Critical current density and vortex phase diagram in the superconductor $\text{Sn}_{0.55}\text{In}_{0.45}\text{Te}$

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Critical current density and vortex pinning dynamics have been studied in the superconductor $\text{Sn}_{0.55}$ In_{0.45}Te. Analysis of the temperature-dependent lower critical field shows that it has a weakly anisotropic single energy gap. The critical current density $J_c(0)$ and pinning potential $U_0(H)$ values reach as high as 2.56×10^3 A/cm² at 1.8 K and 2.1×10^3 K at $\mu_0 H = 0.01$ T, respectively. Based on the collective pinning model, we demonstrate the coexistence of vortex pinning regimes in $\text{Sn}_{0.55}\text{In}_{0.45}$ Te. One is a δT_c pinning regime induced by the spatial fluctuations of the transition temperature in a low field. The other is a dominantly δ*l* pinning regime associated with the spatial variations of the charge-carrier mean free path in a higher field. This causes a nonconstant exponent of the power-law behavior $J_c(T) \propto H^n$. A very weak vortex fluctuation is unveiled by a narrow separation between the irreversibility field $\mu_0 H_{\text{irr}}(T)$ and upper critical field $\mu_0 H_{\text{c2}}(T)$ in the vortex phase diagram. We discuss the potential application in superconducting electronics like the single-photon detector in thin film form.

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I. INTRODUCTION

SnTe has been widely studied as a prototype topological crystalline insulator with a narrow direct band gap of 0.18– 0.31 eV $[1-3]$. By alloying $[4]$ and strain-engineering $[5]$, it achieves great thermoelectric performance. Through Indium doping [\[6,7\]](#page-6-0), enhanced superconductivity appears with a tunable superconducting transition temperature (T_c) by adjusting the In content, ranging from 24 mK to 5.1 K $[6-8]$ $[6-8]$. At the lower In-doping side for $x < 0.12$ [\[9\]](#page-7-0), the observation of a zero-bias conductance peak in $\text{Sn}_{1-\alpha}\text{In}_{x}\text{Te}$ ($x = 0.045$) with preserved topological surface state indicates that it hosts topological superconductivity [\[10,11\]](#page-7-0). Nevertheless, the nuclear magnetic resonance (NMR) experiment shows that the polycrystalline $Sn_{0.96}In_{0.04}Te$ is a spin-singlet superconductor [\[12\]](#page-7-0). When *x* exceeds 0.12 [\[9\]](#page-7-0), one theoretical work points out that the $Sn_{1-x}In_xTe$ becomes a Dirac semimetal, or a Weyl semimetal through a ferroelectric effect or Jahn-Teller distortion although lacking of further experimental verifications. The combined possible topological nontrivial band structure and superconductivity in Sn_{1−*x*}In_{*x*}Te makes it a fertile plat-

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form for exploring a variety of fermions and new topological states of matter [\[9,11\]](#page-7-0).

Despite intensive studies that have been reported on the bulk $\text{Sn}_{1-x}\text{In}_x\text{Te}$ superconductors, we need to note that the nature of its superconducting state is still under debate. The London penetration depth measurements on $x = 0.10$ and 0.45 [\[13\]](#page-7-0), muon-spin spectroscopy measurements on $x = 0.38 - 0.45$ [\[14\]](#page-7-0), and thermal conductivity measurements on $x = 0.40$ [\[15\]](#page-7-0) have evidenced the fully gapped superconducting state, which can be described by the *s*wave Bardeen-Cooper-Schrieffer (BCS) theory [\[13,14\]](#page-7-0). These results lead to the controversy on its topological superconductivity nature for $Sn_{1-x}In_xTe$ superconductors. Furthermore, based on superconducting density functional theory (SCDFT) and tunneling spectroscopy, Nomoto *et al.* revealed that the spin-orbit coupling (SOC) can largely enhance T_c at a low carrier region and the T_c of Sn_{1−*x*}In_{*x*}Te can be quantitatively described by the conventional *s*-wave BCS scenario [\[16\]](#page-7-0). In addition, the disorder induced by the increase of Sn content in InTe or In content in SnTe can broaden the features of the density of states (DOS) spectral function [\[7,](#page-6-0)[16\]](#page-7-0), which leads to the redistribution of spectral weight and the mild changes of DOS at Fermi level $N(E_F)$ with the dopant amount. It is also shown that the evolution of $N(E_F)$ positively correlates the T_c vs. *x* in $\text{Sn}_{1-\alpha} \text{In}_{\alpha}$ Te superconductors [\[7\]](#page-6-0).

Recently, to explore the junctions or devices applications, thin films of Sn1−*^x*In*x*Te with various In-doping contents have

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been successfully grown on substrates by molecular beam epitaxy and pulsed-laser deposition $[17–20]$. Sn_{1−*x*}In_{*x*}Te thin films grown on a $Bi₂Te₃$ buffer layer show a coexistence of bulk and topological surface superconductivity [\[18\]](#page-7-0). Masuko *et al.* reported superconducting $Sn_{1-x}In_xTe$ thin films with $x = 0.66$ [\[20\]](#page-7-0), beyond the bulk solubility limit of $x = 0.5$. These studies pave the way for their possible device applications. For instance, a combination of the tunable T_c , large normal state resistivity, with the higher slope of upper critical field vs. temperature in superconducting thin films makes them being candidate materials for single-photon detector (SNSPD) application [\[21,22\]](#page-7-0). Also, the avoided large $N(E_F)$ due to In doping also benefits for developing a SNSPD using $Sn_{1-x}In_xTe$ candidates [\[7](#page-6-0)[,16\]](#page-7-0). However, the critical current density and vortex pinning dynamics in Sn1−*^x*In*x*Te remain unexplored, which are important ingredients for applications [\[23\]](#page-7-0). Moreover, there is a discrepancy on the energy gap ratio determined by London penetration depth and muon-spin spectroscopy for the $x = 0.45$ composition [\[13,14\]](#page-7-0), which needs a further investigation using an alternative technique.

Here, we choose $Sn_{0.55}In_{0.45}Te$, the optimal doping level with maximum T_c for the ambient pressure growth, to investigate the critical current density and vortex dynamics in its Dirac semimetal superconductor form [\[6,7\]](#page-6-0). Through the magnetic and electrical transport measurements, we established the vortex pinning phase diagram and derived a few key physical parameters, which are important for superconducting electronics applications.

II. EXPERIMENTAL METHODS

A. Sample synthesis and characterizations

The experimental SnTe and $Sn_{0.55}In_{0.45}Te$ bulk single crystals were synthesized via the melt method in a vertical furnace. The Sn (6N, Alfa Aesar), In (4N, Alfa Aesar), and Te (6N, Alfa Aesar) granules with a stoichiometric ratio were put into an evacuated quartz tube with a conical bottom and then sealed by the hydrogen-oxygen flame. The mixtures were heated at 1123 K for 12 h followed by cooling to 1018 K at a rate 1.5 K/h and annealing at 1018 K for more than 24 h, and the SnTe and $Sn_{0.55}In_{0.45}Te$ ingots were obtained by quenching into an ice water bath. Specially rectangle-shaped single crystals of SnTe were found on the top part of the ingot.

A piece of rectangle-shaped specimen (size: 2.75×2.66× 0.80 mm³) was used for the magnetic and electrical transport measurements. Electrical transport measurements were performed using the standard four-probe method on a DynaCool physical property measurement system (PPMS). Isothermal magnetization hysteresis loops (MHLs) and initial magnetization measurements were conducted on a superconducting quantum interference device vibrating sample magnetometer. The temperature range was between 1.8 and 5.1 K with an interval of $\Delta T = 0.3$ K. The ramping rate for collecting MHLs was 20 Oe/s and 1.6 Oe/s for the initial magnetization.

The phase purity was confirmed by the laboratory PANalytical Empyrean x-ray diffraction (XRD) system with a Cu K α 1 radiation source. Single crystal XRD measurements were performed on a Bruker D8 VENTURE diffractometer with Ag target ($\lambda = 0.561$ Å) and Mo target ($\lambda = 0.717$ Å)

for SnTe and In-doped samples. The single crystal XRD data were refined by ARPEX3 software and the structural parameters can be found in Supplemental Material [\[24\]](#page-7-0). The Rietveld refinement of the powder XRD patterns was further processed using the GSAS program package with the user interface EXPGUI [\[25,26\]](#page-7-0). The compositions were determined by energy-dispersive x-ray spectroscopy (EDS) and the inductively coupled plasma atomic emission spectroscopy (ICP-AES).

B. Theoretical calculations

All calculations were done in the framework of DFT [\[27,28\]](#page-7-0) using the projector augmented wave [\[29,30\]](#page-7-0) pseudopotential method, as implemented in the VASP package [\[31,32\]](#page-7-0). The electron exchange-correlation functional was treated in generalized gradient approximation as proposed by Perdew-Burke-Ernzerhof. The valance electrons were considered as $4d^{10}5s^25p^2$ for Sn, $4d^{10}5s^25p^1$ for In and $5s^25p^4$ for Te. The kinetic energy cutoff for the plane-wave basis set expansion was set to 360 eV in all cases to avoid Pullay stress. A Γ -centered Monkhorst-Pack grid of $7 \times 7 \times 7$ and $10\times10\times10$ *k* points were used for the geometry optimization and calculation of DOS, respectively. The energy convergence of 1.0×10^{-8} eV was used for the electronic energy minimization steps. During relaxation, Feynman forces on each atom were minimized until they were less than 0.001 eV/A . SOC was taken into account in all calculations.

III. RESULTS AND DISCUSSIONS

A. Electronic band structures

Band structure calculations were performed to characterize the topological properties of SnTe and $Sn_{0.55}In_{0.45}Te$. For the bulk $Sn_{1-x}In_xTe$ systems, we used a 2×2×2 supercell at the face-centered-cubic primitive cell basis, containing eight cations (Sn/In) and eight anions (Te), as illustrated in Figs. $1(a)$ and $1(b)$. Considering the complexity of calculations and the similar electronic structure of $x = 0.45$ and 0.5 doping [\[9\]](#page-7-0), we used the structure of Sn and In evenly distributed in a 1:1 ratio. In Figs. $1(d)$ and $1(e)$ we plotted the electronic band structure and projected DOS by considering SOC effect. The high symmetry *k* path in the Brillouin zone for both structures is shown in Fig. $1(c)$. The bulk band structure of SnTe shows a gap of less than 0.2 eV at the Fermi energy level (E_F) , with the band inversion occurring at the Γ point, a good ingredient for a nontrivial topological surface state [\[2](#page-6-0)[,9\]](#page-7-0). A previous report of band inversion at the L point does not contradict our results as the L point is folded onto the Γ point with supercells consideration [\[33\]](#page-7-0). After alloying In atoms, the E_F is significantly lowered and $N(E_F)$ increases, leading to a metallic state for $Sn_{0.55}In_{0.45}Te$. Moreover, band inversion can still be observed at the Γ point but at a higher level near 0.9 eV. The *p* electrons of Sn, *s* electrons of In, and p electrons of Te hybridize around E_F , giving a large $N(E_F)$ that may lead to an increase in the T_c . Actually, previous reports show that the doped of In in SnTe will lead to the increase of $N(E_F)$ with adding more In atoms for $x \le 0.5$, which correlates with the evolution of T_c with In content and

FIG. 1. Supercell structures at the face-centered cubic primitive cell basis for (a) SnTe and (b) $Sn_{0.5}In_{0.5}Te$. (c) High symmetry *k* path in the Brillouin zone. Band structures include SOC and projected DOS for (d) SnTe and (e) $Sn_{0.5}In_{0.5}Te$. The Fermi energy (blue dashed lines) is set to zero.

may account for the enhancement of T_c in In-doped SnTe materials [\[7](#page-6-0)[,34\]](#page-7-0).

B. Structure, composition, and superconducting properties

The cubic symmetry and phase purity are confirmed by powder XRD patterns, as shown in Fig. $2(a)$. Note that the weak reflection at ∼2θ angle 33 degree is likely from the

FIG. 2. (a) XRD patterns. (b) EDS spectrum. (c) and (d) show the temperature-dependent magnetic susceptibility and resistivity for $Sn_{0.55}In_{0.45}Te$.

tetragonal InTe phase [\[6\]](#page-6-0). The Rietveld refinement yields the lattice parameter $a = 6.3092(1)$ and $a = 6.2507(2)$ for SnTe and $Sn_{0.55}In_{0.45}Te$, respectively. Representative EDS profiles are plotted in Fig. $2(b)$. Comparing to the pristine sample, the appearance of the excitation peaks clearly indicates the alloying In atoms in the $Sn_{1-x}In_xTe$ ($x = 0.45$) sample. To accurately analyze the composition, we have used ICP-AES technique, which gives $Sn_{0.98}Te$ and $Sn_{0.54}In_{0.45}Te$. Hereafter, we will use the nominal composition for discussions and comparison. The temperature-dependent resistivity and magnetic susceptibility at low temperature were measured to detect the superconductivity for $Sn_{0.55}In_{0.45}Te$, as shown in Fig. 2(c). The sharp diamagnetic transition below 5.10 K and the high shielding volume of 94% at 1.8 K indicate a homogeneous bulk superconductivity. This is further corroborated by the sharp zero-resistivity transition at 5.04 K shown in Fig. $2(d)$.

C. Lower critical field

Initial magnetization versus applied field $M(H)$ curves was measured to verify the temperature-dependent lower critical field $\mu_0 H_{c1}$ that determines the superfluid density, as displayed in Figs. $3(a)$ and $3(b)$. The demagnetization effect was taken into account. The lower critical field μ_0H_{c1} can be obtained by [\[35\]](#page-7-0)

$$
\mu_0 H_{\rm cl}(T) = \frac{\mu_0 H_{\rm pl}(T)}{1 + N \cdot \chi_{\rm int}},\tag{1}
$$

where $\mu_0 H_{\text{pl}}(T)$ is the first penetration field. *N* is the demagnetization factor as is calculated by $N = [1 +$ $(3/4) \cdot (c/a)(1 + (a/b))]^{-1}$ [\[35\]](#page-7-0), where *a*, *b*, *c* are dimensions for a cuboid. Here, *N* is calculated to be 0.69 for our sample. The first $\mu_0 H_{p1}(T)$ is determined by the deviation from the linear behavior of the initial $M(H)$, as indicated in Fig. [3\(b\)](#page-3-0) and its inset for the plots of $\Delta M = M_{\rm exp} - M_{\rm fit}$ versus field. Here, the $M_{\text{fit}} = s \cdot \mu_0 H$ with *s* being the slope of the initial magnetization. The $\mu_0 H_{c1}(T)$ data follows on an empirical

FIG. 3. (a) Initial magnetization curves vs. applied field of $Sn_{0.55}In_{0.45}Te$ at various temperatures. (b) Initial diamagnetization at 1.8 K. Solid line presents a linear fitting curve. Inset shows the difference between experimental and linear fitting data. (c) Temperature-dependent $\mu_0 H_{c1}$. Solid lines are theoretical fitting curves.

parabolic line [\[36\]](#page-7-0)

$$
\mu_0 H_{c1}(T) = \mu_0 H_{c1}(0) \bigg[1 - \bigg(\frac{T}{T_c}\bigg)^2 \bigg].
$$
 (2)

Moreover, as shown later, our sample satisfies the local limit since $\xi(0)$ is much smaller than London penetration depth $\lambda(0)$. In the local London model, the temperature-dependent superfluid density [\[37,38\]](#page-7-0)

$$
n_s(T) = \frac{\lambda^{-2}(T)}{\lambda^{-2}(0)} = \frac{\mu_0 H_{c1}(T)}{\mu_0 H_{c1}(0)},
$$
\n(3)

can be analyzed using the $\mu_0 H_{c1}(T)$. Given the metallic behavior and assuming a spherical Fermi surface, the mean free path can be estimated by $l = \hbar k_F / \rho_0 n e^2$ with the Fermi wave number $k_F = (3\pi^2 n)^{1/3}$ and residual resistivity ρ_0 , yielding $l = 127(3)$ nm and 0.6(2) nm for SnTe and Sn_{0.55}In_{0.45}Te [\[24\]](#page-7-0). The *l* value is much smaller than its superconducting coherence length at the 0 K limit $\xi(0) = 13.8(1)$ nm, indicating the dirty limit in $Sn_{0.55}In_{0.45}Te$. The significant decrease of product $k_F l$ that quantifies the disorder in $\text{Sn}_{0.55}\text{In}_{0.45}$ Te compared with SnTe implies the prominent disorders scattering due to the introduction of Indium atoms. In the dirty limit [\[37,39\]](#page-7-0), the $n_s(T)$ is given by

$$
n_s(T) = \frac{\Delta(T)}{\Delta_0} \tanh\left(\frac{\Delta(T)}{2k_B T}\right).
$$
 (4)

The gap function can be approximately expressed as [\[40\]](#page-7-0)

$$
\Delta(T) = \Delta_0 \tanh\left(1.82 \left[1.018\left(\frac{T_c}{T} - 1\right)\right]^{0.51}\right),\qquad(5)
$$

where Δ_0 is the energy gap at zero temperature. According to the α model [\[41\]](#page-7-0), the quantity $\alpha = \Delta_0 / k_B T_c$ can be treated as an adjustable parameter. We obtained $\alpha = 1.18 \pm 0.12$ from the best fitting, smaller than the standard BCS value α_{BCS} = 1.764 in a weak coupling limit [\[42\]](#page-7-0). Both fitting results are shown in Fig. $3(c)$. The obtained ratio is in good agreement with the London penetration depth measurements ($\alpha = 1.18$

for $Sn_{0.55}In_{0.45}Te$ [\[13\]](#page-7-0), but smaller than the value from muonspin rotation or relaxation (μ SR) measurements [$\alpha = 1.89(3)$ for $Sn_{0.55}In_{0.45}Te$ in clean limit] [\[14\]](#page-7-0). In Ref. [14], Saghir *et al.* have analyzed their $\lambda^{-2}(T)$ data using the BCS *s*-wave model in both clean and dirty limit. Also, they show that the Δ_0 magnitudes in the clean limit are larger than those of dirty limit, indicating a smaller gap ratio in dirty limit as demonstrated for $x = 0.45$ herein. In addition, one would expect a variable gap ratio if further considering the gap anisotropy during analyzing the $\lambda^{-2}(T)$ data by μ SR measurements. Nevertheless, this calls for further studies. Furthermore, the results by thermal conductivity probes indicate that the $Sn_{0.6}In_{0.4}Te$ has a full superconducting gap $[15]$. Consequently, the small gap ratio of the $Sn_{0.55}In_{0.45}Te$ sample can be attributed to its weakly anisotropic single gap [\[13,42\]](#page-7-0).

D. Critical current density and vortex pinning mechanism

The critical current density $J_c(H)$ is the maximum unimpeded current that a superconductor can transmit per unit cross section at a given temperature and magnetic field, one of the key parameters to evaluate their superconducting applications. The extended Bean critical state model was adopted to estimate the $J_c(H)$ of $\text{Sn}_{0.55}\text{In}_{0.45}\text{Te}$ [\[43\]](#page-7-0). Figure [4\(a\)](#page-4-0) shows the MHLs between 1.8 and 5.1 K with a temperature interval of $\Delta T = 0.3$ K. For a rectangle-shaped sample with field parallel to the shortest edge, the field-dependent J_c is given by

$$
J_c(H) = \frac{20\Delta M}{w\left(1 - \frac{w}{3l}\right)},\tag{6}
$$

where *w* and *l* are the width and length of the sample, and ΔM is the difference between the magnetization values for the increasing and decreasing field, the sample's thickness *t* is set as $t < w < l$ and the magnetization is in the unit of emu/cm³. The $J_c(0)$ at $T = 1.8$ K is extracted to be $J_c(0) = 2.56 \times 10^3$ A/cm², comparable to those iron-based high-temperature superconductors FeS [\[44\]](#page-7-0), FeSe [\[45\]](#page-7-0), and

FIG. 4. (a) Magnetic hysteresis loops below 5.1 K of Sn0.55In0.45Te. (b) Log-log plot of critical current density *Jc* versus applied field. The J_c is extracted by Bean's model. The inset shows the exponent versus temperature. (c) Normalized $j_c(t)$. The inset shows the percentage of the pinning regime. (d) The normalized pinning energy $g(t) = U(t)/U(0)$ versus temperature.

Mn-doped $K_xFe_{2-y}Se_2$ [\[46\]](#page-7-0). Above a crossover field H_{cr} at 32 Oe, as indicated in Fig. $4(b)$, the $J_c(H)$ curves follow a power-law behavior $J_c \propto H^n$, indicative of a collective pin-ning signature [\[47–49\]](#page-7-0). However, the $J_c(H)$ deviates from the power-law scaling in a high field. The power-law behavior implies a fairly dense distribution of weak pins exists in $Sn_{0.55}In_{0.45}Te$ [\[50\]](#page-7-0), analogous to that in $Cu_{0.10}Bi_2Se_3$ [\[49\]](#page-7-0). According to the strong pinning theory $[23]$, the $J_c(H)$ is nearly constant for $H < H_{cr}$ while $J_c(H) \propto H^{-5/8}$ for $H >$ H_{cr} , where H_{cr} is a crossover field in the $J_c(H)$ curve. At 1.8 K, the fitting exponent is $n = -0.682(2)$. In principle, this could approach the theoretical $n = -0.625$ by further decreasing temperature, as shown in the inset of Fig. $4(b)$. However, the *n* value has an apparent temperature dependence and tends to decline at higher temperature. This is unexpected in the strong pinning theory, which predicts a temperature-independent exponent [\[23\]](#page-7-0).

To understand this puzzling behavior in $J_c(H)$, we conducted a pinning regime analysis for $Sn_{0.55}In_{0.45}Te$. Given it belongs to a large- κ superconductor with Ginzburg-Landau (GL) parameter $\kappa = 25(1)$, this leads us to safely assume that its core interaction plays a major role in the vortex pinning dynamics [\[49\]](#page-7-0). Two categories of vortex pinning, δl pinning and δT_c pinning, were considered. Based on the collective pinning theory [\[51\]](#page-7-0), the temperature-dependent normalized $J_c(T)$ in a single-vortex pinning regime can be described by $J_c^{\delta l}(t)/J_c^{\delta l}(0) = (1 - t^2)^{5/2}(1 + t^2)^{-1/2}$ for δl pinning, and $J_c^{\delta T_c}(t)/J_c^{\delta T_c}(0) = (1 - t^2)^{7/6}(1 + t^2)^{5/6}$ for δT_c pinning with $t = T/T_c$. In addition, the functions $g(t) =$ $U(t)/U(0)$ characterizing the temperature-dependent pinning energy are given by $g^{\delta l}(t) = 1 - t^4$ for δl pinning, and $g^{\delta T_c}(t) = (1 - t^2)^{1/3} (1 + t^2)^{5/3}$ for δT_c pinning. Also, taking the mixed regimes into account, we used a combined relation, $J_c(t) \propto (1 - p) J_c^{\delta l}(t) + p J_c^{\delta T_c}(t)$, to fit the experimental $J_c(T)$ data, where p is the contribution percentage by δT_c pinning. Finally, we plot the $J_c(t)$ and $g(t)$ data at different magnetic fields in Fig. $4(c)$. Intriguingly, the pinning is governed by the δT_c pinning in a low field and gradually transformed to a δ*l* dominant pinning in a higher field. The inset of Fig. 4(c) shows the evolution of the contribution of δT_c pinning with simultaneously increasing temperature and field. Under 1000 Oe, the vortex pinning is governed by δ*l* pinning, and the normalized pinning energy $g(t)$ follows the δl pinning regime well, as shown in Fig. $4(d)$. This result supports that the peculiar nonconstant exponent above H_{cr} arises from the synergetic interplay of the δT_c pinning and δl pinning in Sn_{0.55}In_{0.45}Te. Benefiting from the close-by δT_c pinning, the Sn_{0.55}In_{0.45}Te material may find high *Jc* applications at a low temperature and small field.

From the $J_c(H)$ data, the irreversibility critical field $\mu_0 H_{irr}$ can be determined as the point at extrapolated zero J_c value from Kramer plot [\[52\]](#page-7-0), $J_c^{1/2}H^{1/4}$ as a function of *H*. Figure 5 shows the Kramer plot and the linear extrapolation of solid lines lead to the $\mu_0 H_{irr}$ at $J_c = 0$. Inset of Fig. 5 presents the temperature dependent $\mu_0 H_{irr}(T)$, which can be fitted by the $\mu_0 H_{irr}(T) = \mu_0 H_{irr}(0)(1 - T/T_c)^n$. The best fitting yields $n = 1.24(3)$, comparable to that of Mn-doped K_xFe_{2−*y*}Se₂ [\[46\]](#page-7-0).

E. Thermally activated flux flow behavior and pinning potential

We conducted temperature-dependent resistivity $\rho(T)$ measurements under various fields to determine the irreversibility critical field $\mu_0 H_{irr}$, as shown in Fig. [6\(a\).](#page-5-0) The criterion for determining the T_c and irreversibility temperature

FIG. 5. Kramer plot of the $J_c(H)$ data: $J^{1/2}H^{1/4} \sim H$. Inset presents the temperature dependent $\mu_0 H_{irr}(T)$ and solid line indicates the fitting curve.

FIG. 6. (a) Temperature dependence of resistivity under various fields for $\text{Sn}_{0.55}\text{In}_{0.45}\text{Te}$. (b) Criterion determining the T_c and T_{irr} . (c) Arrhenius plot: $\ln[\rho(T, H)] \sim 1/T$. (d) Field dependence of pinning potential U_0 . Solid lines are power-law fitting results.

 T_{irr} is displayed in Fig. 6(b), where the $\rho(T)$ drops to 10% of the normal state resistivity [\[53\]](#page-7-0). The energy dissipation due to vortex motion causes the broadened resistivity at around T_c under magnetic fields, which can be described by the thermally activated flux flow (TAFF) model. In the framework of the TAFF model [\[54\]](#page-7-0), the temperature and field-dependent resistivity $\rho(T, H)$ can be described by

$$
\rho(T, H) = \rho_0 \exp\bigg(-\frac{U(T, H)}{T}\bigg),\tag{7}
$$

where ρ_0 is a parameter, and $U(T, H)$ is the activation energy for the vortex bundle hopping. The Arrhenius relation

$$
\ln[\rho(T, H)] = \left[\ln \rho_0 + \frac{U_0(H)}{T_c} \right] - \frac{U_0(H)}{T}
$$
 (8)

can be derived by assuming $U(T, H) = U_0(H)(1 - t)$ with $t = T/T_c$ and U_0 is the pinning potential [\[55\]](#page-7-0). The U_0 can be extracted from the Arrhenius-plot, $ln[\rho(T, H)]$ ∼ $1/T$, as shown in Fig. 6(c). The obtained $U_0(H)$ is plotted in Fig. $6(d)$, which reaches 2.1×10^3 K at $\mu_0 H =$ 0.01 T. This value is comparable to the reported values of FeS_{1−*x*}Se_{*x*} [\[56\]](#page-7-0), K_{0.8}Fe_{1.97}Mn_{0.03}Se₂ (*U*₀(0.25 T) = 5.1×10³ K) [\[57\]](#page-7-0), NdO_{0.82}F_{0.18}FeAs ($U_0(0.2 \text{ T}) = 4 \times 10^3 \text{ K}$) [\[58\]](#page-7-0), and $Bi_2Sr_2CaCu_2O_8$ ($U_0 = 3 \times 10^3$ K) [\[54\]](#page-7-0). Generally, the $U_0(H)$ obeys a power law $U_0 \propto H^n$ with *n* as a characteristic parameter for collective flux creep. The best fitting yields $n_1 = 0.21(5)$ below $\mu_0 H_{\text{int}} = 0.24$ T while $n_2 = 0.83(7)$ beyond $\mu_0 H_{\text{int}}$, at which the two fitting lines intersect, as displayed in Fig. $6(d)$.

F. Vortex phase diagram

Figure 7 summarizes the vortex phase diagram for the $Sn_{0.55}In_{0.45}Te superconductor, which divides into a Meissner$ state, vortex solid, vortex liquid, and normal states. Hereafter, we will focus on the irreversibility field that locates fairly close to the $\mu_0 H_{c2}(T)$. To calculate the GL parameter κ and London penetration depth λ , we employed a relation consid-

FIG. 7. Vortex phase diagram of $Sn_{0.55}In_{0.45}Te$. The solid and dashed lines are the fitting results. Meissner state: a fully diamagnetic state that does not allow the magnetic field to penetrate inside; Vortex solid: forming the vortex lattice with a well pinning by defects (pinning centers); Vortex liquid: a state that transforms from melting the vortex lattice cannot easily be pinned by defects.

ering vortex core energy as given by [\[59\]](#page-7-0)

$$
\mu_0 H_{\rm cl}(T) = \frac{\Phi_0}{4\pi\lambda^2} (\ln \kappa + 0.5),\tag{9}
$$

and

$$
\frac{\mu_0 H_{c2}(T)}{\mu_0 H_{c1}(T)} = \frac{2\kappa^2}{\ln \kappa + 0.5}.
$$
 (10)

To obtain $\mu_0 H_{c2}$ at zero-temperature, we fitted the $\mu_0 H_{c2}(T)$ data with the modified GL formula [\[60\]](#page-7-0),

$$
\mu_0 H_{c2}(T) = \mu_0 H_{c2}(0) \left(\frac{1 - t^2}{1 + t^2}\right)^{n_{GL}},\tag{11}
$$

where n_{GL} is a constant. The best fitting yields $\mu_0 H_{c2}(0) =$ 1.73(1) T with $n_{GL} = 0.90(1)$. Combining with $H_{c1}(0) =$ 53(3) Oe, we obtained $\kappa = 25(1)$ and $\lambda(0) = 338(11)$ nm. The $\lambda(0)$ value is smaller than those of other reports, e.g., $\lambda(0) = 425$ nm and 578(2) nm for Sn_{0.55}In_{0.45}Te [\[13,14\]](#page-7-0). Moreover, by taking the initial slope of $d[\mu_0 H_{c2}(T)]/dT =$ $-0.438(2)$ T/K and normal state resistivity $\rho_0 = 139 \mu \Omega$ cm [\[61\]](#page-7-0), the electronic Sommerfeld coefficient and DOS are estimated to be $\gamma = 2.47(4)$ mJ/mol K² and $N(E_F) = 1.05(2)$ states/eV spin−¹ f.u. The theoretical DOS value of 0.56 states/eV spin⁻¹ f.u. can be seen in Fig. [1\(e\).](#page-2-0) According to the relation $N(E_F) = (1 + \lambda_{ep})N_b(E_F)$ with $N_b(E_F)$ is the band DOS, we can further obtain the electron-phonon coupling constant $\lambda_{ep} = 0.88(3)$. We note that previous reports of heat capacity studies of Sn1−*^x*In*x*Te by Haldolaarachchige *et al.* [\[62\]](#page-7-0) and Kobayashi *et al.* [\[7\]](#page-6-0) show that as increasing the Indoping content, the electronic Sommerfeld coefficient γ tends to increase. The γ value derived here is consistent with the results of the heat capacity measurements on the $Sn_{0.6}In_{0.4}Te$ [\[62\]](#page-7-0).

A narrow critical fluctuation region width is a good ingredient for the practical use of superconducting materials. Therefore, we evaluated its thermal fluctuations effect, which can be quantified by the Ginzburg parameter $G_i =$ $(2\pi k_B T_c \mu_0 \lambda_0^2 / \Phi_0^2 \xi)^2 / 2$, where μ_0 is the permeability of the vacuum, k_B is the Boltzmanns constant, and ξ is coherence length, respectively. Here, $\xi(0) = 13.8(1)$ nm is calculated using GL relation $\mu_0 H_{c2}(0) = \Phi_0/2\pi \xi^2(0)$. We obtained $G_i = 5.9(8) \times 10^{-7}$ for Sn_{0.55}In_{0.45}Te, which is five orders of magnitude smaller than that of high- T_c cuprates (10⁻²) [\[55\]](#page-7-0), but comparable to that of low- T_c materials (10⁻⁸) [\[55,63\]](#page-7-0). For more comparisons, its G_i value is smaller than other super-conductors like 2H-NbSe₂ (~10⁻⁴) [\[63\]](#page-7-0), CeRu₂ (~5×10⁻⁴) [\[64\]](#page-7-0), Ca₃Rh₄Sn₁₃ (∼3×10⁻⁷) [\[48\]](#page-7-0), and MgB₂ (∼10⁻⁶) [\[65\]](#page-8-0). Thus, it is expected that the small G_i value causes quite a narrow width of critical fluctuation region with Ginzburg temperature $|T_c - T| < T_c G_i = 3.0(4) \times 10^{-6}$ K. This is experimentally supported by the $\mu_0 H_{irr}(T)$ at low temperature but the critical fluctuation region broadening as it approaches the T_c , as guided by the dashed line in Fig. [7.](#page-5-0) At temperatures close to the T_c , the interval is about 0.09 K between $\mu_0 H_{irr}(T)$ and $\mu_0H_{c2}(T)$, supporting a very weak vortex fluctuation [\[53\]](#page-7-0). The $\text{Sn}_{0.55}\text{In}_{0.45}$ Te superconductor has a relatively narrow region between $\mu_0 H_{irr}(T)$ and $\mu_0 H_{c2}(T)$ compared to other low T_c superconductors, as reported in FeS_{1−*x*}Se_{*x*} [\[56\]](#page-7-0), Cu_{*x*}TaS₂ [\[66\]](#page-8-0), $SrPd₂Ge₂$ [\[67\]](#page-8-0) and thin film TiO [\[68\]](#page-8-0). Therefore, $\text{Sn}_{0.55}\text{In}_{0.45}\text{Te}$ is a promising candidate for superconducting devices applications below 5 K [\[17–20\]](#page-7-0). Furthermore, using the $\xi(0)$ and λ_0 values extracted before, the depairing current density [\[55\]](#page-7-0) (the maximum current that the Cooper pairs can carry before breaking up) can be further calculated by $J_0 =$ $\Phi_0/3\sqrt{3}\pi\mu_0\lambda^2\xi = 6.4(4) \text{ MA/cm}^2$. This results in a small ratio $J_c/J_0 \sim 4.62 \times 10^{-4} \ll 1$. In principle, this indicates that increasing J_c is further accessible by inducing artificial pinning centers for δT_c pinning.

Finally, we turn to discuss the potential applications as a SNSPD candidate for the $Sn_{1-x}In_xTe$ superconductors, which can be operated in liquid helium temperature. According to the photon detection model $[21]$, several physical parameters determine the smallest photon energy ϵ that is adequate for the single photon detection regime, $\epsilon = n\Delta(T)N(E_F)k_BT_cD d\tau_{th}$, where n is a factor for the energy lost due to the generation of subgap phonons, $\Delta(T)$ is the superconducting gap, *D* is the electron diffusivity, *d* is the thin film thickness, and τ_{th} is the electron thermalization time counting the time scale of the quasiparticles multiplication. Compared to the NbN $(T_c =$ 9 − 15 K) [\[69\]](#page-8-0), TaN (T_c = 8.16 K) [\[70\]](#page-8-0), NbSi (T_c = 2 K) [\[71\]](#page-8-0), α -W_{*x*}Si_{1−*x*} ($T_c \le 5$ K) [\[72\]](#page-8-0) superconductors for SNSPD applications, the availability of high-quality superconducting thin films with a broad tunable T_c value below 4.2 K, uniformity as indicated by the sharp superconducting transition, higher upper critical field in the two-dimensional limit and large normal state resistivity provide practical ingredients for developing a single-photon detector in wide wavelength range for Sn_{1−*x*}In_{*x*}Te materials [\[20\]](#page-7-0). On one hand, in the dirty and local limit [\[69\]](#page-8-0), the electron diffusivity is given by $D = 1.097(-d\mu_0H_{c2}/dT)^{-1}$ with *D* and $-(d\mu_0H_{c2})/dT$ in the units of cm² s⁻¹ and T K⁻¹. For the bulk Sn_{0.55}In_{0.45}Te crystal, the *D* is calculated to be 2.51(1) cm² s^{-1}. However, as a rough estimation, the *D* decreases to 1.2(1) cm² s^{-1} and $0.17(5)$ cm² s⁻¹ for out-of-plane and in-plane configuration in a 25 nm thick $\text{Sn}_{0.58}\text{In}_{0.42}$ Te thin film [\[20\]](#page-7-0), which are comparable to that of typical NbN material ($D \sim 0.5-0.7$ cm² s⁻¹) [\[69,73\]](#page-8-0). As a SNSPD candidate, the dimension-reduction provides a practical route to access smaller *D* to extend the cut-off wavelength for Sn_{1−*x*}In_{*x*}Te superconductors [\[21,22\]](#page-7-0). On the other hand, the quantum yield is given by $K = (1/n)(h\nu/\Delta)$ with v is photon frequency [\[21\]](#page-7-0). It indicates that the lower Δ would yield a higher *K* ratio and thus more Cooper pairs are to be broken per photon absorption, resulting in a higher quantum efficiency. Considering the small gap ratio determined here and the other reported for the $Sn_{0.55}In_{0.45}Te$ superconductor [\[13\]](#page-7-0), this provides a beneficial ingredient for it being as a SNSPD candidate in thin film form.

IV. CONCLUSIONS

In summary, we investigated the critical current density and vortex pinning dynamics in the superconductor $Sn_{0.55}In_{0.45}Te$. Our $H_{c1}(T)$ data supports that it has a weakly anisotropic single gap. The $J_c(T)$ curves show a nonuniversal power-law behavior with increasing field, which is ascribed to the coexistence of δT_c pinning and δl pinning regimes with field and temperature dependence. The high $J_c(0)$ and $U_0(H)$ values and very weak vortex fluctuation were demonstrated from the magnetization and electrical transport measurements. We show that the $Sn_{0.55}In_{0.45}Te$ superconductor has its potential applications in superconducting electronics like single-photon detectors in its thin film form.

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