the lower critical field versus temperature for CBS418 for \( H \parallel ab \). The solids lines represent fitting curves by equations (16) and (14).

Table 2. Physical parameters derived from the analysis of the superconducting transitions in \( M(T) \) under magnetic field and initial \( M(H) \).

| Sample          | Properties | Units | Values     | Remarks                             |
|-----------------|------------|-------|------------|-------------------------------------|
| CBS418          | \( T_{c1} \) | K     | 4.18       | GL empirical formula, fitted by equation (10) |
|                 | \( \mu_0 H_{c1}^{ab} \) | Oe    | 10.6(1)    | In dirty limit, fitted by equation (16) |
|                 | \( \mu_0 H_{c1}^{ab} \) | Oe    | 6.82(6)    | In dirty limit, fitted by equation (16) |
|                 | \( \mu_0 H_{c1}^{ab} \) | Oe    | 10.2(4)    | In dirty limit, fitted by equation (16) |
|                 | \( \mu_0 H_{c1}^{ab} \) | Oe    | 6.1(3)     | In clean limit, fitted by equation (14) |
|                 | \( \mu_0 H_{c1}^{ab} \) | Oe    | 9.7(3)     | In clean limit, fitted by equation (14) |
|                 | \( \alpha_{ab}^{I} \) |       | 1.79       | In dirty limit, equation (16)         |
|                 | \( \alpha_{ab}^{II} \) |       | 1.759      | In clean limit, equation (14)         |
|                 | \( \alpha_{ab}^{ll} \) |       | 2.029      | In dirty limit, equation (16)         |
|                 | \( \Delta_{I}^{0} \) | meV   | 0.61(7)    | In dirty limit, equation (14)         |
|                 | \( \Delta_{I}^{0} \) | meV   | 0.73       | In clean limit, equation (14)         |
|                 | \( \mu_0 H_{c1}^{ab} \) | T     | 3.60       | In dirty limit, equation (16)         |
|                 | \( \mu_0 H_{c1}^{ab} \) | T     | 2.22       | In clean limit, equation (14)         |
|                 | \( \mu_0 H_{c2}^{ab} \) | T     | 1.20(4)    | In dirty limit, equation (14)         |
|                 | \( \mu_0 H_{c2}^{ab} \) | T     | 0.73(3)    | In clean limit, equation (14)         |
|                 | \( \mu_0 H_{c}^{ab} \) | Oe    | 211.23     | In dirty limit, equation (16)         |
|                 | \( \Gamma \) |       | 1.62       | In clean limit, equation (14)         |
| \( \xi_{ab}(0) \) | nm         |       | 12.18      | In dirty limit, equation (16)         |
| \( \xi_{c}(0) \) | nm         |       | 9.56       | In clean limit, equation (14)         |
| \( r_{ab}^{0} \) | nm         |       | 126.58     | In dirty limit, equation (16)         |
| \( r_{ab}^{0} \) | nm         |       | 1541.57    | In clean limit, equation (14)         |
| \( \mu_0 H_{c1}^{ab} \) | T     | 3.02    | [55]       | In dirty limit, equation (16)         |
| \( \mu_0 H_{c1}^{ab} \) | T     | 1.71    | [43]       | In clean limit, equation (14)         |
| \( \mu_0 H_{c1}^{ab} \) | Oe    | 4.5     | [43]       | In dirty limit, equation (16)         |
| \( \mu_0 H_{c1}^{ab} \) | Oe    | ~167    | [43]       | In clean limit, equation (14)         |
| \( \Delta_{I}^{0} \) | meV      | 0.52    |            | In dirty limit, equation (16)         |
| \( \Delta_{I}^{0} \) | meV      | 1.9     |            | In clean limit, equation (14)         |
| \( \Gamma \) |       | 1.77    |            | In dirty limit, equation (16)         |

\[
\mu_0 H_k = \sqrt{\frac{\mu_0 H_{c1}^{ab} \mu_0 H_{c1}^{cd}}{\ln \kappa_1^{ab}}} \text{, which yields 211.23 Oe at 0 K. This value is close to the reported value 167 Oe in Cu}_x\text{Bi}_2\text{Se}_3\text{ prepared by an electrochemical technique [43].}
\]

3.5. Superfluid density

To study the temperature dependence of the superfluid density \( n_s(T) \), we have analyzed the \( \mu_0 H_{c1}(T) \) extracted in criterion I. The \( n_s(T) \) is related to the \( \lambda \) as well as equivalently the \( \mu_0 H_{c1} \), which is
given by [56, 57]

\[ n_s(T) = \frac{\lambda^2(T)}{\lambda^2(0)} = \frac{\mu_0 H_{c1}(T)}{\mu_0 H_{c1}(0)}. \]

(13)

For an isotropic BCS superconductor in the clean limit, the \( n_s(T) \) is given by [58]

\[ n_s(T) = 1 - 2 \int_{\Delta}^{\infty} \frac{dE}{\Delta} \left( 1 - \frac{f(E)}{f(0)} \right) \frac{E}{\sqrt{E^2 - \Delta^2(T)}}, \]

(14)

where \( f(E) = (1 + e^{E/k_B T})^{-1} \) is Fermi distribution function, \( \Delta(T) \) is the superconducting gap function, and the quasiparticle DOS is \( N(E) = \frac{E}{\sqrt{E^2 - \Delta^2(T)}} \). In a good approximation, the gap function versus the temperature can be expressed as [58]

\[ \Delta(T) = \Delta_0 \tanh \left\{ 1.82 \left[ 1.018 \left( \frac{T_c}{T} - 1 \right) \right]^{0.51} \right\}, \]

(15)

where \( \Delta_0 \) is the energy gap at zero temperature. According to the \( \alpha \) model [59], the quantity \( \alpha \equiv \frac{\Delta_0}{k_B T_c} \) is treated as an adjustable parameter in equation (15), and the BCS value in weak coupling limit is \( \alpha_{BCS} = 1.764 \) [60]. In the dirty limit [43, 57], the \( n_s(T) \) is given by

\[ n_s(T) = \frac{\Delta(T)}{\Delta_0} \tanh \left( \frac{\Delta(T)}{2k_B T} \right). \]

(16)

To find the difference of fitting parameters between the directly determined \( \mu_0 H_{c1}^{\alpha \beta}(T) \) (criterion I) and extrapolated \( \mu_0 H_{c1}^{\alpha \beta,0}(T) \) (criterion II), we plot the \( \mu_0 H_{c1}(T) \) for \( H \parallel ab \) in figure 10. The fitted \( \mu_0 H_{c1}(T) \) curves by equations (14) and (16) are presented in insets of figure 10. The \( n_s(T) \) is obtained by normalizing \( \mu_0 H_{c1}(T) \). We find that while the values of \( \mu_0 H_{c1} \) for criterion II in the dirty limit are slightly larger than those obtained by criterion I in the clean limit, the difference between the values of \( \alpha \) derived from two criterions is more than 20%. Independent of the fitting strategy, the obtained \( \alpha \) is always smaller than \( \alpha^I \), indicating the criterion for determination of \( \mu_0 H_{c1} \) has strong influence on the magnitude of \( \Delta_0 \), i.e., the paring strength. Here, we use the criterion I to evaluate the superconducting critical parameters. Moreover, the previous reports on \( Cu_{0.10}Bi_2Se_3 \) indicates the superconducting state is in the clean limit [15], which also holds here. We obtain \( \alpha = 2.029 \pm 0.124 \), suggesting \( Cu_{0.09}Bi_2Se_3 \) is a strong-coupling superconductor. Our result from lower critical field measurements is in line with the conclusion as inferred from the specific heat experiment in \( Cu_{0.20}Bi_2Se_3 \) (\( \alpha = 1.9 \)) [43]. The physical parameters obtained in this work and from literatures are summarized in table 2.

3.6. Electrical transport properties in superconducting and normal states

To reduce the effect of inhomogeneity, we have measured a smaller piece cleaved from the bigger one used for the magnetic susceptibility measurement. Figure 11(a) shows the temperature dependence of the in-plane resistivity \( \rho_{xx}(T) \) under zero field. Different magnitudes of excitations were applied to examine whether the superconductivity is robust or filamentary. As seen in figure 11(a), we clearly observe there are two superconducting transitions with the higher one (\( T_{c1} \)) and the lower one (\( T_{c2} \)), similar to the magnetic susceptibility measurements. This double \( T_c \) phenomenon is reminiscent of the well-known intercalated iron–selenides superconductor \( MFe_2Se_2 \) (\( M = K, Rb, Tl \)) [61–63]. Actually, two systems share additional common features such as the non-bulk superconducting phase and the intercalation induced superconductivity. Interestingly, by applying high pressure [64], Sun et al discovered the reemergence of superconducting phase with a higher \( T_c \) up to 48 K in \( MFe_2Se_2 \). This high-pressure phase might be the minority phase as observed in early reports at ambient condition [61–63]. In this context, it is worth exploring whether the reemergent superconductivity with a higher \( T_c \) exists in \( Cu_0Bi_2Se_3 \). Recently, Zhou et al reported the pressure-induced reemergent superconductivity with \( T_c = 3.6 \) K at 6 GPa in its cousin material \( Sr_{0.05}Bi_2Se_3 \) [65]. As seen in figure 11(a), by applying different excitations, \( T_{c1}^{\alpha} \) is still observable up to 5 mA, indicating the robust superconductivity. In figure 11(b), we plot the current dependence of the two \( T_c \) values. By extrapolation to zero current limit, we identify two superconducting transitions occur at around 4.2 K and 3.6 K, which are in consistent with the magnetic susceptibility measurements.

Furthermore, we evaluate the \( l \) of CBS418 sample, which is critical to discern whether the specimen is in the clean or the dirty limit. To evaluate this issue, we employ the slopes of upper critical field \( d\mu_0 H_{c2}/dT \) near \( T_c \) [47, 66], Sommerfeld coefficient \( \gamma \), residual resistivity \( \rho_0 \). The relation between \( \gamma \) and the \( \mu_0 H_c \) can be found in the discussion section. In the dirty limit [66], the initial slope \( d\mu_0 H_{c2}/dT \) at \( T_c \) is given by

\[ d\mu_0 H_{c2}/dT \bigg|_{T_c} = -4.48 \times 10^4 \gamma / \rho_0, \]

(17)
where the $d\mu_0H_c^2/dT|_{T_c}$ is in unit of Oe K$^{-1}$, $\gamma$ is in unit of erg cm$^{-3}$ K$^{-2}$ and $\rho_0$ is in unit of $\Omega$ cm. Our calculation shows $d\mu_0H_c^2/dT|_{T_c} = 0.25$ T K$^{-1}$, much smaller than the experimental values $d\mu_0H_c^2/dT|_{T_c} = 1.20$ T K$^{-1}$ and $d\mu_0H_c^2/dT|_{T_c} = 0.73$ T K$^{-1}$. This suggests that the sample is not in the dirty limit.

In the clean limit [66], the $l$ can be calculated by the following relation:

$$l = 1.27 \times 10^4 \left[ \rho_0 \left( \frac{9.55 \times 10^{24} \gamma^2 T_c^3}{d\mu_0H_c^2/dT|_{T_c}} \right)^{1/2} \right]^{-1}, \quad (18)$$

where $l$ is in unit of cm. For $H || ab$ and $H || c$, we obtain $l_{ab}(l_c) = 39.86(31.09)$ nm, which are larger than the $\xi_{ab}(0)$ and $\xi_c(0)$, indicating our CBS418 sample is in the clean limit. Taking the $k_F^2 = 0.97$ nm$^{-2}$ [2], we further estimate $k_F \approx 39$, which is used as the quantity for parametrizing disorders [14, 53]. The value of $k_Fl$ is close to the lower critical Cu doping level for inducing superconductivity as prepared by electrochemical synthesis route [18], implying the CBS418 sample is weakly disordered. This may be important for observing a higher $T_c$ since the strong disorders can suppress the $T_c$ [14, 53].

### 3.7. Discussion

To find the behind reason why $T_c$ increases to 4.18 K in the CBS418 crystal, it will be helpful to estimate the DOS from the $\mu_0H_c^2$ and the $\mu_0H_c$. Due to the difficulty of measuring the heat capacity jump in non-bulk superconducting Cu$_{0.09}$Bi$_2$Se$_3$, we here indirectly estimate the electronic coefficient and the DOS at $E_F$ by the relation [67]

$$\gamma_e = \mu_0H_c^2(0)/2\pi T_c^2 = \frac{1}{3}\pi^2 k_B^2 N(E_F), \quad (19)$$

where $\gamma_e$ is the electronic Sommerfield coefficient. Using the parameters determined before, we obtain $\gamma_e = 3.46$ mJ mol$^{-1}$ K$^{-2}$ and $N(E_F) = 3.76$ states/eV atoms spin$^{-1}$ per f.u., much larger than those of ambient Cu$_{0.25}$Bi$_2$Se$_3$ obtained from the specific heat measurement [43]. We infer the enhancement of e$-$p interactions can account for the large DOS.

Before applying the BCS theory to further discuss the unusual 'high' $T_c$ for Cu$_{0.09}$Bi$_2$Se$_3$, we have examined if the adiabatic limit satisfies or not in Cu$_{0.09}$Bi$_2$Se$_3$. The BCS solution is only a good description for the systems in the adiabatic limit [68, 69], which requires the phonon energy $\hbar\omega$ must be much smaller than the carrier relaxation rate $1/\tau$, namely, $\omega \tau \ll 1$. To make such an estimation, taking the highest phonon frequency $\omega_0$ mode 175.4 cm$^{-1}$ at 3 K for Bi$_2$Se$_3$ [70] and the scattering time $\tau = 5.2 \times 10^{-14}$ s for Cu$_{0.25}$Bi$_2$Se$_3$ [2], we obtain $\omega \tau = 0.27$ as an upper bound. This indicates the Cu$_{0.25}$Bi$_2$Se$_3$ is close to violate the adiabatic limit for optical branch but is satisfied for acoustic branches due to its lower frequency in the long-wavelength limit. A brief discussion on the enhancement of $T_c$ within BCS theory can be found in supplementary material stacks.iop.org/NJP/22/053026/mmedia. Nevertheless, according to our calculation of $E_F$ for Cu$_{0.25}$Bi$_2$Se$_3$ in the following, the Debye frequency ($\omega_D$) [43] is actually comparable with the order of the $E_F$, suggesting the antiadiabatic effect needs to be considered. If the optical phonons are at play in
mediating superconductivity in Cu$_x$Bi$_2$Se$_3$, the standard BCS theory has to be modified by including the antiadiabatic effect [68, 71, 72]. Theoretically, Wan and Savrasov showed the phonon dispersions of electron-doped Bi$_2$Se$_3$ has an unusual large linewidth for both highest optical and acoustic phonons along Γ–Z direction [27]. In the antiadiabatic limit, the enhanced e–p interaction by optical phonons may be crucial generating the unusual ‘high $T_c$' for Cu$_{0.09}$Bi$_2$Se$_3$ although the acoustic phonon-mediated exotic superconductivity cannot be fully ruled out [73]. In this context, further experimental and theoretical efforts shall be made to clarify the unusual ‘high $T_c$' for Bi$_2$Se$_3$-based superconductors.

According to the discussion in supplementary material, the electron–electron (e–e) scattering dominates over the e–p scattering at low temperatures for Cu$_0.29$Bi$_2$Se$_3$. Actually, the evaluated values of $k_{\alpha}l > 1/2\pi \approx 0.16$ for various Cu doping range ($0.11 \leq x \leq 0.50$) and $k_{\alpha}l \gg 0.16$ in CBS418 sample assures the Fermi-liquid description is appropriate [53, 74]. To further estimate the Kadowaki–Woods (KW) ratio [75], $A/\gamma^2$, being a measure of the magnitude of the e–e correlation, we fit the $\rho(T)$ data using the simple formula, $\rho(T) = \rho_0 + A \cdot T^2$, to extract the e–e scattering contribution. As shown in figure 12, the fitting yields $\rho_0 = 133.41(1) \mu\Omega \cdot cm$ and $A = 0.00363(7) \mu\Omega \cdot cm \cdot K^{-2}$. For many heavy-fermion compounds [75], the KW ratio is found to follow a universal value, $A/\gamma^2 = a_0 \approx 1.0 \times 10^{-3} \mu\Omega \cdot cm \cdot mol^2 \cdot K^{-2}$, while for a lot of transition metals [76], $A/\gamma^2 \approx 4.0 \times 10^{-7} \mu\Omega \cdot cm \cdot mol^2 \cdot K^{-2}$. For the heavy-fermion, high-$T_c$ cuprates, organics, doped fullerene [77, 80, 81], the KW ratio is found to follow a universal value, $A/\gamma^2 \approx 30.3a_0$. For iron-based superconductors also lie in the similar range.

Taking the values of $a = 4.143 \text{ Å}$ and $c = 28.636 \text{ Å}$ for Bi$_2$Se$_3$ [17], $m^* = 0.194m_e$ and $k_F(k_F) = 0.97(1.26) \text{ nm}^{-1}$ [2], we obtain $A^2(A^2) = 0.00638(0.00290) \mu\Omega \cdot cm \cdot K^{-2}$, which is consistent with the fitted value. This indicates the quasi-2D conduction feature in Cu$_{0.29}$Bi$_2$Se$_3$, as expected from its layered crystal structure. Thus the 2D Fermi liquid dominates at low temperatures. From the analysis mentioned above, we infer that the large KW ratio may either result from the anisotropy as proposed in Sr$_2$RuO$_4$ [79] or the proximity to the charge density instability [7].

To classify the unconventional and conventional superconductors, Uemura et al plotted the $T_c$ versus the effective Fermi temperature $T_F$ [80, 81]. For the heavy-fermion, high-$T_c$, cuprates, organics, doped fullerene and other unconventional superconductors, the ratio $T_c/T_F$ falls within the range of 0.01 $\leq T_c/T_F \leq 0.1$. Recently, it is reported that the ratios $T_c/T_F$ for iron-based superconductors also lie in the similar range [82–86]. For the 3D superconducting systems in the clean limit [80, 81], the $T_F$ can be calculated by using the equation

$$k_F T_F = \frac{\hbar^2}{2m^*} \left( \frac{3\pi^2 n}{2} \right)^{2/3}.$$  \hspace{1cm} (21)

While in the 2D case, one can convert the volume carrier density $n$ into an areal density on the conducting planes, $n^{2D} = n \cdot d_{\text{int}}$, where $d_{\text{int}}$ is the average inter-plane distance. Therefore, one can obtain the relation

![](image)

**Figure 12.** In-plane electrical resistivity versus temperature under zero field at low temperatures for CBS418. The solid line is a fit of $\rho(T) = \rho_0 + A \cdot T^2$. 

$$
\rho_0, T = \text{fitting yields }

\mu

\text{mediating superconductivity in Cu}

\crucial generating the unusual 'high } \Gamma \text{ compounds in figure13. I nas i n g l e - b a n d 2 D m e t a l[77, 78], the coefficient planes,}

\text{a larger KW ratio would be expected according to Kriener[82–86]. For the 3D superconducting systems in the clean limit[80, 81], the effective Fermi temperature}

\text{is the average inter-plane distance. Therefore, one can obtain the relation}

\text{in CBS418 sample assures the Fermi-liquid description is appropriate [53, 74]. To further estimate the Kadowaki–Woods (KW) ratio [75], } \frac{A}{\gamma^2}, \text{ being a measure of the magnitude of the e–e correlation, we fit the } \rho(T) \text{ data using the simple formula, } \rho(T) = \rho_0 + A \cdot T^2, \text{ to extract the e–e scattering contribution. As shown in figure 12, the fitting yields } \rho_0 = 133.41(1) \mu\Omega \cdot cm \text{ and } A = 0.00363(7) \mu\Omega \cdot cm \cdot K^{-2}. \text{ For many heavy-fermion compounds [75], the KW ratio is found to follow a universal value, } \frac{A}{\gamma^2} = a_0 \approx 1.0 \times 10^{-3} \mu\Omega \cdot cm \cdot mol^2 \cdot K^{-2}, \text{ while for a lot of transition metals [76], } \frac{A}{\gamma^2} \approx 4.0 \times 10^{-7} \mu\Omega \cdot cm \cdot mol^2 \cdot K^{-2}. \text{ For the heavy-fermion, high-} T_c \text{, cuprates, organics, doped fullerene [77, 80, 81], the KW ratio is found to follow a universal value, } \frac{A}{\gamma^2} \approx 30.3a_0. \text{ For iron-based superconductors also lie in the similar range [82–86]. For the 3D superconducting systems in the clean limit [80, 81], the } T_F \text{ can be calculated by using the equation}

$$
\frac{A}{\gamma^2} = \frac{\hbar^2}{2m^*} \left( \frac{3\pi^2 n}{2} \right)^{2/3}. \hspace{1cm} (21)
$$

While in the 2D case, one can convert the volume carrier density $n$ into an areal density on the conducting planes, $n^{2D} = n \cdot d_{\text{int}}$, where $d_{\text{int}}$ is the average inter-plane distance. Therefore, one can obtain the relation
Figure 13. $T^2$ coefficient of electrical resistivity $A$ versus Sommerfeld coefficient $\gamma$ of specific heat of several systems. Experimental data beyond Cu$_x$Bi$_2$Se$_3$ was obtained from other literatures, including CDW materials [100, 101], oxides [79, 102–107], A-15 superconductors [104], heavy fermions [75, 107, 108], transition metals [76].

using the formula of non-interacting 2D electron gas

$$k_B T_F^{2D} = (\hbar^2 \pi) \frac{n^{2D}}{m^*}. \tag{22}$$

Using our previously reported value of $n$ and $m^*$ [2, 15], we calculate the $T_F^{2D} = 122.89$ K, resulting in $T_c/T_F^{2D} = 0.034$, falling into the range of $0.01 \leq T_c/T_F \leq 0.1$ for unconventional superconductors. We plot $T_c$ versus $T_F$ for various superconducting families in figure 14. This indicates the unconventional superconductivity in Cu$_x$Bi$_2$Se$_3$.

To further address the unconventional superconductivity in Cu$_x$Bi$_2$Se$_3$, we also plot the $T_c$ versus $\lambda^{-2}(0)$, another alternative type of Uemura plot [80, 81], which is directly proportional to the $n_s/m^*$. Before calculating the ratio, it is meaningful to mention another powerful technique, i.e., $\mu$SR spectroscopy, which is widely used to determine the $\lambda$ value [87]. The muon spin relaxation rate $\sigma$ is proportional to $\Delta B$, which is the inhomogeneous width of local fields. In general, the $\lambda$ is a function of the $n_s$, the $m^*$, the $\xi$, and the $l$, such as [80, 81]

$$\sigma \propto \Delta B \propto \frac{1}{\lambda^2} \propto \frac{4\pi n_s e^2}{m^* c^2} \frac{1}{1 + \xi/l}. \tag{23}$$

In the clean limit, $\xi/l \ll 1$, the relation becomes $\sigma \propto \frac{1}{\lambda^2} \propto \frac{n_s}{m^*}$. In detail, $\lambda(T)$ is related to the relaxation rate by [87]

$$\frac{\sigma_{\mu}(T)}{\gamma_\mu} = 0.06091 \frac{\Phi_0}{\lambda^2(T)}, \tag{24}$$

where $\gamma_\mu = 2\pi \times 135.5$ MHz T$^{-1}$ is the gyromagnetic ratio of muon. Some of ratios in figure 15 for other superconductors are obtained from equation (24). Taking the calculated $\lambda_{ab}(0)$, an unprecedented value of $T_c/\lambda^{-2}(0) \cong 9.86$ is obtained for Cu$_{0.09}$Bi$_2$Se$_3$, again signifying the unconventional superconductivity. The electronic parameters obtained in this work and from the literature are summarized in table 3. We note that the ratio $T_c/\lambda^{-2}(0)$ has an apparent dependence on the magnitude of the disorders in superconductors like NbN thin films [14]. In NbN, the phase fluctuations are proposed to be important for forming a pseudogap state [14]. For Cu$_x$Bi$_2$Se$_3$, we also find a similar trend according to the results reported by Kriener.
Figure 14. Plot of $T_c$ versus $T_F$. Data points of other materials were obtained from literatures, including $\beta$-Bi$_2$Pd [109], layered transition metal dichalcogenides NbSe$_2$ [80], quasi-skutteridite cubic superconductors Ca$_3$Ir$_4$ Sn$_{13}$ and Sr$_3$Ir$_4$ Sn$_{13}$ [110, 111], heavy fermions superconductors $\text{V}_x\text{Si}$ [80], doped fullerene superconductor $K_x\text{C}_{60}$ [80], Chevrel-phase superconductors LaMo$_6$S$_8$, LaMo$_6$Se$_8$ and PbMo$_6$S$_8$ [80, 81], organic superconductors (BEDT-TTF)$_2$ Cu(NCS)$_2$ (BEDT) and (TMTSF)$_2$ClO$_4$ (TMTSF) [80, 81], heavy fermions superconductors [80–82, 112], A-15 superconductor V$_3$Si [80], doped fullerene superconductor K$_3$C$_{60}$ [80], Chevrel-phase superconductors LaMo$_6$S$_8$, LaMo$_6$Se$_8$ and PbMo$_6$S$_8$ [80, 81], organic superconductors (BEDT-TTF)$_2$ Cu(NCS)$_2$ (BEDT) and (TMTSF)$_2$ClO$_4$ (TMTSF) [80, 81], bismuthates superconductors Ba$_{1-x}$K$_x$BiO$_3$ [80, 81], high-$T_c$ cuprates [80, 81, 112], high-$T_c$ iron-based superconductors [82–86], noncentrosymmetric superconductors [56, 57, 113, 114], conventional metal superconductors [81, 113], and $^4$He [82].

et al [53]. With increasing $k_Fl$, although the $T_c$ value increases, the ratio $T_c/\lambda^{-2}(0)$ tends to decrease but is still larger than those of high-$T_c$ cuprates and iron-based superconductors. The large ratio $T_c/\lambda^{-2}(0)$ was also observed in LaMo$_6$Se$_8$ [81] and in one-dimensional molybdenum cluster compounds Tl$_2$Mo$_6$Se$_8$ and In$_2$Mo$_6$Se$_8$ [88], which may arise from the considerable anisotropy or the proximity to CDW instability. Also, the Ba$_{1-x}$K$_x$BiO$_3$ with maximum $T_c = 34$ K ($x = 0.35$) derives from a diamagnetic semiconductor BaBiO$_3$ with the CDW [10], and the superconductivity occurs by suppressing the CDW through doping and accompanying a cubic–tetragonal structural phase transition. The large ratio $T_c/\lambda^{-2}(0)$ was also reported in Ba$_{1-x}$K$_x$BiO$_3$ [89–92].

Another factor to enhance the $T_c/\lambda^{-2}(0)$ is the chemical doping, as demonstrated in SnTe. Before In doping, SnTe is a superconducting narrow-band semiconductor with a low carrier density ($n_c$ up to $2 \times 10^{21}$ cm$^{-3}$, $T_c = 0.21$ K) and the $T_c/\lambda^{-2}(0)$ value is 0.03–0.15 [93]. However, the $T_c$ can be greatly improved to 4.5 K by In doping despite the decrease of $\lambda(0)$ compared to SnTe [94]. Interestingly, the $T_c/\lambda^{-2}(0)$ for Sn$_{1-x}$In$_x$Te is enhanced to the range of 1.12–1.60, implying the unconventional superconductivity. Lastly, it is meaningful to compare with other unusual superconducting semiconductors such as SrTiO$_3$ [93] and La$_3$Se$_4$ [95]. The latter one has a relatively high $T_c$ largely depending on the electron carrier density ($n_c = 1.2$ to $3.67 \times 10^{11}$ cm$^{-2}$, $T_c = 2.60$–6.86 K). Both materials undergo a lattice instability at low temperatures. For instance, La$_3$Se$_4$ has a lattice instability at 60 K, where a structure distortion from cubic to tetragonal occurs [96], while the same crystallographic change takes place at 110 K in SrTiO$_3$ [93]. It implies the higher $T_c$ and larger $T_c/\lambda^{-2}(0)$ shall be relevant to this lattice distortion. Altogether, the route of tuning the disorders or the chemical doping in semiconductors can be adopted to search more exotic superconductors.

Experimentally, the small negative susceptibility in Cu$_{0.09}$Bi$_2$Se$_3$ indicates the very weak magnetic correlations. One can further estimate the decrease of DOS from the $\chi(T)$ by the following relation:

$$\Delta \gamma = \Delta \chi \cdot (\pi k_B)^2 / (\mu_0\mu_B^2) .$$
Figure 15. Plot of $T_c$ versus $\lambda^{-2}(0)$. The data of Cu$_{0.09}$Bi$_2$Se$_3$ were extracted from our magnetization measurement, and other data points of Cu$_x$Bi$_2$Se$_3$ were obtained from magnetization and $\mu$SR measurements in previous reports [53, 55]. The data of various materials were obtained from other reports, including putative topological superconductors $\alpha$-Bi$_2$Pd and $\beta$-Bi$_2$Pd [109, 115], doped superconducting topological crystalline insulators Sn$_{1-x}$In$_x$Te [94], superconducting semiconductors SnTe, GeTe, SrTiO$_3$ [93] and La$_2$Se$_3$ [95], bismuthates superconductors Ba$_{1-x}$K$_x$BiO$_3$ [89–92], layered Bi$_2$-based superconductors of Bi$_2$O$_2$S and La$_{1+y}$Ba$_{2-x}$La$_x$O$_{2+y}$ [116–118], layered transition metal dichalcogenides NiSe$_2$, MoTe$_2$, Cu$_3$Te$_2$ and Ho$_{1.50}$Pt$_{0.50}$Te$_2$ [119–122], layered cobalt oxide Na$_{0.35}$CoO$_2 \cdot 1.3$H$_2$O [123], quasi-skutteridite cubic superconductors Ca$_3$Ir$_4$Sn$_13$ and Sr$_3$Ir$_4$Sn$_13$ [110, 124], Chevrel-phase superconductors LaMo$_6$S$_8$, LaMo$_6$Se$_8$ and PbMo$_6$S$_8$ [81], one-dimensional molybdenum cluster compounds $\text{Ti}_3\text{Mo}_6\text{Se}_{13}$ and In$_5$Mo$_6$Se$_{13}$ [88], high-$T_c$ cuprates [80, 81, 125–130], high-$T_c$ iron-based superconductors [83–85, 131–142], noncentrosymmetric superconductors Re$_3$Ta [57], Nb$_{1.18}$Re$_{0.82}$ [143], Nb$_{0.3}$Os$_{0.7}$ [114], Re$_6$Zr [144], Mo$_3$P [113], and BeAu [145], thin films of transition-metal nitrides NbN [14, 91], heavy fermions superconductors UPt$_3$, UBe$_{13}$ and $\text{Cu}_1\text{.3Hb}_{\text{Coln}}$ [146–148], and conventional metal superconductors [82, 119].

The change of the $\chi$ across the charge density anomaly transition can be calculated by the following relation:

$$\Delta \chi_{\text{Pauli}} = (1.37 \times 10^{-3}) \times 2.359 \Delta N(E_F) = (3.2318 \times 10^{-5}) \Delta N(E_F),$$

which yields $\Delta N(E_F) = 0.072 99$ states per eV f.u. for $H \parallel c$ and 0.006 16 states per eV f.u. for $H \parallel ab$. This indicates the strong anisotropy of the FS. The previous results from the ARPES [29] and quantum oscillations [97] showed that the Cu$_{0.09}$Bi$_2$Se$_3$ has a 2D cylinder FS with high carrier density. This potentially brings the FS nesting. As is well known, the FS nesting could be a crucial ingredient for the CDW [98], which again enhances the e–p interactions. In a previous theoretical work, Wan and Savrasov proposed there existed a strong FS nesting at small wave vectors $q$ along the $\Gamma(000) - Z(\pi\pi\pi)$ direction in Cu$_{0.09}$Bi$_2$Se$_3$ [27], which could greatly enhance the e–p coupling constant $\lambda$. Actually, with increasing the carrier density, the FS exhibits a tendency to become a 2D-like one [29]. In principle, the FS nesting can assist the formation of CDW state [98]. However, the strong spin–orbit coupling and/or the e–p coupling shall also be considered for the origin of the charge density anomaly in CBS418 [8, 99]. In any case, the nature of the observed charge density anomaly needs to be further clarified, which is believed to relate to the ‘high-$T_c$’ superconductivity in doped topological insulators.
Table 3. Electronic properties of CBS418 sample in present work (PW).

| Properties | Units | Values | References |
|------------|-------|--------|------------|
| $m$ | $m_e$ | 0.194 | [2] |
| $k_{F1}$ | nm$^{-1}$ | 0.97 | [2] |
| $k_{F1}/k_{F2}$ | | 1.3 | [2] |
| $l_{bb}$ | nm | 39.86 | PW |
| $l_{ll}$ | nm | 31.09 | PW |
| $k_{F1} \pm 0.39$ | | PW [2] |
| $A_{ee}$ | $\mu\Omega$ cm K$^{-2}$ | 0.00363(7) | PW |
| $A_{ee}^{\rm cal}$ | $\mu\Omega$ cm K$^{-2}$ | 0.00638 | PW |
| $A_{cal}$ | $\mu\Omega$ cm K$^{-2}$ | 0.00290 | PW |
| $T^D_{l}$ | K | 122.89 | PW [2] |
| $T^D_{e}$ | | 0.034 | PW |
| $\gamma_s$ | mJ mol$^{-1}$ K$^{-2}$ | 3.46 | PW |
| $N(E_F)$ | states/eV atoms spin$^{-1}$ per f.u. | 3.76 | PW |
| $\gamma_s$ | mJ mol$^{-1}$ K$^{-2}$ | 1.95 | [43] |
| $N(E_F)$ | states/eV atoms spin$^{-1}$ per f.u. | 1.34 | [43] |
| $T_c/\lambda_{ab}^2(0)$ | K $\mu$m$^{-2}$ | 9.94 | PW |
| $T_c/\lambda_{eff}^2(0)$ | K $\mu$m$^{-2}$ | 10.24 | [55] |

4. Summary and conclusions

To summarize, we report the investigations of the physical properties of Cu$_{0.09}$Bi$_2$Se$_3$ single crystal with $T_{\text{onset}} = 4.18$ K. According to the magnetic susceptibility, a charge density anomaly at 96 K was observed. This leads to a conjecture of the possible coexistence of charge density wave and superconductivity in Cu doped Bi$_2$Se$_3$ topological insulator. We determined the $\mu_0H_c^2(T)$ and $\mu_0H_c^1(T)$ from the magnetization curves for Cu$_{0.09}$Bi$_2$Se$_3$. In the clean limit, the $\mu_0H_c^1(0)$ was extracted to be 6.0 Oe for $H \parallel ab$. The experimental energy gap ratio was $\Delta_0/k_B T_c = 2.029 \pm 0.124$, indicating it is a strong-coupling superconductor. From electrical transport measurements, the 2D Fermi liquid behavior is found at low temperature and a high Kadowaki–Woods ratio is determined to be $A/\gamma^2 = 30.3$. Furthermore, we revealed an unprecedented high ratio $T_c/\lambda_{ab}^2(0) \cong 9.86$ with $\lambda_{ab}(0) = 1541.57$ nm as well as a high ratio of $T_c/T^D_{l} = 0.034$ as comparable to those of high-$T_c$ cuprates and iron-based superconductors. These results demonstrate the unconventional superconducting mechanism in Cu$_{0.09}$Bi$_2$Se$_3$ according to Uemura's regime. Finally, we propose the enhanced $T_c$ up to 4.18 K results from the increased density of states at $E_F$ and the strong electron–phonon interaction induced by the charge density instability. Our results suggest the higher $T_c$ in Cu$_x$Bi$_2$Se$_3$ could be further achieved by gating-technique or high pressure technique, as realized in iron–selenides superconductors.

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