Supercconductivity in noncentrosymmetric Ag$_2$Pd$_3$S

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We have successfully synthesized the single crystal of Ag$_2$Pd$_3$S, which exhibits superconductivity with the transition temperature of $T_c = 2.25$ K. Ag$_2$Pd$_3$S crystallizes in the space group $P4_132$ with the filled $\beta$-Mn structure, which has no inversion symmetry. The value of the Ginzburg-Landau parameter $\kappa_{GL}$ indicates that Ag$_2$Pd$_3$S is a type-II superconductor. $\Delta C(T_c)/\gamma T_c = 1.50$ and $2\Delta/k_B T_c = 3.48$ from the heat-capacity analyses indicate that Ag$_2$Pd$_3$S is a weak-coupling Bardeen-Cooper-Schrieffer (BCS) superconductor with an isotropic superconducting gap. On the other hand, the violation of the Werthamer-Helfand-Hohenberg curve in the $H$-$T$ phase diagram implies Ag$_2$Pd$_3$S is not a typical BCS superconductor.

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Noncentrosymmetric superconductors (NCSCs), whose crystal structures lack inversion symmetry, have recently attracted much attention due to their potential to realize unconventional superconductivity [1]. The Cooper-pair wave function of conventional superconductors generally consists of the product of orbital (i.e., $s$-, $p$-, or $d$-wave) and spin (singlet or triplet) parts. The combinations of these parts are strictly determined by the parity symmetry; if the orbital part is $s$ wave, the spin part is definitely singlet (even parity). However, this rule does not apply to NCSCs. The large antisymmetric spin-orbit coupling (ASOC) of NCSCs induces the mixing of orbital and spin parts, resulting in a mixture of the spin-singlet and spin-triplet states [2,3].

Triggered by the discovery of exotic superconductivity in CePt$_3$Si [4], extensive experimental and theoretical investigations on NCSCs have been carried out to understand their advantageous properties. CePt$_3$Si has an unusually high upper critical field $H_{c2} = 4 \sim 5$ T, which significantly exceeds the Pauli limit for singlet superconductivity ($H_{c2} \sim 1$ T). In addition, NCSCs are predicted to exhibit some other exotic phenomena, such as anisotropic superconducting gap structures and unique magnetoelectric effects that derive from their polar symmetry [5–10].

Recently, a number of compounds have been investigated as candidates for NCSCs [11–24], such as CePt$_3$Si [4] and CeRhSi$_3$ [11]. These two compounds have been extensively studied by many researchers, leading to the discovery of an unusual symmetry present in the Cooper pairs. CePt$_3$Si and CeRhSi$_3$ belong to the category of heavy fermion superconductors. In these superconductors, not only the broken inversion symmetry but also the quantum criticality affect the superconductivity of the compounds.

On the other hand, Li$_2$M$_3$B ($M = \text{Pd, Pt}$) [14–24] are two of the appropriate compounds for forming NCSCs to investigate effects of the broken spatial inversion symmetry only. Li$_2$M$_3$B ($M = \text{Pd, Pt}$) do not contain any magnetic elements; therefore, effects of magnetic ordering can be completely ignored. In this system, superconducting properties are directly dependent on the magnitude of ASOC. Li$_2$Pt$_3$B exhibits unconventional superconductivity with an anisotropic superconducting gap structure; $^{11}$B NMR experiments and magnetic penetration depth measurements revealed that the Cooper pair included about 60% of the spin-triplet pairing and 40% of the spin-singlet pairing, and the energy gap had a line node. In contrast, the gap structure of Li$_2$Pd$_3$B has been under debate. An isotropic full gap structure is suggested by various bulk measurements such as heat capacity [16,17], whereas Yuan et al. claimed by magnetic penetration depth measurements that a multigap structure consist of two gaps in which one is dominant and the other is small enough not to affect the obtained gap structure by bulk heat-capacity measurements [18]. On the other hand, Yuan et al. also pointed out by precise fitting of the temperature dependence of magnetic penetration depth below $T_c$ that the gap is slightly anisotropic and there existed two gaps which had almost the same size [20]. The magnitude of ASOC of Pt is somewhat larger than that of Pd, and this should cause the two compounds to demonstrate different types of superconductivity.

In this study, we focused on the ternary compound Ag$_2$Pd$_3$S. The superconductivity of Ag$_2$Pd$_3$S ($T_c = 1.13$ K) was first reported by Khan et al. in 1973 [25]. They synthesized a polycrystalline sample of Ag$_2$Pd$_3$S and found superconductivity by measuring an electrical resistivity. Thirty years later, Schaak et al. reported that Ag$_2$Pd$_3$S could be synthesized by the rapid low-temperature solution method [26]. However, details of the crystal structure have not been completely resolved due to the difficulty of structural analysis of polycrystalline samples, although the basic crystal structure of Ag$_2$Pd$_3$S was expected to be the $\beta$-Mn-type structure [25,26]. In addition, superconducting properties also remain poorly understood.

We have successfully synthesized single crystals of Ag$_2$Pd$_3$S. Detailed structural analyses were performed using the single-crystal samples, and it was revealed that Ag$_2$Pd$_3$S crystallized in a filled $\beta$-Mn-type structure with space group $P4_132$. This is isostructural to Li$_2$M$_3$B ($M = \text{Pd, Pt}$; space group $P4_132$), but these two structures have different types of crystallographic chirality: Ag$_2$Pd$_3$S has a right-handed structure, and Li$_2$M$_3$B has a left-handed structure. The superconductivity in Ag$_2$Pd$_3$S is of interest due to the similarity of its crystal structure with that of Li$_2$Pt$_3$B, since Li$_2$Pt$_3$B exhibits unconventional superconductivity in terms of the absence of the inversion symmetry. The studies of Ag$_2$Pd$_3$S,
TABLE I. Details of single-crystal x-ray diffraction data collection and crystal refinement results for Ag$_2$Pd$_3$S.

<table>
<thead>
<tr>
<th>Compound</th>
<th>Formula weight</th>
<th>Color, shape</th>
<th>Crystal dimension</th>
<th>Temperature</th>
<th>Space group</th>
<th>a</th>
<th>Volume</th>
<th>Z</th>
<th>D$_0$</th>
<th>F(000)</th>
<th>Absorption coefficient</th>
<th>Goodness of fit</th>
</tr>
</thead>
<tbody>
<tr>
<td>Ag$_2$Pd$_3$S</td>
<td>567.1 g mol$^{-1}$</td>
<td>Metallic luster, cuboid</td>
<td>0.1 $\times$ 0.08 $\times$ 0.06 mm$^3$</td>
<td>293 K</td>
<td>P4$_1$32</td>
<td>7.2292(2) Å</td>
<td>377.81(3) Å</td>
<td>4</td>
<td>9.968 g m$^{-3}$</td>
<td>992</td>
<td>24.491 mm$^{-1}$</td>
<td>1.31</td>
</tr>
</tbody>
</table>

An appropriate candidate compound for a NCSC, can be of significant importance for understanding physical properties of superconductors without inversion symmetry. In this paper, we investigate details of the crystal structure and superconducting properties of Ag$_2$Pd$_3$S and discuss similarities to other NSCSs, especially Li$_2$M$_6$B$_3$ where M = Pd, Pt.

The single-crystal samples were prepared by a self-flux method. The starting materials of Ag$_2$S and a fine powder of Pd metal were mixed in the molar ratio 1:3 and then pressed into a pellet. The pellet was sealed in an evacuated quartz tube and heated at 1000 °C for 48 h. After slowly cooling down to 500 °C at a rate of 5 °C/h, it was then quenched to room temperature. The obtained samples were in the form of a chunk of single crystals with a metallic luster. No interdomain impurity was confirmed by powder x-ray diffraction using pulverized crystals.

A single-crystal x-ray diffraction experiment at 293 K was carried out with a Bruker SMART APEX single-crystal diffractometer with a charge-coupled device area detector using Mo K$_\alpha$ radiation. The intensity data were integrated using SAINT and SAEDABS was employed to correct for absorption effects [27]. The structural parameters were refined using the program SHELXL-2013 based on the full-matrix method [28]. The final reliability indexes ($R$ values) were $R_{int}(F) = 1.81\%$ and $wR_{int}(F^2) = 3.24\%$. The details of the single-crystal x-ray diffraction measurement are summarized in Table I. The chemical composition of the crystal was confirmed to be almost stoichiometric using electron probe microanalysis.

The electrical resistivity was measured with a conventional dc four-probe method on a shaped lump of the single crystals in the temperature range 0.5–300 K using a Quantum Design (QD) Physical Properties Measurement System (PPMS). Heat-capacity measurements were carried out between 0.4 and 300 K using the heat-relaxation method on the lump of the single crystals in the QD PPMS. Magnetic susceptibility data were collected using a powder sample obtained from the pulverized crystal in the temperature range 1.8–300 K using a superconducting quantum interference device magnetometer (QD Magnetic Properties Measurement System).

First-principles calculations with spin-orbit coupling were carried out using the full-potential linearized augmented plane-wave (FLAPW) method, as implemented in the WIEN2K program package [29]. The Perdew-Burke-Ernzerhof generalized gradient approximation (GGA) potential was used for the exchange-correlation potential [30]. Experimentally obtained values for lattice constants and internal coordinates (see Table II) were used in our calculations. The linearized augmented plane-wave space radii of Ag$_2$Pd$_3$S were set to 2.50 a.u. for Ag, 2.28 a.u. for Pd, and 2.02 a.u. for S. The cutoff parameter $R_{MT}$ $K_{\text{max}}$ was chosen as 8.0, where $R_{MT}$ is the radius of the Ag sphere and $K_{\text{max}}$ is the magnitude of the largest reciprocal lattice vector. Self-consistency was imposed on $k$-point meshes of 94 points (12 $\times$ 12 $\times$ 12; we used approximately 1800 special $k$ points) in the irreducible part of the Brillouin zone.

The crystal structure of Ag$_2$Pd$_3$S is illustrated in Fig. 1(a). The details of structural parameters and the anisotropic atomic displacement parameters are shown in Tables II and III. Ag$_2$Pd$_3$S crystallizes in space group P4$_1$32, and the lattice constant is $a = 7.229 \AA$. The crystal structure is well described by the filled $\beta$-Mn structure. There are many intermetallic compounds that have a $\beta$-Mn structure, which originally possesses the $A_2B_3$ composition for the Mn metal, as shown in Fig. 1(b). In the $A_2B_3$-type structure, the 8c and 12d sites can be occupied by atoms, whereas the interstitial 4a site.

FIG. 1. (a) Crystal structure of Ag$_2$Pd$_3$S. Each shaded polyhedron represents a Pd$_6$S octahedron, in which sulfur is located at the center. The interstitial space in the Ag network is filled by corner-sharing Pd$_6$S octahedra. (b) Crystal structure of the $A_2B_3$ $\beta$-Mn structure. The solid line represents a cubic unit cell.
is basically vacant. Notably, the \( \beta \)-Mn structure exhibits a flexible composition due to an accommodating atom C at the 4a site, such as in \((A, B)_2C\), \((A, B)_2C\), or \((A, B)_2C\), including defects at the 8c and 12d sites to satisfy the Hume-Rothery rules.

The filled \( \beta \)-Mn structure is the crystal structure in which all 4a, 8c, and 12d sites are fully occupied. In many cases, small atoms, such as carbon, boron, and nitrogen, can only be located at the interstitial 4a site. To the best of our knowledge, \( Ag_2Pd_3S\) is the first compound that includes a sulfur atom at the 4a site. A feature of the structure is a three-dimensional network consisting of Ag atoms with corner-sharing Pd\(_6\)S octahedra that penetrate into the interspaces of the Ag network, as shown in Fig. 1. In the Ag network, all Ag atoms have three neighboring Ag atoms with the interatomic distance 2.74 Å, which is 0.15 Å shorter than that of Ag metal (2.89 Å). This suggests that an electron transfer occurs from Ag to the Pd/S framework.

The crystal structure of \( Ag_2Pd_3S\) is very similar to that of \( Li_2Pd_3B\). However, the space group of \( Ag_2Pd_3S\) is \( P4_132\), which is different from that of \( Li_2Pd_3B\) (\( P4_132\)). The corresponding crystallographic sites are as follows: S and B at 4a, Ag and Li at 8c, and Pd at the 12d site. The difference between these crystal structures lies in the helicity around the spiral axis. In \( Ag_2Pd_3S\), atoms have right-handed helicity, while in \( Li_2Pd_3B\), atoms are arranged with a left-handed helicity; in other words, the \( \beta \)-Mn structure has a degree of freedom of the structural chirality. It is known that the structural chirality sometimes affects the magnetic properties of localized spin systems [31,32]. On the other hand, it is unclear whether the chirality degree of freedom influences physical properties of metals, itinerant spin systems or not.

As shown in Fig. 2(a), the temperature dependence of the resistivity shows typical metallic behavior. The resistivity decreases with decreasing temperature, and the curve shows an upward convex feature at low temperatures. Zero resistivity was clearly observed below 2 K at zero magnetic field, indicating an occurrence of a superconducting transition. The transition temperature was determined from a zero and an onset of resistivity. The onset of resistivity is defined as a point where the linear extrapolation of the slope in the transition region intersects the linear extrapolation of the plateau region, as shown in the inset of Fig. 2(a). By applying magnetic fields, the transition is gradually broadened. Zero resistivity was not observed at or below 0.5 K under a magnetic field of \( \mu_0H = 1 \) T, as shown in Fig. 2(b), resulting in the residual resistivity \( \rho_0 = 6.21 \times 10^{-5} \Omega \text{ cm} \) and the residual resistivity ratio, \( RRR = 5 \) at \( T = 0.55 \) K.

Figure 3(a) shows the temperature dependence of the magnetic susceptibility under a magnetic field of 10 Oe. Single crystals were well ground to obtain powder samples, and measurements were performed on the polycrystalline samples. As shown in Fig. 3(a), the magnetic susceptibility shows diamagnetism below 2.2 K. This is evidence for superconductivity. The Meissner volume fraction estimated at 1.8 K under the field-cooled process is about 40%, indicating that the superconductivity is a bulk phenomenon.

The superconducting \( M-H \) curve measured at 1.8 K is shown in Fig. 3(b). The magnetic shielding effect clearly occurs below about 900 Oe at 1.8 K. The \( M-H \) curve strongly indicates that \( Ag_2Pd_3S\) is a type-II superconductor, which is a typical superconducting behavior with weak vortex pinning. The inset shows a magnified version of the low-magnetic-field region. The lower critical field \( H_c1 \), at which the magnetization

<table>
<thead>
<tr>
<th>Atom</th>
<th>( U_{11} )</th>
<th>( U_{22} )</th>
<th>( U_{33} )</th>
<th>( U_{12} )</th>
<th>( U_{13} )</th>
<th>( U_{23} )</th>
</tr>
</thead>
<tbody>
<tr>
<td>Ag</td>
<td>0.01438(9)</td>
<td>0.01438(9)</td>
<td>0.01438(9)</td>
<td>0.01438(9)</td>
<td>0.00043(8)</td>
<td>0.00043(8)</td>
</tr>
<tr>
<td>Pd</td>
<td>0.01288(8)</td>
<td>0.01071(14)</td>
<td>0.01397(10)</td>
<td>0.00329(10)</td>
<td>0.00178(7)</td>
<td>-0.0004(3)</td>
</tr>
<tr>
<td>S</td>
<td>0.0101(3)</td>
<td>0.0101(3)</td>
<td>0.0101(3)</td>
<td>0.0101(3)</td>
<td>-0.0004(3)</td>
<td>-0.0004(3)</td>
</tr>
</tbody>
</table>

\[
\begin{align*}
\beta & = 0 \\
\rho & = 6.21 \times 10^{-5} \Omega \text{ cm} \\
RRR & = 5 \text{ at } T = 0.55 \text{ K}
\end{align*}
\]
deviates from a linear response, was 3.5 Oe at 1.8 K, indicating that vortices can penetrate the compound at very low magnetic fields.

The temperature dependence of total heat capacity (electronic part and lattice part) $C_{\text{total}}/T$ vs $T^2$ under $\mu_0 H = 0$ and 1 T is depicted in Fig. 4(a). A peak observed at zero magnetic field was completely suppressed at $\mu_0 H = 1$ T. The Sommerfeld coefficient $\gamma$ and the phononic specific heat coefficient $\beta$ were estimated as $\gamma = 8.55 \text{ mJ/(mol K}^2)$ and $\beta = 0.96 \text{ mJ/(mol K}^4)$. These values were obtained by a linear extrapolation to the normal heat capacity $C_{\text{total}}$ under $\mu_0 H = 1$ T, assuming $C_{\text{total}} = \gamma T + \beta T^3$ equation. The Debye temperature $\Theta_D = 230$ K was also determined from the relation $\beta = (12/5)\pi^4 N_A N_A k_B / \Theta_D^3$. Here, $N_A$ is the Avogadro constant, $N_A$ is the number of atoms per formula unit, and $k_B$ is the Boltzmann constant. We obtain the electron-phonon coupling constant $\lambda_{\text{el-ph}} = 0.48$ from McMillan’s formula $T_c = (\Theta_D/1.45)\exp[-1.04(1 + \lambda_{\text{el-ph}}/\mu^*(1 + 0.62\lambda_{\text{el-ph}}))]$ [33] by substituting $\Theta_D = 230$ K and the Coulomb pseudopotential $\mu^* = 0.13$ [34]. The small value of $\lambda_{\text{el-ph}}$ suggests that a weak-coupling superconductivity is realized in this compound.

The temperature dependence of the electronic specific heat divided by temperature, $C_{\text{el}}/T = C_{\text{total}}/T − \beta T^2$, is illustrated in Fig. 4(b). We estimated the transition temperature $T_c = 2.25$ K and the jump of heat capacity $\Delta C(T_c)/T_c = 10.52$ at $T_c$ to satisfy the entropy conservation around $T_c$. Figure 4(c) shows the temperature dependence of $C_{\text{el}}$ at zero magnetic field. The thermodynamic critical field $\mu_0 H_c(0)$ estimated from the equation $\mu_0 H_c(0)/2 = −\gamma T_c^2/2 + \int_0^{T_c} C_{\text{el}}(T) dT$ is 0.018 T. We analyzed $C_{\text{el}}$ below $T_c$ to reveal the nature of a superconducting gap. The solid line shows the fitting result assuming the specific heat expression for a conventional superconductor with temperature-linear term $C_{\text{el}} = \gamma T + a\exp[−(\Delta(0)/k_B T)]$, where $\Delta(0)$ is the energy gap. The estimated gap value is $\Delta(0)/k_B = 3.91$ K, and thus $2\Delta/k_B T_c$ is calculated to be 3.48. This is close to the value of 3.53 that is expected from the BCS theory. We obtained a finite value of $\gamma = 1.53 \text{ mJ/(mol K}^2)$ at $T = 0$ K, suggesting that part of the domains in the Ag$_2$Pd$_3$S crystal does not participate in the superconductivity. It is probably due to internal defects in the Ag$_2$Pd$_3$S crystal, such as local inhomogeneity deviated slightly from the stoichiometric composition. Accordingly, the Sommerfeld coefficient of normal state intrinsic to the superconducting fraction should be $\gamma_n = \gamma − \gamma_0 = 7.02 \text{ mJ/(mol K}^2)$. Thus, the experimental value $\Delta C(T_c)/\gamma_n T_c = 1.50$ is consistent with that of the BCS weak-coupling limit value 1.43.

Figure 4(d) shows the temperature dependence of $C_{\text{el}}/T$ under several magnetic fields. Transition temperatures under each field were determined by assuming entropy balance around $T_c$. Applying magnetic fields causes each peak to shift to a lower temperature. Finally, the peaks vanish above $\mu_0 H = 0.5$ T, indicating that the superconducting transition is suppressed under those magnetic fields. Note that the onset of the superconductivity was still observed in the temperature dependence of the resistivity at $\mu_0 H = 0.5$ T.

Figure 5 shows the total and partial density of states (DOS) of Ag$_2$Pd$_3$S including the spin-orbit coupling calculated by the FLAPW method. As shown in Fig. 5, we found that Ag$_2$Pd$_3$S possesses a finite DOS at the Fermi level, which is mainly composed of the Pd 3$d$ orbital, while the contributions from Ag and S are small. The DOS structure we obtained is similar to that of Li$_2$Pd$_3$B calculated by Lee et al. [35]. The calculated DOS at $E_F$ per formula unit (f.u.) is 9.51 states/eV f.u., and the Sommerfeld coefficient $\gamma_{\text{el}}$ evaluated from the DOS is 5.61 mJ/(mol K$^2$). This value is slightly smaller than that obtained experimentally: $\gamma_n = 7.02 \text{ mJ/(mol K}^2)$. The mass enhancement may be attributed to the electron-phonon and electron-electron interactions.

Figure 6 summarizes the temperature dependence of the upper critical field $H_{c2}$ in Ag$_2$Pd$_3$S deduced from resistivity and heat-capacity measurements. $T_c$ values under several magnetic fields are estimated from both onset and zero resistivity, and from the $C_{\text{el}}/T$ vs $T$ plot. $T_c$ estimated from the heat capacity nearly corresponds to that from the onset of resistivity. Thus, we estimate the upper critical field using the plots of the onset of resistivity. The Werthamer-Helfand-Hohenberg (WHH) model for the $H_{c2}$ in the dirty
limit, $\mu_0H_c^{\text{WHH}}(0) = -0.693 T_c |d(\mu_0H_c)/dT|_{T_c} = 0.43$ T, is used to fit the $H$-$T$ curve. However, we were unable to reproduce the experimental results, especially in low-temperature and high-magnetic-field regions. Thus, we determined an approximated value of the upper critical field $\mu_0H_c(0) \sim 0.63$ T from a linear extrapolation at around $T_c$. According to the BCS theory, the $H_c(0)$ value is related to the coherence length $\xi$ by $H_c(0) = \phi_0/2\pi\xi(0)^2$, where $\phi_0$ is the magnetic flux quantum. Using this formula, we obtain the value $\xi(0) = 23$ nm. The Ginzburg-Landau parameter $\kappa_{GL}$ and the penetration depth $\lambda(0)$ are estimated from the relations $\mu_0H_c(0) = \sqrt{2}\kappa_{GL}\mu_0H_\lambda(0)$ and $\kappa_{GL} = \lambda(0)/\xi(0)$. We obtain the values $\kappa_{GL} = 25$ and $\lambda(0) = 568$ nm, respectively. These parameters from the experiments are summarized in Table IV.

The value $T_c = 2.25$ K for Ag$_2$Pd$_3$S is about two times larger than values previously reported [25,26]. Since we were unable to prepare samples that have a lower transition temperature in this research, it is hard to understand experimentally why the $T_c$ values are different from each other. One possible explanation is an influence of the sulfur deficiency of the previous studies, which may be caused by the large surface area due to the polycrystalline samples. The DOS structure of Ag$_2$Pd$_3$S shown in Fig. 5 may shed light on the difference in the $T_c$ values of those two samples. The DOS of Ag$_2$Pd$_3$S obtained from the band calculation exhibits steep peaks around $E_F$. If the system obeys a rigid band scheme, a small amount of sulfur deficiency decreases the carrier electron content and leads to the reduction of the DOS at $E_F$ ($E_F$ shifts to a lower energy). In the BCS theory, $T_c$ directly relates to the DOS at $E_F$; therefore, this reduction of DOS may significantly affect the magnitude of $T_c$. On the other hand, we used single-crystalline samples in our experiment. The chemical analysis and the single-crystalline x-ray diffraction measurement ensure that there is negligible sulfur deficiency in our samples. This may cause the larger DOS at $E_F$ compared to those of previous samples, resulting in our higher $T_c$.

The results of bulk measurements of Ag$_2$Pd$_3$S suggest that superconducting properties of the compound are described by the conventional BCS model; thermodynamic analyses of $\Delta C(T_c)/\gamma_n T_c = 1.50$ and $\Delta/k_B T_c = 3.48$ correspond to the usual single full-gap picture. Comparing with related compounds Li$_2$M$_2$B($M$ = Pd,Pt), the superconductivity of Ag$_2$Pd$_3$S is close to that of Li$_2$Pd$_3$B. Previous reports reveal...
FIG. 5. Calculated DOS of Ag2Pd3S with spin-orbit interaction. The total and partial DOS for each atom is illustrated. The broken line at $E = 0$ indicates the Fermi energy.

that admixture of the spin-singlet and the spin-triplet states with a line node in a superconducting gap is realized in Li2Pt3B, while the detailed gap structure of Li2Pd3B is unsolved; the possibilities of an isotropic single gap and a two-gap structure has been under discussion \[18,20\]. The observed $T$ dependence rather looks like a $T$-linear behavior (heat-capacity data) or an upward tendency (resistivity onset-observed $T_c$ data) even at $0 K \sim 0.3 T_c$. At a glance, it resembles the $h^*$ curve for strongly coupled superconductors \[39\]; however, it is unlikely because the experimental results of specific heat jump at $T_{C1} [\Delta C(T_c)/\gamma_c T_c \sim 1.50]$ and electron-phonon constant ($\lambda_{el-ph} \sim 0.48$) clearly suggests weak electron-phonon coupling in Ag2Pd3S. Werthamer Helfand, and Hohenberg expand their theory of $h^*$ by taking a Pauli spin paramagnetic pair-breaking effect with a Maki parameter $\alpha$ and a spin-orbit scattering frequency $\lambda_{SO}$ as perturbation into the orbital pair-breaking effect \[38\]. Their improved theory could successfully give more realistic $H_{c2}-T$ curve having a more plateau-like shape under high magnetic fields for some alloy superconductors; however, a curve thus obtained cannot reproduce such a straight data-set line observed in Ag2Pd3S even if $\alpha$ and $\lambda_{SO}$ are adjusted to modest values. It can be expected that for Ag2Pd3S, orbital pair breaking essentially limits the actual $H_{c2}$, because the experimental $H_{c2}(0)$ value ($\sim 0.63 T$) is much lower than the Pauli limit ($\mu_B H_{c2}(0) = 1.86 \times T_c \sim 4.2 T$), suggesting a small Maki parameter. The paramagnetic spin correction does not seem so much applicable to the case of Ag2Pd3S. A more realistic model in describing such a deviation of $H_{c2}(T)$ from the WHH theory, e.g., Fermi surface topology, gap anisotropy, and/or multigap structure, might be key factors to understand superconducting properties moderately, which superficially makes the system behave as a normal BCS superconductor like Li2Pt3B.

Contrarily, the observed $H_{c2}$ vs $T$ relationship (Fig. 6) does not seem to obey the BCS theory with simplified model assumptions. Although the theoretical calculation of the orbital limit of the critical field by Helfand and Werthamer \[37\] should give a convex curvature of $T$-dependent reduced critical field $h^*(T/T_C)$ around low temperatures ($T \ll T_c$), the actually observed $T$ dependence rather looks like a $T$-linear behavior (heat-capacity data) or an upward tendency (resistivity onset) even at 0.7 K ($\sim 0.3 T_c$). At a glance, it resembles the $h^*$ curve for strongly coupled superconductors \[39\]; however, it is unlikely because the experimental results of specific heat jump at $T_{C1} [\Delta C(T_c)/\gamma_c T_c \sim 1.50]$ and electron-phonon constant ($\lambda_{el-ph} \sim 0.48$) clearly suggests weak electron-phonon coupling in Ag2Pd3S. Werthamer Helfand, and Hohenberg expand their theory of $h^*$ by taking a Pauli spin paramagnetic pair-breaking effect with a Maki parameter $\alpha$ and a spin-orbit scattering frequency $\lambda_{SO}$ as perturbation into the orbital pair-breaking effect \[38\]. Their improved theory could successfully give more realistic $H_{c2}-T$ curve having a more plateau-like shape under high magnetic fields for some alloy superconductors; however, a curve thus obtained cannot reproduce such a straight data-set line observed in Ag2Pd3S even if $\alpha$ and $\lambda_{SO}$ are adjusted to modest values. It can be expected that for Ag2Pd3S, orbital pair breaking essentially limits the actual $H_{c2}$, because the experimental $H_{c2}(0)$ value ($\sim 0.63 T$) is much lower than the Pauli limit ($\mu_B H_{c2}(0) = 1.86 \times T_c \sim 4.2 T$), suggesting a small Maki parameter. The paramagnetic spin correction does not seem so much applicable to the case of Ag2Pd3S. A more realistic model in describing $h^*$ may be required for understanding the $T$-linear behavior of $H_{c2}(T)$ in Ag2Pd3S. Presumably, additional factors unconsidered in the WHH theory, e.g., Fermi surface topology, gap anisotropy, and/or multigap structure, might be key factors to understand such a deviation of $H_{c2}(T)$ from the WHH theory. Kita and Arai have reported that $h^*$ is enhanced significantly over the curve for the spherical Fermi surface (assumed in the WHH theory) with marked upward curvature \[40,41\]. This might be the most plausible case for Ag2Pd3S. At present, there is no definitive answer for the question of the unsat-

FIG. 6. $H-T$ phase diagram of Ag2Pd3S deduced from the temperature dependence of the resistivity and heat capacity. The solid and open circles indicate the $T_c$ values from onset and zero resistivity, respectively. The solid triangles show the $T_c$ values from the heat capacity. The solid and dashed lines are respectively the linear fit and the WHH fit to the data around the low-magnetic-field region of the diagram.

| $T_c$ | 2.25 K |
| $\gamma$ | 8.55 mJ/(mol K$^2$) |
| $\gamma_n$ | 7.02 mJ/(mol K$^2$) |
| $\gamma_0$ | 1.53 mJ/(mol K$^2$) |
| $\Theta_n$ | 230 K |
| $\mu_B H_{c2}(0)$ | 0.63 T |
| $\mu_B H_c(0)$ | 0.018 T |
| $\kappa_{GL}$ | 25 |
| $\xi(0)$ | 23 nm |
| $\lambda(0)$ | 568 nm |
| $\Delta C(T_c)/\gamma_c T_c$ | 1.50 |
| $2\Delta/k_B T_c$ | 3.48 |
| $\lambda_{el-ph}$ | 0.48 |
uration of $H_{c2}$. Further experimental and theoretical studies on Fermi surface shape and Maki parameters are expected to be helpful for deeper understanding of the critical field in Ag$_2$Pd$_3$S.

To clarify the detailed superconducting properties of Ag$_2$Pd$_3$S, microscopic experiments such as photoemission spectroscopy, NMR, or magnetic penetration depth measurements should be carried out. A variety of superconductors without inversion symmetry is desired in order to reveal the detailed physical properties of NCSCs. In particular, the fact that Ag$_2$Pt$_3$S, which is strongly related to the exotic superconductor Li$_2$Pt$_3$B, has a much larger ASOC than that of Ag$_2$Pd$_3$S must be a clue to help understand the unconventional properties of the NCSCs; actually, the preparation of the compound has been under way for some time. The theoretical and experimental studies on various NCSCs will reveal the novel and interesting physical properties of superconductors without inversion symmetry.

We successfully synthesized single crystals of the non-centrosymmetric superconductor Ag$_2$Pd$_3$S with a filled $\beta$-Mn structure and revealed the superconducting properties in terms of resistivity, magnetic susceptibility, heat capacity, and band calculations. Ag$_2$Pd$_3$S exhibits the superconducting transition at $T_c = 2.25$ K. The obtained superconducting parameters summarized in Table IV indicate that a type-II conventional weak-coupling BCS superconductivity with an isotropic gap is realized in Ag$_2$Pd$_3$S. On the other hand, temperature and field dependence of the $T_c$ indicates that the superconductivity of Ag$_2$Pd$_3$S deviates from the normal BCS picture. Although an origin of the non-WHH shape in the $H$-$T$ curve is still unclear, the understanding of superconducting properties of Ag$_2$Pd$_3$S plays an important role to study the weak ASOC effect on the compounds without inversion symmetry.

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