Direct evidence for a pressure-induced nodal superconducting gap in the Ba$_{0.65}$Rb$_{0.35}$Fe$_2$As$_2$ superconductor

Z. Guguchia$^1$, A. Amato$^1$, J. Kang$^2$, H. Luetkens$^1$, P.K. Biswas$^1$, G. Prando$^3$, F. von Rohr$^4$, Z. Bukowski$^5$, A. Shengelaya$^6$, H. Keller$^4$, E. Morenzoni$^1$, Rafael M. Fernandes$^2$ & R. Khasanov$^1$

The superconducting gap structure in iron-based high-temperature superconductors (Fe-HTSs) is non-universal. In contrast to other unconventional superconductors, in the Fe-HTSs both $d$-wave and extended $s$-wave pairing symmetries are close in energy. Probing the proximity between these very different superconducting states and identifying experimental parameters that can tune them is of central interest. Here we report high-pressure muon spin rotation experiments on the temperature-dependent magnetic penetration depth in the optimally doped nodeless $s$-wave Fe-HTS Ba$_{0.65}$Rb$_{0.35}$Fe$_2$As$_2$. Upon pressure, a strong decrease of the penetration depth in the zero-temperature limit is observed, while the superconducting transition temperature remains nearly constant. More importantly, the low-temperature behaviour of the inverse-squared magnetic penetration depth, which is a direct measure of the superfluid density, changes qualitatively from an exponential saturation at zero pressure to a linear-in-temperature behaviour at higher pressures, indicating that hydrostatic pressure promotes the appearance of nodes in the superconducting gap.

1 Laboratory for Muon Spin Spectroscopy, Paul Scherrer Institute, CH-5232 Villigen, Switzerland. 2 School of Physics and Astronomy, University of Minnesota, Minneapolis, Minnesota 55455, USA. 3 Leibniz-Institut für Festkörper- und Werkstoffforschung (IFW) Dresden, D-01171 Dresden, Germany. 4 Physik-Institut der Universität Zürich, Winterthurerstrasse 190, CH-8057 Zürich, Switzerland. 5 Institute of Low Temperature and Structure Research, Polish Academy of Sciences, 50-422 Wroclaw, Poland. 6 Department of Physics, Tbilisi State University, Chavchavadze 3, GE-0128 Tbilisi, Georgia. Correspondence and requests for materials should be addressed to Z.G. (email: zurab.guguchia@psi.ch).
After 6 years of intensive research on the Fe-based high-temperature superconductors (Fe-HTSs), no consensus on a universal gap structure has been reached. There is evidence that small differences in electronic or structural properties can lead to a strong diversity in the superconducting (SC) gap structure. On the one hand, nodeless isotropic gap functions were observed in optimally doped $\text{Ba}_1 - x \text{K}_x \text{Fe}_2\text{As}_2$, $\text{Ba}_1 - x \text{Rb}_x \text{Fe}_2\text{As}_2$ and $\text{BaFe}_2 - x \text{Ni}_x \text{As}_2$ as well as in $\text{BaFe}_2 - x \text{Co}_x\text{As}_2$, $\text{K}_x\text{Fe}_2 - x \text{Se}_2$ and $\text{FeTe}_1 - x \text{Se}_2$ (refs 1–8). On the other hand, signatures of nodal SC gaps were reported in $\text{LaOFeP}$, $\text{LiFeP}$, $\text{KFe}_2\text{As}_2$, $\text{BaFe}_2(\text{As}_1 - x \text{P}_x)$, $\text{BaFe}_2 - x \text{Ru}_x\text{As}_2$, $\text{FeSe}$ as well as in overdoped $\text{Ba}_1 - x \text{K}_x\text{Fe}_2\text{As}_2$ and $\text{BaFe}_2 - x \text{Ni}_x\text{As}_2$ (refs 7,9–17). Understanding what parameters of the systems control the different SC gap structures observed experimentally is paramount to elucidate the microscopic pairing mechanism in the Fe-HTSs and, more generally, to provide a deeper understanding of the phenomenon of high-temperature superconductivity. On the theoretical front, it has been proposed that both the $s^\pm$-wave and $d$-wave states are close competitors for the SC ground state$^{18–25}$, although there is no clear consensus on a universal gap structure. On the experimental front, it has been proposed that both the $d$-wave and $s$-wave state may be realized on removing electron or hole pockets. 

On the experimental front, we observe that while $\text{ARPES}$ measurements$^{1,30,31}$, in this article, we report on quantum critical point in $\text{Ba}_0.65\text{Rb}_{0.35}\text{Fe}_2\text{As}_2$. More importantly, we find a qualitative change in the low-temperature behaviour of $\Delta^2(T)$ as pressure is increased. While at $p = 0$ an exponential suppression characteristic of a nodeless superconductivity is observed, for $p = 2.25$ GPa a clear power-law behaviour is found. Because pressure does not affect the impurity concentration, which could promote power-law behaviour even for a nodeless system$^{33}$, our findings are suggestive of a nodeless to nodal SC transition. Our fittings to microscopic models reveal that this behaviour is more compatible with a $d$-wave state rather than an $s^\pm$ state with accidental nodes, suggesting that pressure promotes a change in the pairing symmetry.

**Results**

**Probing the vortex state as a function of pressure.** Figure 1a,b exhibit the transverse-field $\mu$SR time spectra for $\text{Ba}_{0.65}\text{Rb}_{0.35}\text{Fe}_2\text{As}_2$, measured at ambient $p = 0$ GPa and maximum applied pressure $p = 2.25$ GPa, respectively. The spectra above (45 K) and below (1.7 K) the SC transition temperature $T_c$ are shown. Above $T_c$ the oscillations show a small relaxation due to the random local fields from the nuclear magnetic moments. Below $T_c$ the relaxation rate strongly increases with decreasing temperature due to the presence of a non-uniform local magnetic field distribution as a result of the formation of a flux-line lattice in the SC state. Figure 1c,d show the Fourier transforms of the $\mu$SR time spectra shown in Fig. 1a,b, respectively. At $T = 5$ K the narrow signal around $\mu_0H_{0,SC} = 50$ mT (Fig. 1c,d) originates from the pressure cell, while the broad signal with a first moment $\mu_0H_{0,NS}$ marked by the solid arrow in Fig. 1c, arises from the SC sample.

Below $T_c$ a large diamagnetic shift of $\mu_0H_{0,SC}$ experienced by the muons is observed at all applied pressures. This is evident in Fig. 2a, where we plot the temperature dependence of the diamagnetic shift $\Delta B_{0,SC} = \mu_0[H_{0,SC} - H_{0,NS}]$ for $\text{Ba}_{0.65}\text{Rb}_{0.35}\text{Fe}_2\text{As}_2$ at various pressures, where $\mu_0H_{0,SC}$ denotes the internal field measured in the SC state and $\mu_0H_{0,NS}$ the internal field measured in the normal state at 45 K. Note that $\mu_0H_{0,NS}$ is temperature independent. This diamagnetic shift indicates the bulk character of superconductivity and excludes the possibility of field-induced magnetism$^{34}$ in $\text{Ba}_{0.65}\text{Rb}_{0.35}\text{Fe}_2\text{As}_2$ at all applied pressures. The SC transition temperature $T_c$ is determined from the intercept of the linearly extrapolated $\Delta B_{0,SC}$ curve to its zero line (we used the same criterion for determination of $T_c$ from $\Delta B_{0,NS}(T)$ as from the susceptibility data $\chi(T)$, presented in Supplementary Fig. 1a). It is found to be $T_c = 36.9(7)$ K and $35.9(5)$ K for $p = 0$ and 2.25 GPa, respectively. The ambient pressure value of $T_c$ is in perfect agreement with $T_c = 36.8(5)$ K obtained from susceptibility and specific heat measurements (Supplementary Note 1 and Supplementary Figs 1a,b and 2). At the highest pressure of $p = 2.25$ GPa applied, $T_c$ decreases only by $\approx 1$ K, indicating only a small pressure effect on $T_c$ in $\text{Ba}_{0.65}\text{Rb}_{0.35}\text{Fe}_2\text{As}_2$. The temperature dependence of the muon spin depolarization rate $\sigma_c$ of $\text{Ba}_{0.65}\text{Rb}_{0.35}\text{Fe}_2\text{As}_2$ in the SC state at selected pressures is shown in Fig. 2b; note that $\sigma_c$ is proportional to the second moment of the field distribution, which was extracted using the equations described in the Method section. At all applied pressures, below $T_c$ the relaxation rate $\sigma_c$ starts to increase from zero with decreasing temperature due to the formation of the flux-line lattice (we note that no pressure-induced magnetism is observed in $\text{Ba}_{0.65}\text{Rb}_{0.35}\text{Fe}_2\text{As}_2$, as shown in Supplementary Note 2 and Supplementary Figs 3a,b and 4). It is interesting that the low-temperature value $\sigma_c(5 \text{ K})$ increases substantially under pressure (Fig. 2b): $\sigma_c(5 \text{ K})$ increases about 30% from $p = 0$ to 2.25 GPa. Interestingly, the form of the temperature dependence of $\sigma_c$ (which reflects the topology of the SC gap, changes as a function of pressure. The most striking change is in the low-temperature behaviour of $\sigma_c(T)$. At ambient pressure $\sigma_c(T)$ shows a flat behaviour below $T/T_c = 0.4$, whereas the high-pressure data exhibit a steeper (linear) temperature dependence of $\sigma_c(T)$ below $T/T_c = 0.4$. We show in the following how these behaviours indicate the appearance of nodes in the gap function.
moments for the signals of the pressure cell and the sample, respectively. The solid lines in Figure 1 (after field cooling the sample from above $T_c$): (a,c) $p = 0$ GPa and (b,d) $p = 2.25$ GPa. Error bars are the s.e.m. in about $10^6$ events. The error of each bin in $A(t)$ are then calculated by s.e. propagation. The solid lines in a and b represent fits to the data by means of equation (3). The solid lines in c and d are the Fourier transforms of the fitted time spectra. The dashed and solid arrows indicate the first moments for the signals of the pressure cell and the sample, respectively.

Pressure-dependent magnetic penetration depth. To investigate a possible change of the symmetry of the SC gap, we note that $\lambda(T)$ is related to the relaxation rate $\sigma_{sc}(T)$ by the equation:

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

where $\gamma_{\mu}$ is the gyromagnetic ratio of the muon and $\Phi_0$ is the magnetic-flux quantum. Thus, the flat $T$ dependence of $\sigma_{sc}$ observed at $p = 0$ for low temperatures (Fig. 2b) is consistent with a nodeless superconductor, in which $\lambda^{-2}(T)$ reaches its zero-temperature value exponentially. On the other hand, the linear $T$ dependence of $\sigma_{sc}$ observed at $p = 2.25$ GPa (Fig. 2b) indicates that $\lambda^{-2}(T)$ reaches $\lambda^{-2}(0)$ linearly, which is characteristic of line nodes. This is the main result of this communication: pressure in an optimally doped Fe-HTS can tune a nodeless gap into a nodal gap. Although this qualitative analysis is robust, and independent of any fitting models for the gap function, it does not elucidate whether these nodes arise due to a nodal $s^{+}$ state or a $d$-wave state.

To proceed with a quantitative analysis, we consider the local (London) approximation ($\lambda \gg \xi$, where $\xi$ is the coherence length) and first employ the empirical $\lambda$-model. The latter, widely used in previous investigations of the penetration depth of multiband superconductors, assumes that the gaps occurring in different bands, besides a common $T_c$, are independent of each other. Then, the superfluid density is calculated for each component separately and added together with a weighting factor. For our purposes a two-band model suffices yielding:

$$\frac{\lambda^{-2}(T)}{\lambda^{-2}(0)} = \frac{\omega_1 \lambda^{-2}(T, \Delta_{01})}{\lambda^{-2}(0, \Delta_{01})} + \frac{\omega_2 \lambda^{-2}(T, \Delta_{02})}{\lambda^{-2}(0, \Delta_{02})},$$

where $\omega_1$ and $\omega_2$ are the weight factors.
where $\lambda(0)$ is the penetration depth at zero temperature, $\Delta_{0j}$ is the value of the $i$-th SC gap ($i=1, 2$) at $T=0$ K and $\omega_j$ is the weighting factor, which measures their relative contributions to $\lambda^{-2}$ (that is, $\omega_1 + \omega_2 = 1$).

The results of this analysis are presented in Fig. 3a–f, where the temperature dependence of $\lambda^{-2}$ for Ba$_{0.65}$Rb$_{0.35}$Fe$_2$As$_2$ is plotted at various pressures. We consider two different possibilities for the gap functions: either a constant gap, $\Delta_{0j} = \Delta_h$, or an angle-dependent gap of the form $\Delta_{0j}(\mathbf{p}) = \Delta_h \cos 2\varphi$, where $\varphi$ is the polar angle around the Fermi surface. The resulting functions $\lambda(T)$ are shown in the Methods section. The data at $p=0$ GPa are described remarkably well by two constant gaps, $\Delta_1 = 2.7(5)$ meV and $\Delta_2 = 8.4(3)$ meV. These values are in perfect agreement with our previous results and also with ARPES experiments, pointing out that most Fe-HTSs exhibit two-gap behaviour.

Remarkably, we find in both cases that the best fit gives $\Delta_h = 8.4(3)$ meV and $\Delta_0 = 2.7(5)$ meV (Supplementary Note 3). For simplicity, the anisotropies of the electron pockets are neglected, the Fermi velocities of the pockets are assumed to be nearly the same and the gaps are expanded in their leading harmonics. Thus, for the nodal $s^\pm$ state we have $\Delta_h = \Delta_{h_0}$ and $\Delta_0 = \Delta_{00}(|r| \cos 2\varphi)$, whereas for the $d$-wave state it follows that $\Delta_h = \Delta_{h_0} \cos 2\varphi$ and $\Delta_0 = \Delta_{00}$. Note the difference in the position of the nodes in each case: while for the $d$-wave case they are always at $\varphi = \pm \pi/4$, for the nodal $s^\pm$ the nodes exist only when $r < 1$ at arbitrary positions $\varphi \neq \pm \pi/4$ (albeit in different Fermi pockets). We note that from the fits one cannot completely rule out the possibility of small but non-vanishing values of $r$. Therefore, at least within our model, a nodal $s^\pm$ state is compatible with the data only if the accidental nodes are fine tuned to lie either at or very close to the diagonals of the electron pockets for a broad pressure range. Since the position of the accidental nodes is expected to be sensitive to the topology of the Fermi surface, and consequently to pressure, it seems more plausible that the gap state is $d$-wave, since in that case the position of the gaps is enforced by symmetry to be along the diagonals of the Brillouin zone.

To understand the implications of the fitting to a constant and an angle-dependent gap for finite pressures, we analyse the two different scenarios in which nodes can emerge: a nodal $s^\pm$ state (with gap functions of different signs in the hole and in the electron pockets) and a $d$-wave state. In the former, the position of the nodes are accidental, that is, not enforced by symmetry, while in the latter the nodes are enforced by symmetry to be on the Brillouin zone diagonals. Schematic representations of both scenarios are shown in Fig. 4, where a density plot of the gap functions is superimposed to the typical Fermi surface of the iron pnictides, consisting of one or more hole pockets at the centre of the Brillouin zone, and electron pockets at the border of the Brillouin zone.

In this figure, we set the accidental nodes of the $s^\pm$ state to be on the electron pockets, as observed by ARPES in the related compound BaFe$_2$(As$_{1-x}$P$_x$)$_2$ (ref. 17). Note that in the $d$-wave state, while nodes appear in the hole pockets, the electron pockets have nearly uniform gaps. Thus, the fact that the fitting to the $\alpha$-model gives a constant and an angle-dependent gap is consistent with a $d$-wave state.

To contrast the scenarios of a nodal $s^\pm$ gap and a $d$-wave gap, we consider a microscopic model (Supplementary Note 3) that goes beyond the simplifications of independent gap functions of the $\alpha$-model discussed above. In this microscopic model, the fully coupled nonlinear gap equations are solved for a hole pocket $h$ and two electron pockets $e_1, e_2$, and the penetration depth is calculated at all temperatures (Supplementary Fig. 5). The free parameters are then the density of states of the pockets, the amplitude of the pairing interaction and the gap functions themselves (Supplementary Note 3). For simplicity, the anisotropies of the electron pockets are neglected, the Fermi velocities of the pockets are assumed to be nearly the same and the gaps are expanded in their leading harmonics. Thus, for the nodal $s^\pm$ state we have $\Delta_h = \Delta_{h_0}$ and $\Delta_0 = \Delta_{00}(|r| \cos 2\varphi)$, whereas for the $d$-wave state it follows that $\Delta_h = \Delta_{h_0} \cos 2\varphi$ and $\Delta_0 = \Delta_{00}$. Note the difference in the position of the nodes in each case: while for the $d$-wave case they are always at $\varphi = \pm \pi/4$, for the nodal $s^\pm$ the nodes exist only when $r < 1$ at arbitrary positions $\varphi \neq \pm \pi/4$ (albeit in different Fermi pockets). We note that from the fits one cannot completely rule out the possibility of small but non-vanishing values of $r$. Therefore, at least within our model, a nodal $s^\pm$ state is compatible with the data only if the accidental nodes are fine tuned to lie either at or very close to the diagonals of the electron pockets for a broad pressure range. Since the position of the accidental nodes is expected to be sensitive to the topology of the Fermi surface, and consequently to pressure, it seems more plausible that the gap state is $d$-wave, since in that case the position of the gaps is enforced by symmetry to be along the diagonals of the Brillouin zone.
Figure 4 | Schematic representation of the nodal $s^\pm$ and $d$-wave states. In both figures, a density plot of the gap function is superimposed to a representative Fermi surface consisting of a hole pocket ($h$) at the centre and an electron pocket ($e$) at the borders of the Brillouin zone. In the nodal $s^\pm$ states (a) the nodes are not enforced by symmetry (here they are located at the electron pockets). In the $d$-wave state (b) the nodes are enforced by symmetry to be on the diagonals of the Brillouin zone, and therefore can only cross the hole pockets.

hole pockets regardless of the value of the pressure (Supplementary Fig. 7a,b).

The pressure dependence of all the parameters extracted from the data analysis within the $\alpha$-model are plotted in Fig. 5a–c. From Fig. 5a a substantial decrease of $\lambda(0)$ with pressure is evident. At the maximum applied pressure of $p = 2.25$ GPa the reduction of $\lambda(0)$ is $\sim 15\%$ compared with the value at $p = 0$ GPa. Both $\lambda_1$ and $\lambda_2$ show a small reduction on increasing the pressure from $p = 0$ to 1.17 GPa, while above $p = 1.17$ GPa the gaps values stay constant. On the other hand, the relative contribution $\omega_2$ of the small gap to the superfluid density increases by approximately factor of 2 for the maximum applied pressure of $p = 2.25$ GPa (Fig. 5c), indicating a spectral weight shift to the smaller gap. The parameters extracted from the microscopic model are discussed in Supplementary Note 3.

Discussion

The main finding of our paper is the observation that pressure promotes a nodal SC gap in Ba$_{0.65}$Rb$_{0.35}$Fe$_2$As$_2$. This conclusion is model independent, as it relies on the qualitative change in the low-temperature behaviour of $\Delta \lambda_{\perp}^{-2}$ from exponential to linear in $T$ on applied pressure. To our knowledge this is the first direct experimental demonstration of a plausible pressure-induced change in the SC gap structure in a Fe-HTS. Two possible gap structures could be realized at finite pressures: a nodal $s^\pm$ state and a $d$-wave state. In the first case, the change from nodeless $s^\pm$ to nodal $s^\pm$ is a crossover rather than a phase transition$^{42,43}$, whereas in the latter it is an actual phase transition that could harbour exotic pairing states, such as $s+id$ (refs 21,23,24) or $s+d$ (ref. 44).

Additional results provide important clues of how pressure may induce either a nodal $s^\pm$ or a $d$-wave state. In the closely related optimally doped compound Ba$_{0.6}$K$_{0.4}$Fe$_2$As$_2$, Raman spectroscopy$^{27}$, as well as theoretical calculations$^{21,20}$, reveal a subdominant $d$-wave state close in energy to the dominant $s^\pm$ state. Pressure may affect this intricate balance, and tip the balance in favour of the $d$-wave state. On the other hand, theoretical calculations have shown that the pnictogen height is an important factor in determining the structure of the $s^\pm$ SC order parameter$^{18,45}$. A systematic comparison of the quasiparticle excitations in the 1111, 122 and 111 families of Fe-HTSs showed that the nodal $s^\pm$ state is favoured when the pnictogen height decreases below a threshold value of $\approx 1.33$ Å (ref. 46). Hydrostatic pressure may indeed shorten the pnictogen height and consequently modify the $s^\pm$ gap structure from nodeless to nodal. Although our fitting of the penetration depth data to both a microscopic model and an effective $\alpha$-model suggest that the $d$-wave state is more likely to be realized than the nodal $s^\pm$ state, further quantitative calculations of the pressure effect are desirable to completely discard a nodal $s^\pm$ state.

Besides the appearance of nodes with pressure, another interesting observation is the reduction of $\lambda(0)$ under pressure, despite the fact that $T_c$ remains nearly unchanged. Interestingly,
in the compound BaFe_{2}As_{2}−xP_{x}, a sharp enhancement of $\lambda(0)$ is observed as optimal doping is approached from the overdoped side\(^\text{5,6}\), which has been interpreted in terms of a putative quantum critical point (QCP) inside the SC dome\(^\text{7–9}\). In Ba_{0.65}Rb_{0.35}Fe_{2}As_{2}, if such a putative QCP is also present, pressure is likely to move the system away from the putative QCP, which, according to the results of BaFe_{2}As_{2}−xP_{x} would explain the observed suppression of the penetration depth at $T = 0$. This scenario does not explain why $T_{C}$ stays nearly constant under pressure, but this could be due to the intrinsic flatness of $T_{C}$ around optimal doping in Ba_{1−x}Rb_{x}Fe_{2}As_{2}. Note that a similar behaviour for $\lambda(0)$ and $T_{C}$ with pressure has been recently observed in LaFeAsO_{1−x}F_{x} (ref. 50), but interpreted in terms of the interplay between impurity scattering and pressure. To distinguish between these two scenarios, pressure-dependent studies of the quasiparticle mass in Ba_{0.65}Rb_{0.35}Fe_{2}As_{2} are desirable to probe whether a putative QCP is present or not in this compound.

In conclusion, the zero-temperature magnetic penetration depth $\lambda(0)$ and the temperature dependence of $\lambda−T$ were studied in optimally doped Ba_{0.65}Rb_{0.35}Fe_{2}As_{2} by means of µSR experiments as a function of pressure up to $p \approx 2.25$ GPa. The SC transition temperature stays nearly constant under pressure, whereas a strong reduction of $\lambda(0)$ is observed, possibly related to the presence of a putative quantum critical point. Our main result is the observation of a qualitative change in the low-temperature behaviour of $\Delta \lambda−T(T)$ from exponential to linear in the investigated Fe-based superconductor as pressure is increased. This most likely indicates that a nodal SC gap is promoted by hydrostatic pressure. Model calculations favour a $d$-wave over a nodal $s^{\pm}$-wave pairing as the origin for the nodal gap. The present results offer important benchmarks for the elucidation of the complex microscopic mechanism responsible for the observed non-universality of the SC gap structure and of high-temperature superconductivity in the Fe-HTSs in general.

**Methods**

**Sample preparation.** Polycrystalline samples of Ba_{0.65}Rb_{0.35}Fe_{2}As_{2} were prepared in evacuated quartz ampoules by a solid-state reaction method. FeAs, BaAs and RbAs were obtained by reacting high-purity As (99.999 %), Fe (99.9%), Ba (99.9%) and Rb (99.95%) at 800, 650 and 500 °C, respectively. Using stoichiometric amounts of BaAs or RbAs and FeAs, the terminal compounds BaFe_{2}As_{2} and RbFe_{2}As_{2} were synthesized at 950 and 650 °C, respectively. Finally, samples of Ba_{1−x}Rb_{x}Fe_{2}As_{2} with $x \sim 0.35$ were prepared from appropriate amounts of single-phase BaFe_{2}As_{2} and RbFe_{2}As_{2}. The components were mixed, pressed into pellets, placed into alumina crucibles and annealed for 100 h under vacuum at 650 °C with one intermittent grinding. Powder X-ray diffraction analysis revealed that the synthesized samples are single-phase materials.

**Method for creation and measurement of high pressures.** Pressures up to 2.4 GPa were generated in a double-wall piston-cylinder type of cell made of MP35N material, especially designed to perform µSR experiments under pressure\(^\text{10,11}\). As a pressure-transmitting medium Daphne oil was used. The pressure was measured by tracking the SC transition of a very small indium plate by a.c. susceptibility. The filling factor of the pressure cell was maximized. The fraction of the muons stopping in the sample was ~40%.

**µSR experiment.** Zero-field and transverse-field (TF) µSR experiments at ambient and under various applied pressures were performed at the µE1 beamline of the Paul Scherrer Institute, Switzerland, using the dedicated general purpose decay channel instrument (GPD) spectrometer, where an intense high-energy ($p_{0} = 100$ MeV/c) beam of muons is implanted in the sample through the pressure cell. A gas-flow He (base temperature ~4 K) and a VARIOX cryostat (base temperature ~1.3 K) were used. High-energy muons ($p_{0} = 100$ MeV/c) were implanted in the sample. Forward and backward positron detectors with respect to the initial muon spin polarization were used for the measurements of the µSR asymmetry time spectrum $A(t)$. The typical statistics for both forward and backward detectors were 6 million. All zero-field and transverse-field µSR experiments were performed by stabilizing the temperature in before recording the µSR time spectra. Note that a precise calibration of the GPD results was carried out at the nM3 beamline using the low-background GPS. The µSR time spectra were analysed using the free software package MUSRFIT\(^\text{36}\). The µSR experiment nearly 100% spin-polarized muons $\mu^{+}$ are implanted into the sample one at a time. The positively charged $\mu^{+}$ thermalize at interstitial lattice sites, where they act as magnetic microprobes. In a magnetic material the muon spin precesses in the local field $B_{\mu}$ at the muon site with the Larmor frequency $\nu_{L} = \gamma_{\mu}/2\pi B_{\mu}$ (muon gyromagnetic ratio $\gamma_{\mu}/(2\pi) = 135.5$ MHz T$^{-1}$). By means of µSR it is possible to measure length scale of superconductor because of the magnetic penetration depth $\lambda$. When a type II superconductor is cooled below $T_{c}$ in an applied magnetic field ranging between the lower ($H_{c1}$) and the upper ($H_{c2}$) critical field, a vortex lattice is formed, which in general is incommensurate with the crystal lattice, and the vortex cores will be separated by much larger distances than those of the unit cell. Because the implanted muons stop at given crystallographic sites, they will randomly probe the field distribution of the vortex lattice. Such measurements need to be performed in a field applied perpendicular to the initial muon spin polarization (so-called transverse-field configuration).

**Analysis of transverse-field-µSR data.** Our zero-field µSR experiments (Supplementary Note 2) reveal a pressure-independent magnetic fraction of about 10% in the sample, caused by the presence of diluted Fe moments as discussed in previous µSR studies. The signal from the magnetically ordered parts vanishes within the first 0.2 µs. Thus, the fits of transverse-field data were restricted to times $t > 0.2$ µs for all temperatures.

The transverse-field µSR data were analysed by using the following functional form\(^\text{36}\):

$$P(t) = A_{\mu} \exp \left[ -\frac{\nu_{L} t}{\nu_{L} + i \phi} \right] 
+ A_{pc} \exp \left[ -\frac{\nu_{L} t}{\nu_{L} + i \phi} \right],$$

where $A_{\mu}$ and $A_{pc}$ denote the initial asymmetries of the sample and the pressure cell, respectively. $\gamma_{\mu}/(2\pi)$ ~ 135.5 MHz T$^{-1}$ is the muon gyromagnetic ratio, $\phi$ is the initial phase of the muon spin ensemble and $B_{\mu}$ represents the internal magnetic field at the muon site. The relaxation rates $\sigma_{\mu}$ and $\sigma_{pc}$ characterize the damping due to the formation of the vortex lattice in the SC state and of the nuclear magnetic dipolar contribution, respectively. In the analysis $A_{\mu}$ was assumed to be constant over the entire temperature range and was fixed to the value obtained above $T_{c}$, where only nuclear magnetic moments contribute to the muon relaxation rate $\phi$. The Gaussian relaxation rate $\sigma_{pc}$ reflects the depolarization due to the nuclear magnetism of the pressure cell. It can be seen from the Fourier transforms shown in Fig. 1c,d that the width of the pressure cell signal increases below $T_{c}$ as shown previously\(^\text{23}\), this is due to the influence of the diamagnetic moment of the SC sample on the pressure cell, leading to a temperature-dependent $\sigma_{pc}$ below $T_{c}$.

To consider this influence, we assume a linear coupling between $\sigma_{pc}$ and the field shift of the internal magnetic field in the SC state: $\sigma_{pc}(T) = \sigma_{pc}(T_{c}) + (T_{c} − T)/C_{pc}$, where $\sigma_{pc}(T_{c})$ is the temperature-independent Gaussian relaxation rate. $\mu_{H_{c1,2}}$ and $\mu_{H_{c2,3}}$ are the internal magnetic fields measured in the normal and in the SC state, respectively. As indicated by the solid lines in Fig. 1a–d, the µSR data are well described by equation (1). The solid lines in Fig. 1c,d are the Fourier transforms of the fitted curves shown in Fig. 1a,b. The model used describes the data rather well.

**Analysis of $\lambda(0)$.** As pointed out in the manuscript, for polycrystalline samples the temperature dependence of the London magnetic penetration depth $\lambda(T)$ is related to the muon spin depolarization rate $\sigma_{\mu}(T)$ by equation (1) (see the main text). Equation (1) is valid when the separation between the vortices is smaller than $\lambda$ and the applied field small with respect to the second critical field $B_{c2}$. In this case according to the London model $\sigma_{\mu}$ is field independent\(^\text{11,12}\). Field-dependent measurements of $\sigma_{\mu}$ at ambient pressure was reported previously\(^\text{23}\). It was observed that first $\sigma_{\mu}$ strongly increases with increasing magnetic field until reaching a maximum at $\mu_{0}H \approx 0.3$ T and then above 0.3 T stays nearly constant up to the highest field (0.64 T) investigated. Such a behaviour is expected within the London model and is typical for polycrystalline HTSs\(^\text{24}\).

$\lambda(T)$ was calculated within the (London) approximation ($\lambda \gg \xi$) by the following expression\(^\text{36,37}\):
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Author contributions

Project planning: Z.G.; sample growth: Z.B.; data analysis: Z.G., R.M.F. and A.V.; interpretation: Z.G., R.M.F. and J.K.; draft writing: Z.G. with contributions and/or comments from all authors.

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