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Superconductivity of MoBe₂₂ and WBe₂₂ at ambient- and under applied-pressure conditions

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MoBe₂₂ and WBe₂₂ compounds belong to the binary $X\text{Be}_{22}$ ($X = 4d$ or $5d$ metal) family of superconductors, whose critical temperature depends strongly on X . Despite the multiphase nature of these samples, it is possible to investigate the superconducting properties of MoBe₂₂ and WBe₂₂ at the macro- and microscopic level. A concurrent analysis by means of magnetization and heat-capacity measurements, as well as muon-spin spectroscopy (μSR) was implemented. At ambient pressure, both compounds enter the superconducting state below 2.6 ± 0.1 K (MoBe₂₂) and 4.1 ± 0.10 K (WBe₂₂) and show modest upper critical fields [$(\mu_0 H_{c2}(0) = 48 \pm 1$ mT and $\mu_0 H_{c2}(0) = 58 \pm 1$ mT, respectively]. In WBe₂₂, the temperature-dependent superfluid density suggests a fully gapped superconducting state, well-described by an s -wave model with a single energy gap. Heat-capacity data confirm that such a model applies to both compounds. Finally, ac magnetic susceptibility measurements under applied pressures up to 2.1 GPa reveal a linear suppression of the superconducting temperature, typical of conventional superconducting compounds.

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I. INTRODUCTION

As one of the lightest elements, beryllium exhibits lattice vibrations of high frequency, a necessary condition for achieving superconductivity with a sizable critical temperature. Nevertheless, the critical temperature of elemental beryllium is only 0.026 K [1]. The value of T_c is also affected by the electron-phonon coupling strength (typically large in elements with covalent-bonding tendencies) and the density of states at the Fermi level $N(E_F)$ (rather low in elemental Be). In particular, $N(E_F)$ depends on the details of crystal structure and atomic volume—two factors which are best illustrated by metal-hydride superconductors under pressure [2]. Here, recent work has demonstrated that, in actinium hydrides at 200 GPa, purely phonon-mediated superconductivity with T_c up to 250 K [3] is likely to occur because these materials contain metal atoms that are close to populating a new electronic subshell—for example, d^1 (Sc, Y, La, and Ac) or p^0 (Be, Mg, and Ca). In this case, the electronic structure becomes highly

sensitive to the positions of the neighboring atoms, resulting in stronger electron-phonon interactions and higher $N(E_F)$. Based on this, Be-rich alloys may achieve a T_c much higher than that of elemental Be, a prediction which is correct for ReBe₂₂ ($T_c \sim 9.6$ K [4,5]). A detailed examination of ReBe₂₂ revealed it to be a multigap type-II superconductor, well-described by an isotropic s -wave model. Here, the enhanced T_c (compared to elemental Be) seems to arise from a *concomitant* increase of both electron-phonon coupling strength and $N(E_F)$ [5]. By investigating the effects of hydrostatic pressure on ReBe₂₂, a 0.5 K decrease of T_c with a pressure of 1 GPa was established. Further increase in pressure does not change T_c , as confirmed by electrical resistivity measurements at even higher pressures—up to 30 GPa [6]. Another possibility to control the superconducting properties of the $X\text{Be}_{22}$ class is to employ the isotope effect and/or the chemical pressure, whose effects are significant in this class of compounds [4]. It is, therefore, of interest to examine the isostructural MoBe₂₂ and WBe₂₂ compounds (with their structure shown in Fig. 1), both under pressure as well as at ambient conditions.

Superconductivity of MoBe₂₂ and WBe₂₂ has been previously reported to occur below $T_c = 2.52$ K and $T_c = 4.14$ K, respectively [4,7,8]. However, given the difficulties in synthesizing and characterizing beryllium-containing materials [9–15], neither system has been investigated thoroughly. Additionally, it has been shown that a combined analysis of specific-heat and μSR data in similar superconducting materials is highly relevant [9,16–18].

Motivated by this, we present an in-depth study of MoBe₂₂ and WBe₂₂ by means of magnetization, ac -susceptibility, heat-capacity, and muon-spin spectroscopy (μSR) measure-

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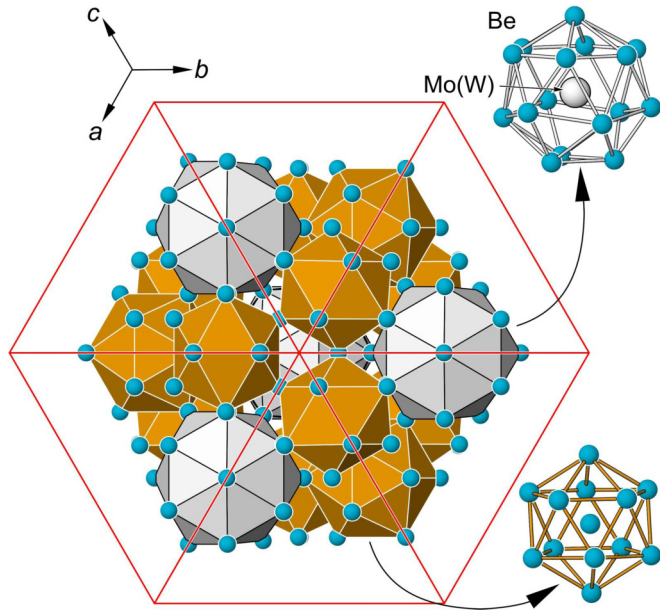


FIG. 1. The arrangement of the coordination polyhedra around Mo(W) and Be₂ atoms (16d site) in the structures of MoBe₂₂ and WBe₂₂. For simplicity, only part of the unit cell is shown.

ments. In addition, we examine the effects of pressure on the superconductivity of both materials. Despite the multiphase nature of the samples, we could establish that both compounds are conventional, fully-gapped type-II superconductors, with a modest electronic specific-heat coefficient, suggestive of a negligible electron-mass enhancement. The low values of the upper critical field and the linear decrease of T_c with pressure are typical of this class of superconductors.

II. EXPERIMENTAL DETAILS

All sample preparation and handling was performed in a specialized laboratory, equipped with an argon-filled glove box system [MBraun, $p(\text{H}_2\text{O}/\text{O}_2) < 0.1$ ppm] [19]. The polycrystalline samples were synthesized by arc melting of Be (Heraeus, >99.9 wt%) with Mo or W (Chempur, 99.99%) in the 99:1 ratio. Since the amount of Be is hard to quantify analytically and given that evaporation losses of Be are unavoidable, the only way to keep the sample composition under control is by following a careful weighing protocol [12–14]. None of the samples exhibit any marked air- or moisture sensitivity. Both MoBe₂₂ and WBe₂₂ phases melt incongruently and have a high melting temperature (1300 °C and 1520 °C, respectively [20,21]). To date, this has prevented the synthesis of single-phase $X\text{Be}_{22}$ materials, let alone of their single crystals. In the current paper, we could successfully synthesize and characterize multiphase polycrystalline samples of both MoBe₂₂ and WBe₂₂, albeit with small amounts of foreign phases (see below).

Given the high mechanical hardness and ductility of both samples, it was not possible to perform reliable powder diffraction experiments. Instead, a metallographic analysis of the polished sample surfaces was performed *via* energy-dispersive x-ray spectroscopy on a Jeol JSM 6610 scanning electron microscope, equipped with an UltraDry

EDS detector (ThermoFisher NSS7). The microstructures of the Mo-Be and W-Be samples show very similar arrangement of the phases. In both cases, the material-contrast images recorded with a backscattered electron detector show bright dendrites embedded in a matrix. The matrix is formed by a mixture of dark grey fibers within a dark, almost globular phase [see Figs. 2(a), 2(b), 3(a), and 3(b) for Mo-Be and W-Be, respectively]. The high transition-metal concentration of the bright dendritic and the fiber-shaped phases, as indicated by their brightness in the BSE images, is confirmed by the qualitative comparison of the EDX spectra. The local EDX spectra are extracted from equal areas of the EDX mapping data and show decreasing intensities of the transition-metal x-ray lines, e.g., Mo L line for the Mo-Be samples and W L and M lines for the W-Be samples for the darker phases. SEM analyses were performed with acceleration voltages of 10 kV or 20 kV for the Mo-Be and W-Be samples, respectively. MoBe₂₂ and WBe₂₂ can be assigned to the dark phase of the respective microstructure due to the significant intensities of the transition-metal x-ray lines. Additionally, solid Be does not show any significant solubility in Mo or W. Based on the published binary phase diagrams of Mo-Be and W-Be [20,21], it is reasonable to assume that the congruent melting phases MoBe₂ and WBe₂ form the respective dendritic phase, while MoBe₁₂ and WBe₁₂ form the fiber-shaped phases with an intermediate transition-metal concentration. We have also established that the impurity phases Mo/WBe₂ and Mo/WBe₁₂ do not show any magnetic transition down to $T = 1.8$ K. Based on Figs. 2 and 3, we estimate that the samples from the current study contain $\sim 40\%$ of the Mo/WBe₂₂, $\sim 30\%$ of the Mo/WBe₁₂, and $\sim 30\%$ of the Mo/WBe₂ phases.

Since the previously reported structures of MoBe₂₂ and WBe₂₂ were obtained from powder diffraction data [22,23], we performed single-crystal x-ray diffraction measurements. For this, a single-phase cuboid, consisting predominantly of either WBe₂₂ or MoBe₂₂ material, was fabricated in a multistep process using a Xe-plasma focused-ion-beam (FEI Helios G4 PFIB). With a size of $\sim 30 \times 30 \times 30 \mu\text{m}^3$, it was ideally suited for single-crystal diffraction experiments [see Figs. 2(c) and 3(c)]. The relevant diffraction data were collected using a Rigaku AFC7 diffractometer, equipped with a Saturn 724+ CCD detector and Mo $K\alpha$ radiation source ($\lambda = 0.71073 \text{ \AA}$). SHELXL software was used for the structural refinement. The results of the crystallographic characterization of MoBe₂₂ and WBe₂₂ are provided in Tables S1 and S2, respectively [24].

The magnetic properties of both compounds were studied using a Quantum Design (QD) Magnetic Property Measurement System in the temperature range from 1.8 K to 300 K, in applied magnetic fields up to 7 T. The specific-heat data were collected on a QD Physical Property Measurement System in the temperature range from 0.4 K to 10 K, in magnetic fields up to 9 T. Both magnetization and specific-heat measurements were performed on bulk samples, showing minimal variation of superconducting features between the different sample batches or pieces. Given the multiphase nature of our samples, it was not possible to carry out reliable electrical resistivity measurements.

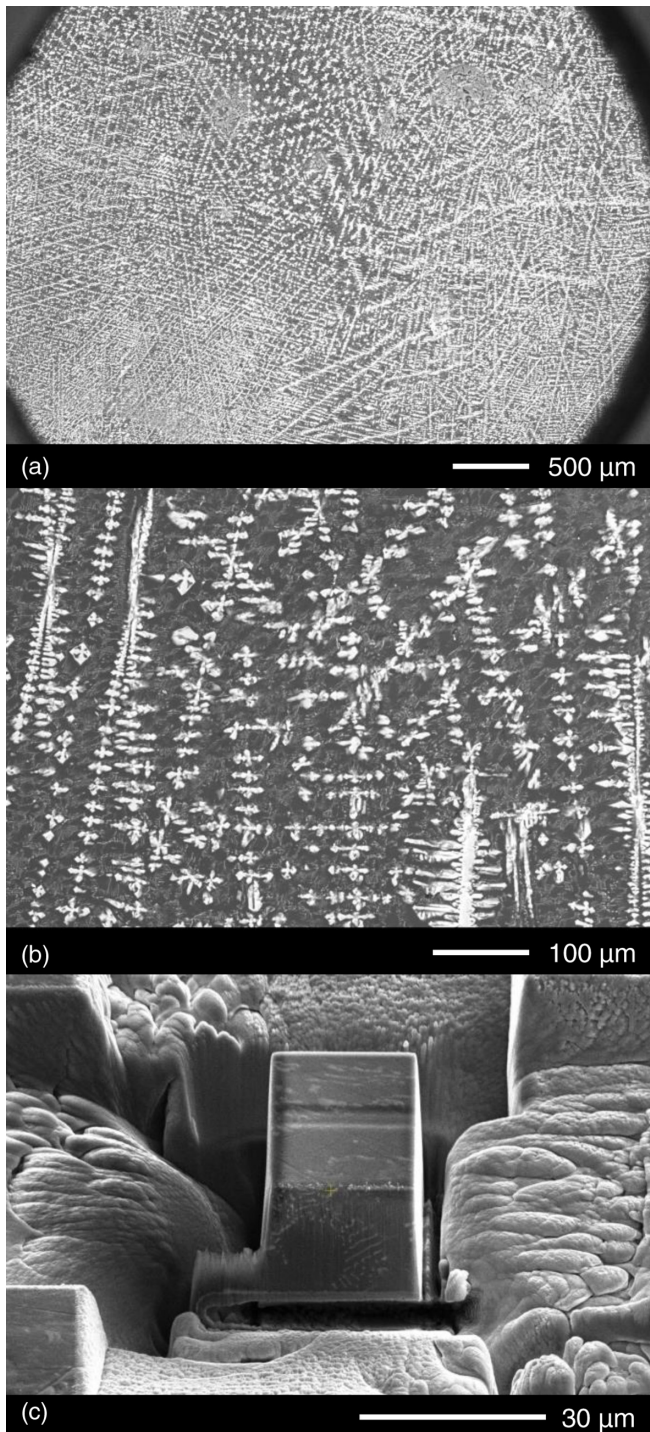


FIG. 2. (a), (b) Backscatter electron micrographs of MoBe₂₂ sample surface at various magnifications. (c) A cuboid of MoBe₂₂, used for the single-crystal diffraction experiments, prepared using focused-ion beam.

The muon-spin spectroscopy measurements were performed at the Dolly (zero pressure) and GPD (under pressure) spectrometers of the Swiss muon source at Paul Scherrer Institut, Villigen, Switzerland. The μ SR data were analyzed by means of the MUSRFIT software package [25]. In the zero-pressure case, three discs of ca. 8 mm in diameter were

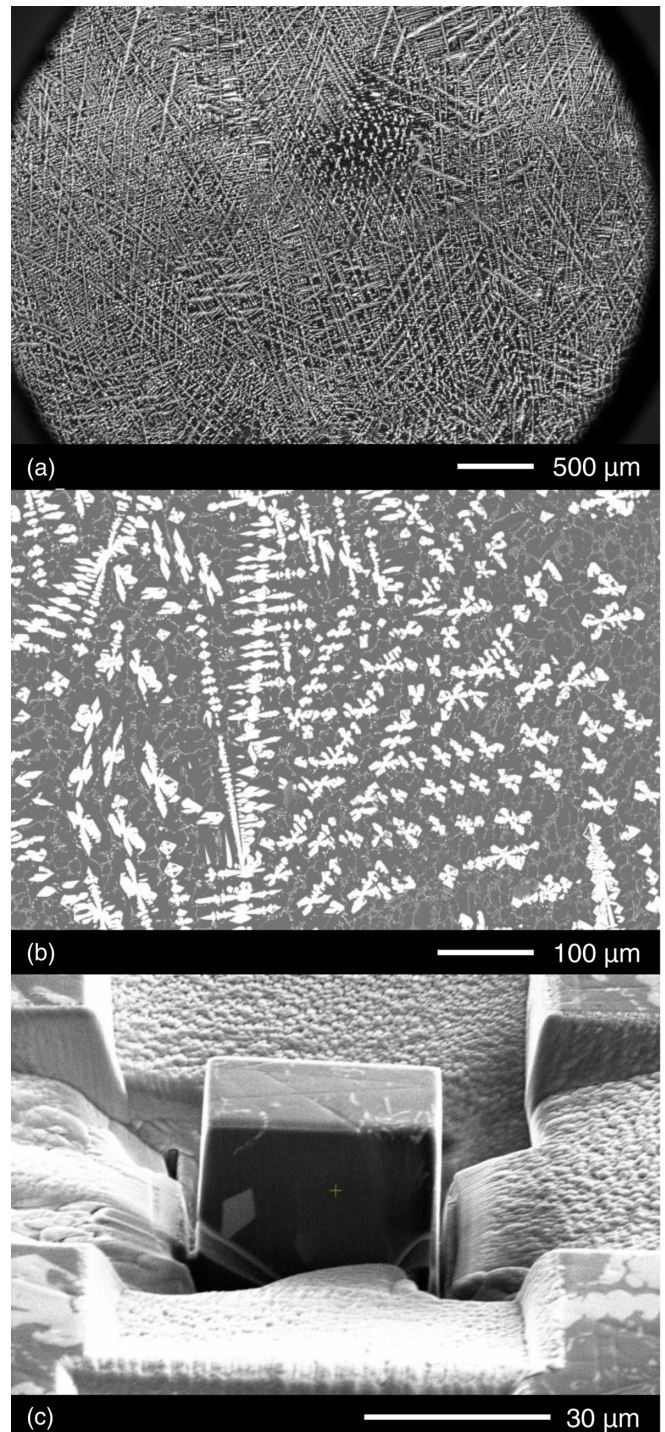


FIG. 3. (a), (b) Backscatter electron micrographs of WBe₂₂ sample surface at various magnifications. (c) A cuboid of WBe₂₂, used for the single-crystal diffraction experiments, prepared using focused-ion beam.

positioned so as to cover more than 85% of the muon-beam cross section, thus resulting in low-background high-quality μ SR signals. The applied field of 15 mT yielded a T_c of 3.5 K for WBe₂₂, significantly lower than $T_c(0) = 4.09$ K, hinting at a rather low H_{c2} value—see below. The same sample was successively used to fill a dual-walled WC pressure cell

(channel diameter 6 mm) up to a height of 12 mm. The dual-wall configuration allows the cell to reach pressures of up to 2.3 GPa, albeit at the expense of a relatively low sample-to-cell μ SR signal ratio. We used Daphne 7373 oil as a pressure transmitting medium, whose properties under a broad range of p - T conditions are well-known [26,27].

ac susceptibility measurements under pressure were performed using the same pressure cell used for the μ SR measurements [28,29]. In our case, the excitation and pick-up coils were wound outside the cell. While this setup generally has a low sensitivity, it is sufficient for successfully detecting the superconducting transition [30]. An Ametek 7270 DSP lock-in amplifier was used to capture the 72-Hz ac signal, successively measured by a Keithley 2000 digital multimeter.

III. RESULTS AND DISCUSSION

A. Crystal structure description

In previous studies [22,23], the structure of MoBe₂₂ and WBe₂₂ was established by comparing the experimental powder diffraction patterns (photographic data) with calculated patterns, obtained on the basis of the Mg₃Cr₂Al₁₈ [31] structure (E phase). The atomic coordinates were taken from the latter model and, therefore, were not refined. For this reason, a reinvestigation of the crystal structure of MoBe₂₂ and WBe₂₂ was performed to obtain a more precise crystallographic description.

Both structures are isotypic with ZrZn₂₂ [32]. The substitution variants were also described for CeCr₂Al₂₀ [33] and Mg₃Cr₂Al₁₈ [31]. The crystal structure of MoBe₂₂ (WBe₂₂) can be represented as an arrangement of Frank-Kasper polyhedra with 16 and 12 apexes (icosahedra) formed by Be species around the Mo and Be₂ positions, respectively. The polyhedra are only connected to each other through their vertices. Each [MoBe₁₆] polyhedron has in its neighborhood four identical groups in tetrahedral arrangement and 12 [BeBe₁₂] icosahedra (see Fig. 1). The described polyhedra resemble the atomic coordination of Mg and Cu, respectively, in the Laves phase MgCu₂ [34].

B. Magnetization and specific heat

The superconductivity of MoBe₂₂ and WBe₂₂ was previously reported to occur below $T_c = 2.52$ K and 4.14 K [4,7,8], respectively. In both materials, the first indication of superconductivity is given by a sharp diamagnetic transition at the respective T_c value, as shown in Figs. 4(a) and 4(b) [35]. For $\mu_0 H > 2$ mT (MoBe₂₂) and $\mu_0 H > 10$ mT (WBe₂₂), the low-temperature plateau is not visible, due to it being below the minimum accessible temperature of 1.8 K. Normally, the lower critical field value H_{c1} is defined as the field at which the $M(H)$ curve deviates from the line having the initial slope of $M(H)$. However, in our case, the lowest experimentally accessible temperature was not sufficient to allow for such an analysis. In addition, in the MoBe₂₂ sample, the origin of the kink around 2.4 K is not yet clear. Nonetheless, since the shielding fraction is above 100% and none of the identified impurities are known to be superconductors, the superconductivity of both MoBe₂₂ and WBe₂₂ is intrinsic.

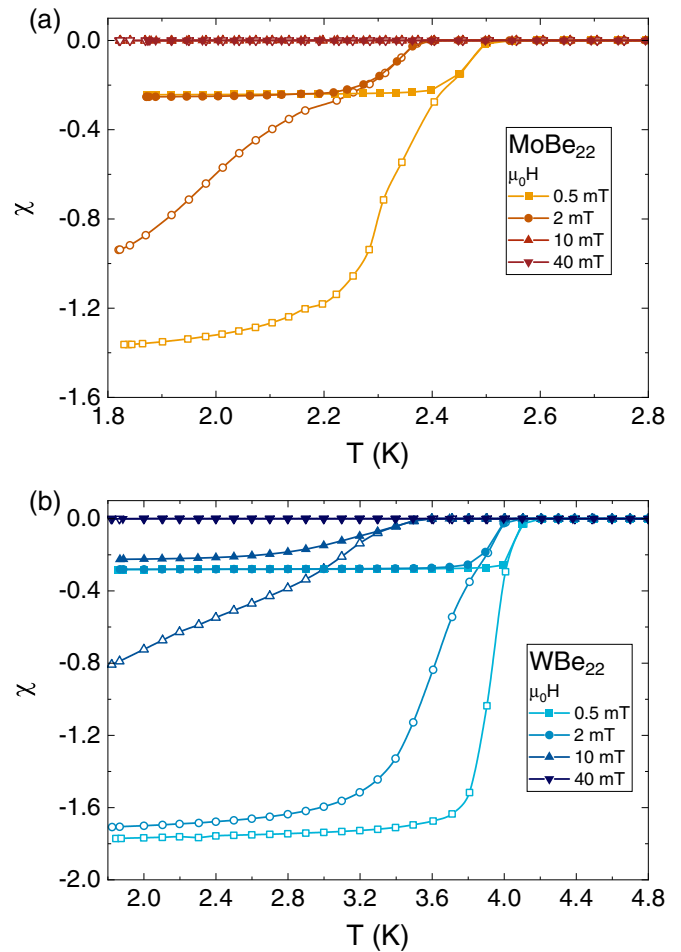


FIG. 4. Zero-field-cooled (open symbols) and field-cooled (full symbols) temperature-dependent magnetic susceptibility data for MoBe₂₂ (a) and WBe₂₂ (b) in an applied field $0.5 \leq \mu_0 H \leq 40$ mT.

The specific-heat data of MoBe₂₂ and WBe₂₂, shown in Fig. 5, also confirm their bulk superconductivity. In the absence of a magnetic field, an anomaly associated with the onset of the superconducting transition is observed at $T_c = 2.55$ K and $T_c = 4.09$ K in MoBe₂₂ and WBe₂₂, respectively. A linear fit to C_p/T versus T^2 data [insets of Figs. 5(b) and 5(d)] provides an upper-bound estimate for the values of the Sommerfeld coefficient $\gamma_n = 6.8 \pm 0.2$ mJ mol⁻¹ K⁻² and $\gamma_n = 10.0 \pm 0.2$ mJ mol⁻¹ K⁻² for MoBe₂₂ and WBe₂₂, respectively. The residual electronic specific-heat coefficient $\gamma_{\text{res}} = C_e/T$ ($T = 0.35$ K, $\mu_0 H = 0$) ≈ 4 mJ mol⁻¹ K⁻² for both MoBe₂₂ and WBe₂₂ samples is consistent with the presence of the secondary phases identified in Figs. 2 and 3. The multiphase nature of the samples is also likely responsible for the large difference in the β coefficients: $\beta = 0.21 \pm 0.05$ mJ mol⁻¹ K⁻⁴ (MoBe₂₂) and $\beta = 0.05 \pm 0.01$ mJ mol⁻¹ K⁻⁴ (WBe₂₂). These β values can, nevertheless, be used for a preliminary estimate of the Debye temperature (600 K $\leq \theta_D \leq 900$ K), as well as of the electron-phonon coupling strength ($\lambda_{e\text{-ph}} \sim 0.4$). The latter value classifies both MoBe₂₂ and WBe₂₂ as weakly-coupled superconductors.

Further evidence of conventional superconductivity in MoBe₂₂ and WBe₂₂ is given by the value of the jump in

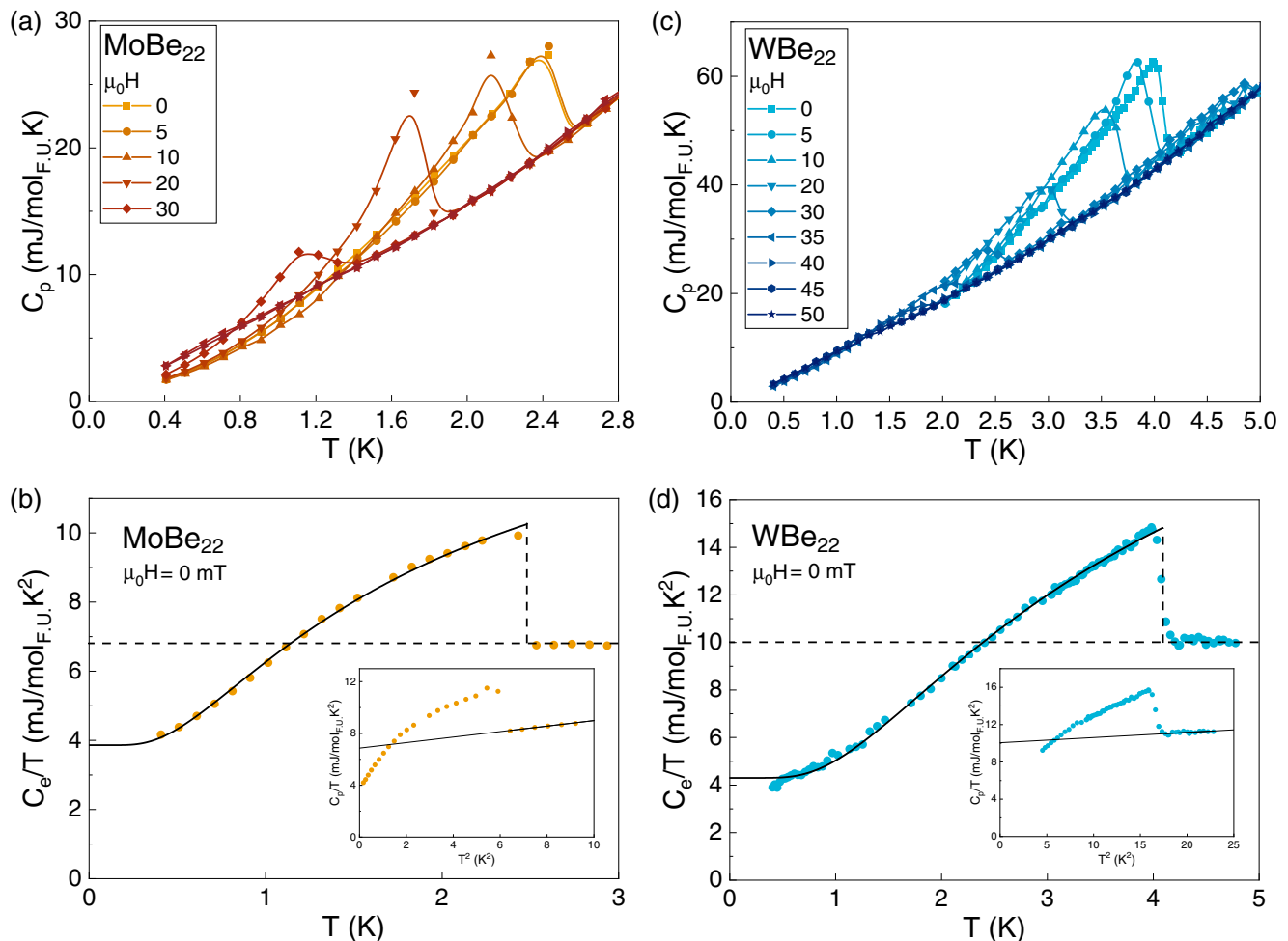


FIG. 5. Total specific heat of MoBe₂₂ (a) and WBe₂₂ (c) as a function of temperature in an applied field $0 \leq \mu_0 H \leq 50$ mT. The electronic specific heat of MoBe₂₂ (b) and WBe₂₂ (d) measured in $\mu_0 H = 0$ mT. Vertical dashed lines represent equal-entropy constructions. The horizontal dashed line corresponds to the Sommerfeld coefficient γ_n . The solid line is a fit to $C_e/T \propto e^{-\Delta/T}$ in the superconducting state. Insets of panels (b) and (d) show C_p/T versus T^2 , from which the values of γ and β were extracted.

the electronic specific heat, $\Delta C_e/\gamma_n T_c \approx 1.39$ for MoBe₂₂ [Fig. 5(b)] and $\Delta C_e/\gamma_n T_c \approx 1.21$ for WBe₂₂ [Fig. 5(d)], both of which are comparable with the standard value of the Bardeen-Cooper-Schrieffer (BCS) theory ($\Delta C_e/\gamma_n T_c = 1.44$). In the superconducting state, the specific-heat data are well fit by the $C_e/T \propto e^{-\Delta/k_B T}$ expression deduced from the BCS theory, yielding $2\Delta(0)/k_B T_c = 3.28$ and $2\Delta(0)/k_B T_c = 3.53$ for MoBe₂₂ and WBe₂₂, respectively. In both cases, the good agreement between the measured data (symbols) and the BCS fit (lines) provides compelling evidence for an *s*-wave isotropic BCS superconducting gap in the electronic density of states, occurring exactly at the Fermi level. The values of $2\Delta(0)/k_B T_c$ are comparable to the weak-coupling value of 3.52, expected from the BCS theory.

Upon applying a magnetic field, the superconducting transitions of both MoBe₂₂ and WBe₂₂ are gradually suppressed, as summarized in Figs. 5(a) and 5(c). As shown in Fig. 6, the corresponding values of T_c are given in the $H - T$ phase diagram, shown Fig. 6 in red (MoBe₂₂) and blue (WBe₂₂). The fit to the Ginzburg-Landau relation $H_{c2}(T) = H_{c2}(0) \frac{1-(T/T_c)^2}{1+(T/T_c)^2}$ (dashed lines) yields $T_c = 2.7$ K and $T_c = 4.3$ K, as well as

$\mu_0 H_{c2}(0) = 48$ mT and $\mu_0 H_{c2}(0) = 58$ mT. Interestingly, the values of $H_{c2}(0)$ are a factor of two larger than those reported previously for MoBe₂₂ and WBe₂₂ [7]. The discrepant $H_{c2}(0)$ values are most likely related to the different level of impurities [36].

C. *ac* susceptibility under applied pressure

In most solid-state systems, the primary effect of a change in temperature is to modify the *occupation* of the energy levels, while the primary effect of applied pressure is to modify the *energy levels* themselves [37]. Clearly, this makes a pressure study very interesting, as it can provide complementary information to temperature. To characterize the behavior of MoBe₂₂ and WBe₂₂ under applied pressure, we measured the *ac* response in pressures of up to 2.1 GPa. A small piece of indium, embedded in the pressure cell, served as a pressure gauge. Indeed, since its $T_c(p)$ dependence is well-known, we can determine p (in GPa) from the measured T_c (in K) of indium [38]:

$$T_c(p) = T_c(0) - 0.3812 p + 0.0122 p^2. \quad (1)$$

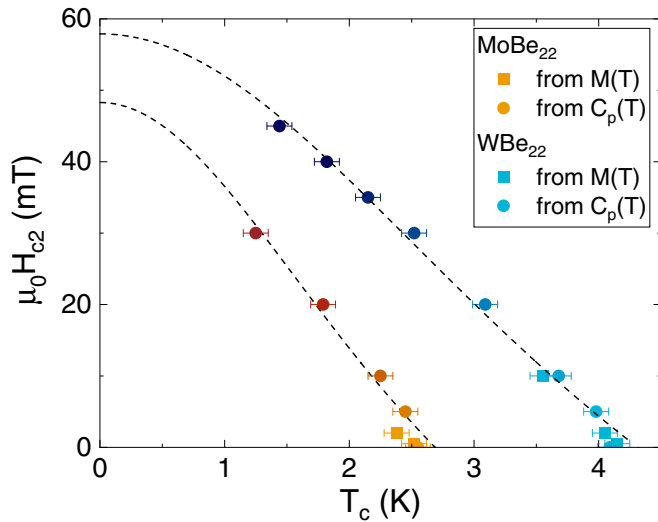


FIG. 6. H - T phase diagram of MoBe_{22} (orange) and WBe_{22} (blue). The values of the critical field $H_{c2}(T)$ are extracted from $M(T)$ (squares) and C_p (circles) data. The dashed lines correspond to a Ginzburg-Landau fit of the data.

Here, $T_c(0) = 3.4$ K, while the coefficients have uncertainties of 0.0002 and 0.0004, respectively. In view of our relatively narrow pressure range (< 2.1 GPa) and due to a 30-fold larger linear coefficient compared to the quadratic one, T_c depends almost linearly on p . Even at the highest pressure, the omission of the quadratic term implies an error smaller than 2%. Similar considerations apply to the investigated samples, as confirmed by the linear behavior of T_c versus p , shown in Fig. 7.

To detect the superconducting transition, we use the real part χ' of the ac susceptibility (Fig. 7, main panels). Here, T_c is defined as the midpoint of the superconducting transition. In both cases, as the pressure increases from 0 to 2.1 GPa, T_c decreases by *ca.* 0.4 K. The raw ac susceptibility data indicate that besides the superconducting transition of In, we

TABLE I. Values of the superconducting temperatures T_c and their pressure dependence for MoBe_{22} and WBe_{22} , compared to In (here used as a pressure gauge).

Compound	T_c [K]	dT_c/dp [mK/GPa]	dt/dp [10^{-3} GPa $^{-1}$]
WBe_{22}	3.975	-192.6	-48.5
MoBe_{22}	2.422	-136.2	-56.2
In	3.403	-354.8	-104.3

also observe a second transition with a lower T_c [see Fig. 7(b)]. Considering its closeness to the T_c of indium and their similar pressure slopes, we attribute the anomaly to a spurious phase of indium (and/or to a mixture of the χ'' and χ' terms of the ac response).

The T_c values under pressure, as determined from ac susceptibility data and shown in the insets of Fig. 7, were fit with a line. The obtained parameters are reported in Table I, with In used as a pressure gauge. Since In is softer than both XBe_{22} compounds, it exhibits a steeper dT_c/dp slope. We note also that, although the absolute slopes of XBe_{22} are rather different [reflecting their different $T_c(0)$ values], the relative slopes, here expressed in terms of the normalized temperature $t = T_c(p)/T_c(0)$, exhibit more regular values. Indeed, in these relative units, one obtains a very similar behavior under pressure for both WBe_{22} and MoBe_{22} , as illustrated in the last column of Table I. This result may be justified by the similar Van der Waals and ionic radii of Mo and Be [39].

Recently, the isostructural ReBe_{22} compound was investigated under applied pressures of up to 30 GPa. The value of its T_c was also shown to decrease linearly with pressure [6]. However, despite the similar structure, its relative slope is a factor of 7 lower than that of WBe_{22} . The reason for such a discrepancy is not yet clear, but it might reflect the different atomic radius of Re, compared to those of Mo and W.

Finally, we note that, in general, the linear dependence of T_c on pressure indicates the absence of a pressure-induced structural transition and indirectly confirms the hydrostaticity

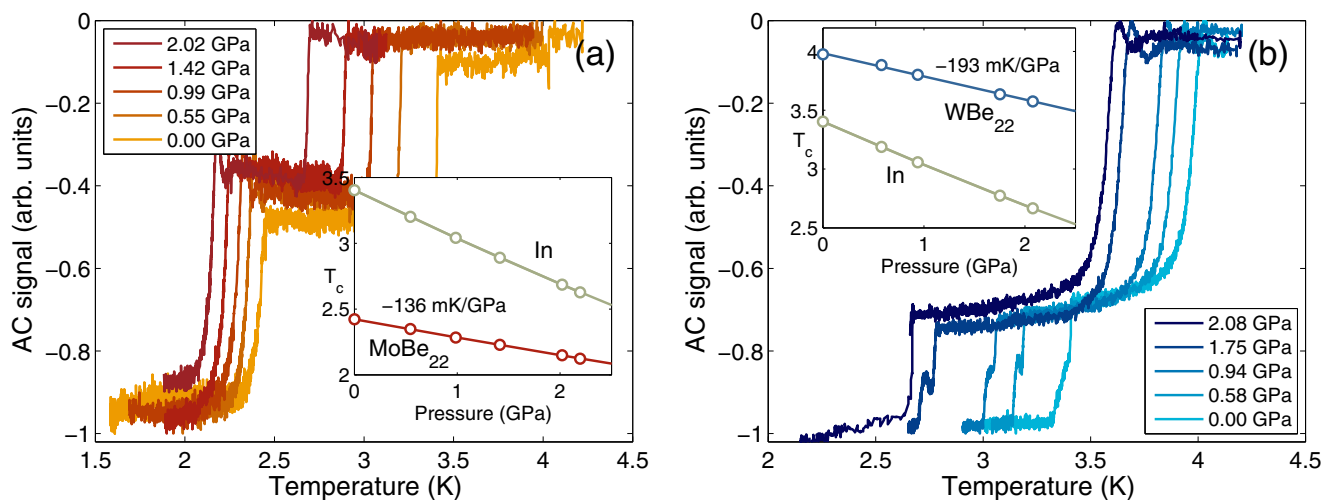


FIG. 7. ac susceptibility data versus temperature, measured at various applied pressures for MoBe_{22} (a) and WBe_{22} (b). In both cases, the steep changes in signal reflect the response of the sample and of indium (used here as a pressure gauge). Insets show the almost linear dependence of T_c on pressure. In both cases, error bars are smaller than the symbols.

of the pressure medium. Normally, the critical temperature of a superconductor depends on both the lattice- and electronic properties. Under very general assumptions (weak coupling, BCS superconductors), one can write [see Eq. (2.9) in Ref. [40]:

$$T_c \sim \sqrt{\frac{k}{M}} e^{\frac{-k}{\eta}}, \quad (2)$$

where k is the spring constant (a lattice term), η is a purely electronic term, and M is the atomic mass. Clearly, dT_c/dp depends on the relative magnitude of the pressure-induced changes in lattice versus electronic properties. Under applied pressure, the lattice stiffens, leading to an increase in k . Additionally, η also increases under pressure, although typically much less than k . Since k inside the exponent outweighs k in the prefactor, in general, an increase in k leads to a decrease in T_c . Since in the majority of conventional superconductors the pressure-induced lattice stiffening dominates over the modest changes in electronic properties, this implies a mostly linear decrease of T_c under pressure.

D. Muon-spin rotation study

As a microscopic technique, μ SR is very suitable for investigating the nature of pairing symmetry in the superconducting phase. In this case, one applies a moderate- to low-magnetic field, transverse to the muon-spin direction (TF- μ SR), and follows the evolution of the muon polarization decay as a function of temperature. Besides the (usually small) depolarization due to the nuclear magnetic moments, below T_c , one observes an increase in the decay rate due to the field modulation from the flux-line lattice (FLL) in the superconducting phase [41]. Such modulation is considerable for fields slightly above H_{c1} and it disappears just below H_{c2} . For optimal results (i.e., to achieve a relatively large decay rate with only a moderate suppression of T_c) one typically applies a field twice H_{c1} , usually above 20 mT. However, in our case, the H_{c2} value of both MoBe₂₂ and WBe₂₂ is quite low (only about 50 mT, see Fig. 6). This makes it rather difficult to study these compounds *via* TF- μ SR, in particular, MoBe₂₂, whose T_c and muon depolarization rate are low. Measurements under applied pressure are even more challenging. This is mostly due to the low fraction of signal coming from the sample (typically 2/3 being due to the pressure cell) and the similarity of the relaxation rates of muons stopped in the pressure cell with those stopped in the sample. For these reasons, the remainder of our paper focuses only on WBe₂₂.

To track the additional field-distribution broadening due to the FLL in the mixed superconducting state, the magnetic field was applied in the normal state, prior to cooling the sample below T_c . After the field-cooling protocol, the TF- μ SR measurements were performed at various temperatures upon warming. Figure 8(a) shows two representative TF- μ SR spectra collected in the superconducting (0.3 K) and the normal state (5 K) in an applied field of 40 mT. Additional TF- μ SR spectra collected at 15 mT exhibit similar features (not shown). The inhomogeneous field distribution due to the FLL is reflected in an enhanced depolarization rate below T_c . In this case, the TF- μ SR spectra can be modeled by the following

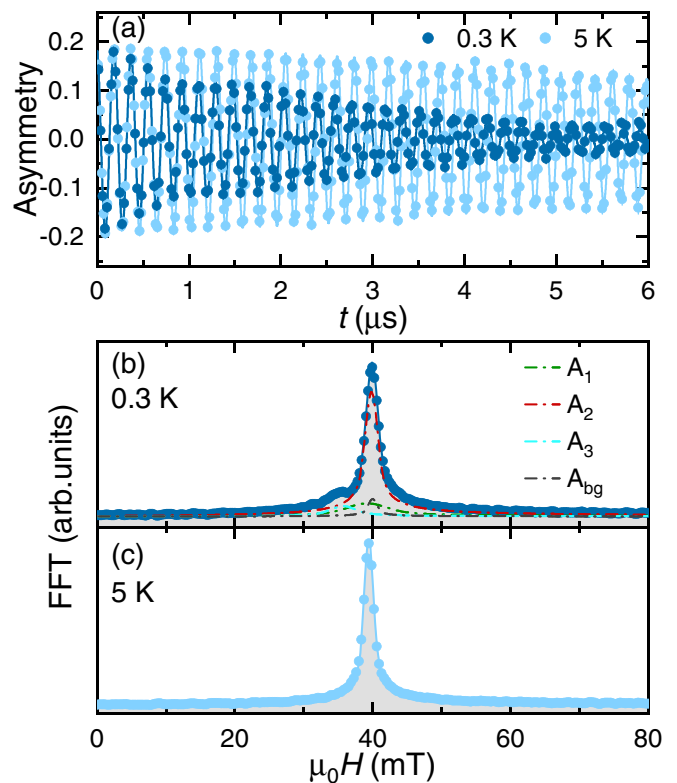


FIG. 8. (a) TF- μ SR time-domain spectra of WBe₂₂ collected at $T = 0.3$ K and $T = 5$ K in an applied field of 40 mT. The respective Fourier transforms are shown in (b) and (c). Below T_c , three oscillations are needed to reproduce the data. Solid lines are fits to Eq. (8) using three oscillations. Note the absence of a diamagnetic shift below T_c in panel (b)

expression:

$$A_{\text{TF}}(t) = \sum_{i=1}^n A_i \cos(\gamma_\mu B_i t + \phi) e^{-\sigma_i^2 t^2 / 2} + A_{\text{bg}} \cos(\gamma_\mu B_{\text{bg}} t + \phi). \quad (3)$$

Here, A_i and A_{bg} represent the initial muon-spin asymmetries for muons implanted in the sample and sample holder, respectively, with the latter not undergoing any depolarization. B_i and B_{bg} are the local fields sensed by implanted muons in the sample and sample holder, $\gamma_\mu = 2\pi \times 135.53$ MHz/T is the muon gyromagnetic ratio, ϕ is a shared initial phase, and σ_i is a Gaussian relaxation rate of the i th component. The number of required components is material dependent [42]. We find that, similarly to the ReBe₂₂ superconductor [5], $n = 3$ is sufficient to describe the TF- μ SR spectra of WBe₂₂. The fast-Fourier-transform (FFT) spectra of the TF- μ SR datasets at $T = 0.3$ K and $T = 5$ K are shown in Figs. 8(b) and 8(c). The solid lines represent fits to Eq. (3) using three oscillations i.e., $n = 3$ in the superconducting state and one oscillation in the normal state. The derived Gaussian relaxation rates as a function of temperature are summarized in the inset of Fig. 9.

Above T_c , the relaxation rate caused by nuclear magnetic moments is small and temperature independent, but below T_c it starts to increase due to the onset of the FLL and the increase of superfluid density. At the same time, a diamagnetic field

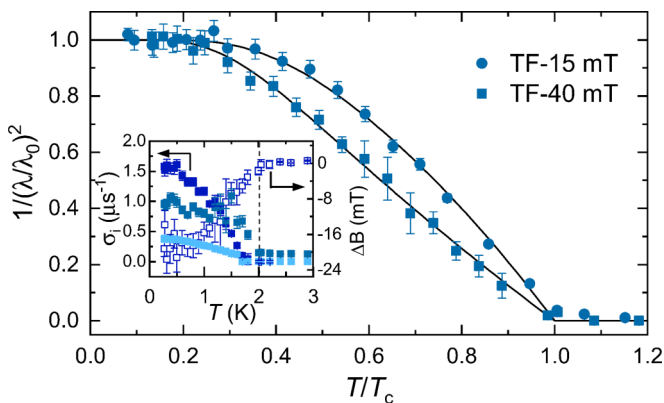


FIG. 9. Superfluid density of WBe₂₂ versus temperature, as determined from TF- μ SR measurements at $\mu_0 H = 15$ and 40 mT. The inset shows the temperature dependence of the three components of the μ SR rate $\sigma_i(T)$ at 40 mT (left axis) and the diamagnetic field shift (right axis). Lines represent fits to a fully gapped s -wave model with a single superconducting gap (see text for details).

shift appears below T_c (see inset of Fig. 9). In the case of multicomponent oscillations, the first-term in Eq. (3) describes the field distribution as the sum of n Gaussian relaxations [42]:

$$P(B) = \gamma_\mu \sum_{i=1}^n \frac{A_i}{\sigma_i} \exp\left[-\frac{\gamma_\mu^2 (B - B_i)^2}{2\sigma_i^2}\right]. \quad (4)$$

Then, the first and second moments of the field distribution can be calculated by

$$\langle B \rangle = \sum_{i=1}^n \frac{A_i B_i}{A_{\text{tot}}} \quad \text{and} \quad (5)$$

$$\langle B^2 \rangle = \frac{\sigma_{\text{eff}}^2}{\gamma_\mu^2} = \sum_{i=1}^n \frac{A_i}{A_{\text{tot}}} \left[\frac{\sigma_i^2}{\gamma_\mu^2} - (B_i - \langle B \rangle)^2 \right], \quad (6)$$

where $A_{\text{tot}} = \sum_{i=1}^n A_i$. The superconducting Gaussian relaxation rate related to the FLL (σ_{sc}) can be extracted by subtracting the nuclear contribution according to $\sigma_{\text{sc}} = \sqrt{\sigma_{\text{eff}}^2 - \sigma_n^2}$, where σ_n is the nuclear relaxation rate.

Since σ_{sc} is directly related to the magnetic penetration depth (and thus to the superfluid density), the superconducting gap value and its symmetry can be investigated by measuring the temperature-dependent $\sigma_{\text{sc}}(T)$. Note that the upper critical field of WBe₂₂ is relatively small (58 mT) compared to the fields used in the TF- μ SR study (15 and 40 mT). Hence, the effects of the overlapping vortex cores with increasing field should be considered when extracting the magnetic penetration depth λ from σ_{FLL} . For WBe₂₂, the effective magnetic penetration depth λ was calculated by means of [41,43]

$$\sigma_{\text{FLL}} = 0.172 \frac{\gamma_\mu \Phi_0}{2\pi} (1-h) [1 + 1.21(1-\sqrt{h})^3] \lambda^{-2}, \quad (7)$$

where $h = H_{\text{appl}}/H_{c2}$ and H_{appl} is the applied magnetic field. The calculated inverse-square of the magnetic penetration depth is proportional to the superfluid density $\lambda^{-2}(T) \propto \rho_{\text{sc}}(T)$ and is shown in Fig. 9 for WBe₂₂. Below $T_c/3$, the two superfluid densities $\rho_{\text{sc}}(T)$, determined *via* TF-15 and TF-40 mT μ SR measurements, are practically independent of temperature. This is in agreement with the specific-heat data,

shown in Fig. 5, indicating a fully gapped superconducting state. Therefore, we applied an s -wave model to analyze the $\rho_{\text{sc}}(T)$ data:

$$\rho_{\text{sc}}(T) = 1 + 2 \int_{\Delta(T)}^{\infty} \frac{E}{\sqrt{E^2 - \Delta^2(T)}} \frac{\partial f}{\partial E} dE, \quad (8)$$

Here, $f = (1 + e^{E/k_B T})^{-1}$ is the Fermi function, and the temperature dependence of the gap is assumed to follow $\Delta(T) = \Delta_0 \tanh\{1.82[1.018(T_c/T - 1)]^{0.51}\}$ [44,45], where Δ_0 is the gap value at zero temperature (here assumed as the only adjustable parameter). As shown by solid lines in Fig. 9, the temperature-dependent $\lambda^{-2}(T)$ is consistent with an s -wave model with a single gap for both applied fields. For TF-15 mT, the zero-temperature magnetic penetration depth is $\lambda_0 = 168(2)$ nm and the estimated gap value is $\Delta_0 = 1.80(5) k_B T_c$. The latter is fully compatible with the standard BCS prediction ($\Delta_0 = 1.76 k_B T_c$), confirming the weakly coupled superconductivity. For TF-40 mT, we find $\lambda_0 = 105(3)$ nm and $\Delta_0 = 1.35(5) k_B T_c$. In this case, due to the low H_{c2} , even an applied field of 40 mT suppresses the superconducting gap significantly.

Finally, although high-pressure μ SR measurements did not provide quantitative results, we can still draw qualitative conclusions. The evolution of the μ SR relaxation rate as a function of temperature suggests a suppression of T_c with applied pressure. In addition, within the experimental error, we could not detect a change in the nature of superconductivity—the s -wave pairing is preserved up to at least $p = 2.1$ GPa. Finally, the lack of significant changes in relaxation down to the lowest temperature is compatible with the lack of a possible pressure-induced magnetic order.

IV. CONCLUSIONS

By combining results of magnetization, specific-heat, and μ SR experiments, we investigated the superconductivity of MoBe₂₂ and WBe₂₂. Both systems exhibit a weakly-coupled BCS-like superconducting state below $T_c = 2.7$ K (MoBe₂₂) and $T_c = 4.3$ K (WBe₂₂). The modest values of the electronic specific-heat coefficient ($\gamma_n = 6.8$ mJ mol⁻¹ K⁻² and 10.0 mJ mol⁻¹ K⁻²), as well as the low upper critical field values [$\mu_0 H_{c2}(0) = 48$ mT and 58 mT] are all compatible with a standard BCS-like behavior. In both cases, high-pressure ac -susceptibility and μ SR results show a linear decrease of T_c with pressure with a similar slope ($dt/dp \sim -50 \times 10^{-3}$ GPa⁻¹) and no change in the nature of superconducting pairing.

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