

Superconductivity in Trilayer Nickelate $\text{La}_4\text{Ni}_3\text{O}_{10}$ under Pressure

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Nickelate superconductors have attracted a great deal of attention over the past few decades due to their similar crystal and electronic structures with high-temperature cuprate superconductors. Here, we report superconductivity in a pressurized Ruddlesden-Popper phase single crystal $\text{La}_4\text{Ni}_3\text{O}_{10}$ ($n = 3$) and its interplay with the density wave order in the phase diagram. With increasing pressure, the density wave order, as indicated by the anomaly in the resistivity, is progressively suppressed, followed by the emergence of superconductivity around 25 K under the $I4/mmm$ space group. The susceptibility measurements confirm bulk superconductivity with a volume fraction exceeding 80%. Moreover, theoretical analysis unveils that antiferromagnetic superexchange interactions can serve as the effective pairing interaction for the emergence of superconductivity in pressurized $\text{La}_4\text{Ni}_3\text{O}_{10}$. Our research provides a new platform for the investigation of the unconventional superconductivity mechanism in Ruddlesden-Popper trilayer perovskite nickelates.

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I. INTRODUCTION

The discovery of superconductivity in cuprates has opened an avenue to explore high-temperature (high- T_c) superconductivity [1,2]. Almost all the discovered cuprate superconductors so far share a common two-dimensional CuO_2 plane and half-filled $\text{Cu } 3d^9$ electronic configuration, which are believed to be essential to the emergence of high- T_c superconductivity [3,4]. Besides cuprates, iron-based superconductors are another prime material family with superconductivity well above the McMillan limit at

ambient pressure [5,6]. In contrast to the cuprates, iron-based superconductors with Fe $3d^6$ electronic configuration have a multiband and multiorbital nature [7,8]. Despite sustained efforts that have been devoted for more than three decades, the full understanding of their superconductivity mechanism remains a challenging and long-sought problem in condensed matter physics.

Nickelates are considered to be an ideal candidate for exploring new cupratelike high- T_c superconductivity due to the proximity of $\text{Ni}^+(3d^9)$ electron configuration [9]. However, superconductivity was not experimentally realized in nickelates until 2019, when Li *et al.* discovered superconductivity within the thin films of infinite-layer nickelates $\text{Nd}_{1-x}\text{Sr}_x\text{NiO}_2$ with $T_c \sim 9\text{--}15$ K [10,11], which quickly promoted renewed interest in the exploration of nickel-based superconductors. Subsequently, T_c has been extended to other infinite-layer nickelates with different rare-earth elements [12–15]. However, superconductivity is limited to thin-film materials. Structurally, whether there exists a nickelate superconductor family with bulk superconductivity is still an open question.

Recently, the experimental breakthrough revealed high- T_c superconductivity above liquid-nitrogen temperature in pressurized bulk $\text{La}_3\text{Ni}_2\text{O}_7$ single crystals, which is one of the Ruddlesden-Popper (RP) phases $\text{La}_{n+1}\text{Ni}_n\text{O}_{3n+1}$ ($n = 2$) with the valence of $\text{Ni}^{2.5+}$ [16]. The synchrotron x-ray diffraction (XRD) patterns and density functional theory calculations indicate that $\text{La}_3\text{Ni}_2\text{O}_7$ undergoes a structural phase transition from Amm to $I4/mmm$ with the emergence of superconductivity, and a series of follow-up experiments have elaborated that oxygen defects and hydrostatic environment are essential for the discovery of superconductivity [17–23]. Moreover, many theoretical works are put forward to understand the underlying mechanisms of high- T_c superconductivity in $\text{La}_3\text{Ni}_2\text{O}_7$ [24–39].

Among RP phase nickelates, the trilayer $\text{La}_4\text{Ni}_3\text{O}_{10}$ ($n = 3$) has drawn more attention due to the metallic ground state and metal-to-metal transition at intermediate temperatures [40–43]. The nominal valence state of Ni is $+2.67$ in $\text{La}_4\text{Ni}_3\text{O}_{10}$, which is usually believed to be given by mixed valence states of Ni^{2+} and Ni^{3+} . Considering the similar crystal structure and Ni valence with $\text{La}_3\text{Ni}_2\text{O}_7$, it is natural to consider investigating the superconductivity and electronic structure in pressurized $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals.

In this work, we grow high-quality $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystal with an optical-image floating zone furnace under high oxygen pressure and perform high-pressure transport measurements using NaCl as the pressure-transmitting medium to improve the hydrostatic environment. We find superconductivity with a T_c of 25 K in $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals under pressure. A distinctive diamagnetic transition consists with the resistance measurements, further confirming the emergence of superconductivity in $\text{La}_4\text{Ni}_3\text{O}_{10}$

single crystals. By combining scanning tunneling microscopy and spectroscopy, angle-resolved photoemission spectroscopy, and *ab initio* calculation, we systematically investigate the electronic structure of trilayer nickelate $\text{La}_4\text{Ni}_3\text{O}_{10}$. These observations broaden and contribute to the superconducting family of nickelates.

II. RESULTS AND DISCUSSION

A. Trilayer nickelate

$\text{La}_4\text{Ni}_3\text{O}_{10}$, one of the RP phase $\text{La}_{n+1}\text{Ni}_n\text{O}_{3n+1}$ ($n = 3$) nickelates, crystallizes in a monoclinic structure ($a = 5.41$ Å, $b = 5.47$ Å, $c = 14.22$ Å), consisting of three layers of perovskite-type LaNiO_3 separated by a single rocksalt-type LaO layer along the crystallographic c -axis direction [Fig. 1(a)]. As shown in Fig. 1(b), high-angle annular dark-field (left) and annular bright-field (right) transmission electron microscopy images [Fig. 1(b)] illustrate that each unit contains well-ordered three layers of inclined NiO_6 octahedral layers stacking along the axis direction. The inclination of the NiO_6 octahedra leads to the bucking of the Ni-O planes. The side-view crystal structure of the La-Ni-O layer in the monoclinic phase of $\text{La}_4\text{Ni}_3\text{O}_{10}$ is shown in Fig. 1(c). As we can see, the Ni-O bonds are tilted in the b - c plane instead of the a - c plane, resulting in a unidirectional structural reconstruction compared to that of the tetragonal phase. Figure 1(d) shows the in-plane unit cells for the tetragonal and the monoclinic phases. The size of the reconstructed monoclinic unit cell is twice that of the tetragonal unit cell due to the structural reconstruction. The unidirectional reconstruction results in a stripelike structure, as illustrated by the light blue ribbons in Fig. 1(d). The periodicity of the stripe is $\sqrt{2}$ times of the tetragonal lattice a_{tet} , as indicated by the $Q_{\sqrt{2}}$ in the Fourier transform (FT) of the NiO lattice [Fig. 1(e)].

The unidirectional reconstruction in $\text{La}_4\text{Ni}_3\text{O}_{10}$ can be confirmed by the low-temperature scanning tunneling microscopy (STM) measurements. Figure 1(f) shows a typical STM topography taken on the large-area LaO-terminated surface. On the LaO surface, we observe an ordered square lattice with some adatomlike impurities and several riplike dark stripes along the diagonal direction of the square lattice. These features can be seen more clearly in the enlarged STM topography as shown in Fig. 1(g). The square lattice is measured to be 0.40 nm, which is consistent with the in-plane LaO tetragonal lattice as shown in Fig. 1(d). This indicates that the riplike dark stripes are along the reconstruction direction. To further quantify the reconstructed surface, we perform an FT on the STM topography [Fig. 1(h)]. In the FT image, apart from the Q_{tet} originated from LaO in-plane tetragonal lattice, there is a set of unidirectional Q_1 vectors, which is $1/\sqrt{2}$ times of the Q_{tet} . The length and direction of the Q_1 wave vector are the same as the $Q_{\sqrt{2}}$ in Fig. 1(e), indicating that there exists a stripelike unidirectional structural reconstruction on the

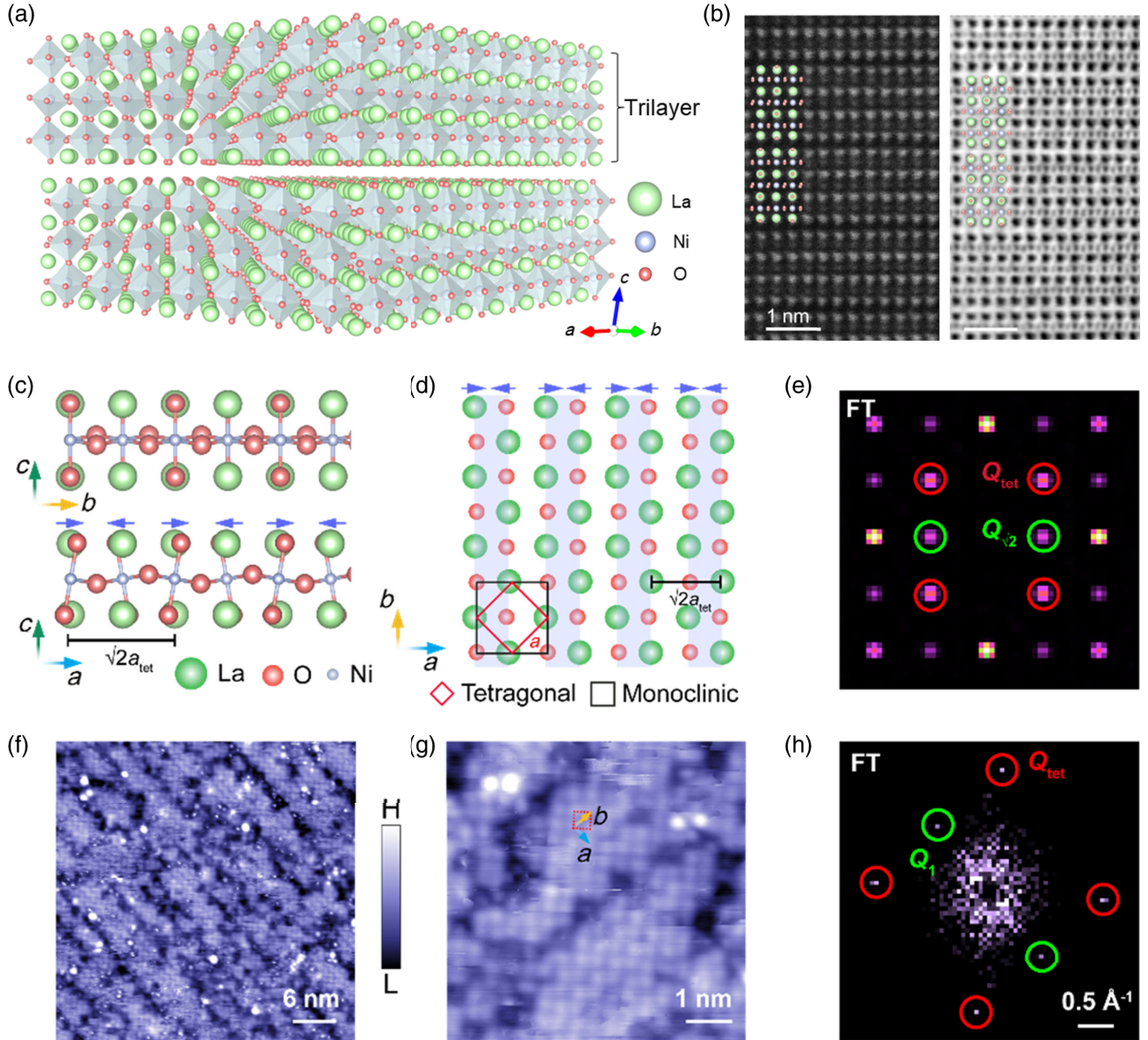


FIG. 1. Crystal structure and surface topographies of $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals. (a) Schematic illustration of the crystal structure of $\text{La}_4\text{Ni}_3\text{O}_{10}$ crystal. (b) High-angle annular dark-field (left) and annular bright-field (right) STEM images of $\text{La}_4\text{Ni}_3\text{O}_{10}$ along the b axis. (c) Side-view crystal structures along the a - c plane (upper) and the b - c plane (lower), respectively. (d) The top-view crystal structure of the LaO surface. The appended red and black squares are the unit cells for the tetragonal and monoclinic phases. The light blue ribbons indicate the unidirectional reconstruction. (e) FT image of the schematic crystal structure shown in (d). The red and green circles indicate the tetragonal Bragg lattice (Q_{tet}) and the unidirectional structural distortion pattern ($Q_{\sqrt{2}}$), respectively. (f),(g) Constant-current STM topographies taken on the LaO surface. The colored arrows in (g) indicate the monoclinic lattice. (h) FT image of the STM topography shown in (g).

LaO surface. Interestingly, the riplike dark stripes in the real space are also along the reconstructed stripe direction, and they can contribute broader signals in the FT images [Fig. 1(g); also see Fig. S3 in Supplemental Material [44]]. This suggests that the riplike dark stripes might be related to the reconstruction in $\text{La}_4\text{Ni}_3\text{O}_{10}$ due to the combination of the unidirectional strainlike effect and the local oxygen vacancies. Although the unidirectional reconstruction has been revealed, we find no clear signs of

the charge density wave pattern in the STM topographies on the LaO-terminated surface.

B. Transport under ambient pressure

Figure 2 show the resistivity, magnetization, and heat capacity measured on $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals. The resistivity decreases with cooling temperature from 300 K, revealing typical metallic behavior. A pronounced

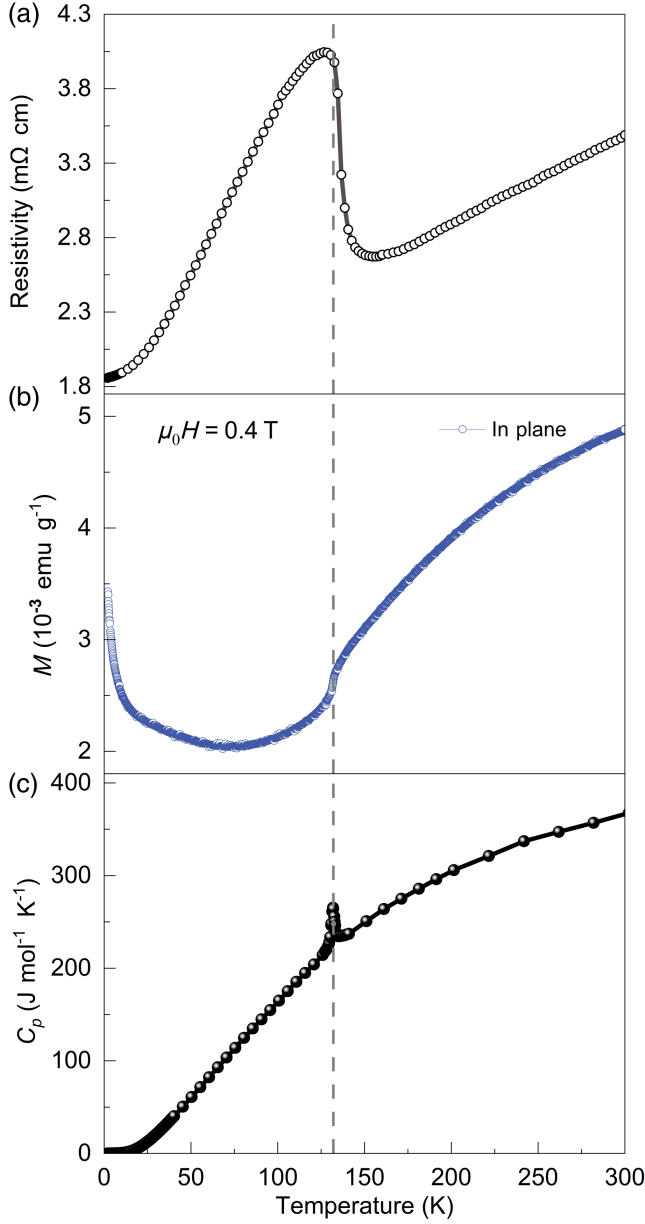


FIG. 2. Physical properties of $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals. (a) Temperature-dependent resistivity. (b) Temperature-dependent magnetization. (c) Temperature-dependent heat capacity.

anomalous peak at around 130 K suggests possible density-wave states. The in-plane magnetization was measured under a magnetic field of 0.4 T [41,44], which showed a sharp decrease at the $T^* \sim 132$ K, similar to typical charge or spin density wave materials such as $\text{K}_{0.3}\text{MoO}_3$ [45], CsV_3Sb_5 [46], and LaFeAsO [5]. The temperature-dependent heat capacity also exhibits an anomalous peak at the same temperature, and it can be well fitted by the formula $C_p/T = \gamma + \beta T^2$, where γ is the Sommerfeld coefficient. The fit yields $\gamma = 13.5$ mJ mole $^{-1}$ K $^{-2}$ and $\beta = 0.34$ mJ mole $^{-1}$ K $^{-4}$, which agrees with the previous

measurement [41]. The Debye temperatures and density of states $N(E_F)$ can be estimated as 459 K and 2.1 states eV $^{-1}$ per Ni, respectively.

Furthermore, we investigate the electronic structure of $\text{La}_4\text{Ni}_3\text{O}_{10}$ using high-resolution angle-resolved photoemission spectroscopy (ARPES) (Supplemental Material Fig. S4 [44]). Overall, the experimental band structure of $\text{La}_4\text{Ni}_3\text{O}_{10}$ is very similar to that of $\text{La}_3\text{Ni}_2\text{O}_7$ [30], despite their different crystal structure, suggesting unified electronic properties of nickelates in Ruddlesden-Popper phases.

C. Superconductivity in pressurized single crystals

The resistivity anomaly indicates that $\text{La}_4\text{Ni}_3\text{O}_{10}$ is located in the vicinity of an electronic instability. It is well known that superconductivity often appears near the critical point of structural or density waves. In this respect, pressure has been proven to be a convenient and effective way to modify lattice parameters and systematically influence the corresponding electronic states without introducing other factors. Motivated by these advantages, we measured the electrical resistance of $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals at various pressures. Figure 3(a) shows the typical resistance curves of $\text{La}_4\text{Ni}_3\text{O}_{10}$ for pressure up to 60 GPa. The resistance anomaly is sensitive to external pressures. With increasing pressure to 2.3 GPa, the temperature of the resistance anomaly shifts to approximately 120 K. As the pressure continues to increase to 14.0 GPa, the anomalous feature in the resistance curve becomes broadened and less pronounced. At 23.1 GPa, the resistance anomaly is suppressed further, and the resistance curve displays a slight upturn at the lower temperatures followed by a downturn feature. When the pressure approaches 32.3 GPa, the temperature dependence of resistance changes to that of a normal metal. Interestingly, a distinct behavior characterized by a resistance drop of about 8.5 K emerges, and this behavior becomes more pronounced as the pressure is increased, indicating the superconducting transition. The superconducting T_c continuously increases and reaches 23.5 K at 56.8 GPa. In particular, with the appearance of superconductivity, T -linear resistance becomes clear gradually in the normal state [Fig. 3(b)]. It is worth noting that the resistance drops by 95% from the normal state to the lowest temperature, which suggests a superconducting transition. Several independent high-pressure transport measurements on $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals provide consistent and reproducible results (Figs. S5 and S6 in Supplemental Material [44]), confirming intrinsic superconductivity under high pressure. Compared with run I, the resistance anomaly peak was suppressed below 10 GPa in run II (Fig. S4 in Supplemental Material [44]), followed by the emergency of sharp drops. The superconducting T_c reaches 25 K at approximately 42.5 GPa.

To further determine whether the observed resistance drop is truly associated with a superconducting transition, we

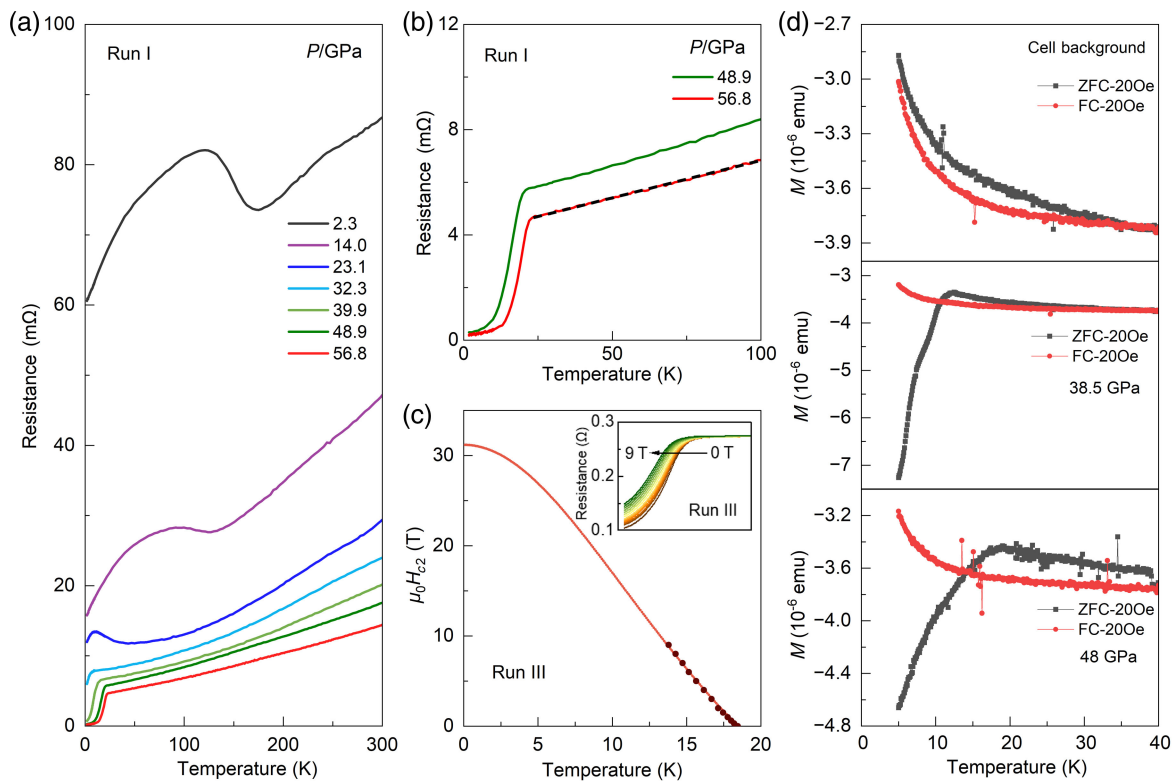


FIG. 3. Transport properties of $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals as a function of the pressure. (a) Electrical resistance of $\text{La}_4\text{Ni}_3\text{O}_{10}$ as a function of the temperature from 2.3 to 56.8 GPa in run I. (b) Linear resistance in the normal state. (c) Temperature dependence of the upper critical field $\mu_0 H_{c2}$ for $\text{La}_4\text{Ni}_3\text{O}_{10}$ at 67.7 GPa in run III. T_c is determined as 90% of the resistance at the onset T_c . The solid lines represent the Ginzburg-Landau fitting. The inset shows the low-temperature resistance under various magnetic fields. (d) dc magnetic susceptibility of $\text{La}_4\text{Ni}_3\text{O}_{10}$ under high pressures.

performed resistance measurements under various magnetic fields. As displayed in Fig. 3(c) and Supplemental Material Fig. S6 [44], the resistance drop of $\text{La}_4\text{Ni}_3\text{O}_{10}$ is progressively suppressed to lower temperatures, which implies a superconducting transition. The onset of superconductivity seems insensitive to the external magnetic fields, and the transition width becomes broad with increasing magnetic field, which is similar to high- T_c cuprates. As shown in the inset of Fig. 3(c), superconductivity survives even under magnetic field $\mu_0 H = 9$ T. The upper critical field $\mu_0 H_{c2}$ can be well fitted using the empirical Ginzburg-Landau formula $\mu_0 H_{c2}(T) = \mu_0 H_{c2}(1 - t^2)/(1 + t^2)$, where $t = T/T_c$. The superconducting transition is robust and extrapolated upper critical field $\mu_0 H_{c2}(0)$ of $\text{La}_4\text{Ni}_3\text{O}_{10}$ can reach 31.1 T at 67.7 GPa in run III, which yields a Ginzburg-Landau coherence length $\xi_{\text{GL}}(0)$ of 3.25 nm. To further validate pressure-induced superconductivity, we utilized a specially designed microfabricated beryllium-copper alloy diamond anvil cell (DAC) to conduct ultrasensitive dc magnetic susceptibility measurements under high pressure [Fig. 3(d) and Supplemental Material Fig. S7 [44]], with the aim of observing the diamagnetic effect characteristic of superconductors. From the raw data of high-pressure dc magnetic susceptibility, no significant transitions were detected in the

background data of the DAC, regardless of whether the measurements were performed in zero-field cooling (ZFC) or field cooling modes. However, when neon was employed as the pressure-transmitting medium and the sample pressure was increased to 38.5 GPa, a distinct diamagnetic transition was observed below 12.3 K in the ZFC mode, confirming bulk superconductivity with a volume fraction exceeding 80%. The superconducting transition temperature derived from these magnetic measurements at this pressure was in agreement with the results obtained from electrical transport measurements, further corroborating the emergence of superconductivity under high pressure. When the pressure was increased to 48.0 GPa, the superconducting transition temperature was enhanced, with diamagnetic below 19 K, consistent with the superconducting transition phase diagram of the sample under pressure. In summary, through high-pressure dc magnetic susceptibility measurements, we directly observed the diamagnetic effect of superconductors under high pressure, thereby substantiating that the sample exhibits superconductivity under such conditions.

To clarify the emergency of superconductivity, we perform *in situ* high-pressure XRD measurements up to 80 GPa. As shown in Figs. 4(a) and 4(b), except for two peaks arising from the Re gasket, all peaks could be

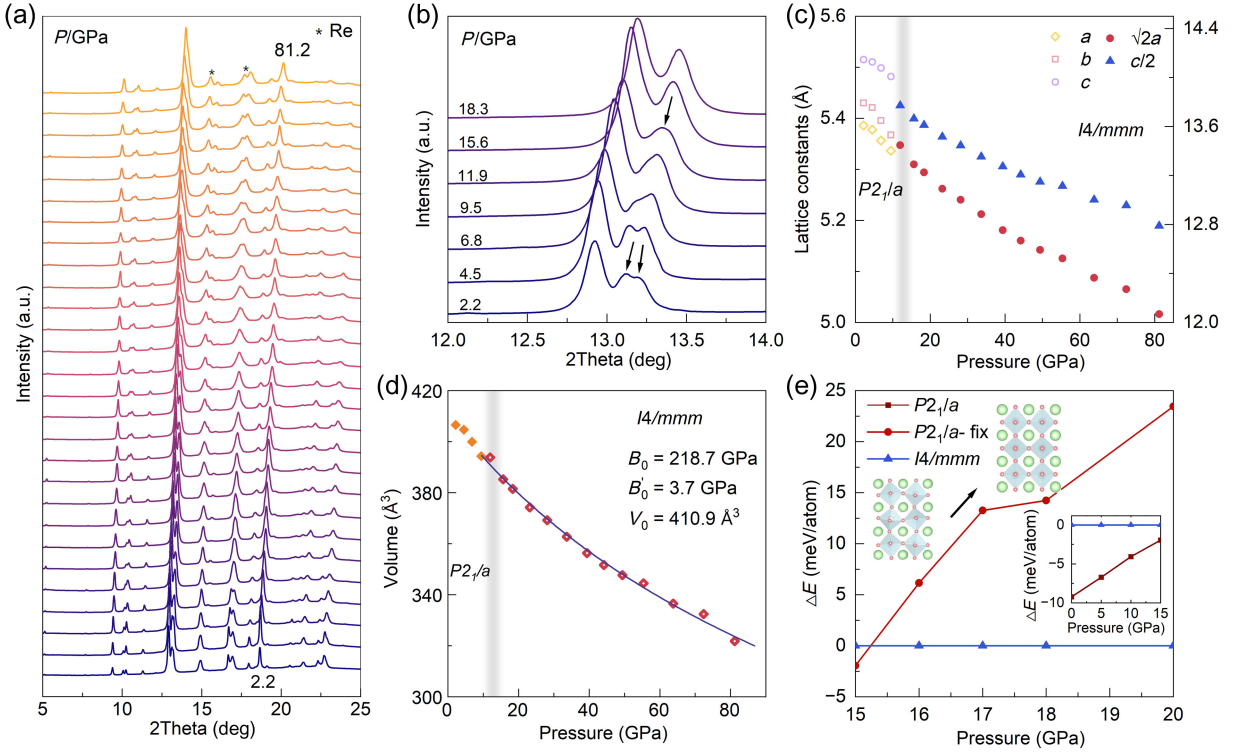


FIG. 4. Structural transition of $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals as a function of the pressure. (a) Synchrotron XRD patterns of powder samples from 2.2 to 81.2 GPa. The asterisk refers to the diffraction peaks of the Re gasket. (b) Details of the evolutions of the (0 2 0) and (2 0 -1) peaks under pressure. (c) The evolution of lattice parameters refined from the XRD patterns. (d) The pressure dependence of volume of the $P2_1/a$ and $I4/mmm$ space group derived from the XRD refinement results. (e) The difference in enthalpy between the space groups $P2_1/a$ and $I4/mmm$ as a function of the pressure calculated using the first-principles method. The inset shows the modification of NiO_6 octahedra stacking mode under pressure.

indexed well to $\text{La}_4\text{Ni}_3\text{O}_{10}$ with the monoclinic structure. With the pressure increasing, both the lattice parameters and unit-cell volume decrease continuously [Figs. 4(c) and 4(d)]. At around 12 GPa, the (0 2 0) and (2 0 -1) peaks merge into one peak, suggesting the occurrence of structural phase transition [Fig. 4(b) and Supplemental Material Fig. S8 [44]]. The structural analysis revealed that *in situ* high-pressure XRD patterns above 12 GPa can be well described with the tetragonal $I4/mmm$ space group (no. 139). The enthalpy difference relative to $I4/mmm$ phase within 20 GPa is plotted in Fig. 4(e). From 0 to 15 GPa [inset of Fig. 4(e)], we optimized the ions' positions, cell shape, and cell volume, and the enthalpy difference between the $P2_1/a$ phase and $I4/mmm$ phase decreased with pressure. When pressure is higher than 15 GPa, the $P2_1/a$ phase transforms to the $I4/mmm$ phase after structural optimization owing to the sensitivity of the NiO_6 octahedra stacking angles. Hence, we compress the volume with the ions' positions and volume shape fixed based on the optimized $P2_1/a$ structure at 15 GPa for the enthalpy calculations above 15 GPa. The enthalpy of $P2_1/a$ phase is higher than that of the $I4/mmm$ phase after 15.2 GPa, illustrating that the $I4/mmm$ phase is more energetically stable than the $P2_1/a$ phase under high pressure, which is in agreement with *in situ* high-pressure

XRD measurements. Accompanied by the phase transition, the Ni-O-Ni angles change within a layer, which leads to the buckled Ni-O planes becoming more flattened [inset of Fig. 4(e)]. Similar to cuprates, the appearance of a quasi-two-dimensional structure under pressure paves the way for the emergence of high-temperature superconductivity in nickelates.

Based on the above high-pressure characterizations, we construct the temperature-pressure phase diagram of $\text{La}_4\text{Ni}_3\text{O}_{10}$ as shown in Fig. 5(a). In the lower-pressure region, the $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystal exhibits metal-to-metal transition with a density-wave resistance anomaly. As the pressure increases, the resistance anomaly is gradually suppressed. Upon further increasing the pressure, superconductivity appears between 10 and 20 GPa, and the superconducting transition temperature T_c reaches 25 K over 50 GPa. In addition, concomitant with the enhancement of superconductivity in this pressure range, the strange-metal behavior becomes gradually evident above 30 GPa. As can be seen from the phase diagram, the competition between the resistance anomaly and superconductivity resembles the situation in some charge or spin density wave materials. All these observations indicate that high-temperature superconductivity can also be achieved in the pressurized $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals.

D. Theoretical analysis

The pairing mechanism in pressurized $\text{La}_4\text{Ni}_3\text{O}_{10}$ is analyzed as follows: As shown in Fig. 5(b), each unit cell contains three Ni atoms located on three different layers, with each Ni atom hosting two $3d$ orbitals, i.e., $3d_{z^2}$ and $3d_{x^2-y^2}$. For this system, a six-orbital tight-binding (TB) model has been established, with the TB hopping integrals given in Refs. [21,48–51]. Among these hopping integrals, the strongest one is the interlayer hopping between the $3d_{z^2}$ orbitals, i.e., t_{\perp}^z , and the next strongest one is the intralayer hopping between the $3d_{x^2-y^2}$ orbitals, i.e., t_{\parallel}^x . Under the strong Hubbard repulsion, these hopping integrals can induce the strongest antiferromagnetic (AFM) interlayer superexchange J_{\perp}^z and the next strongest AFM intralayer superexchange J_{\parallel}^x . In addition, Hund's rule aligns spins in two orbitals on the same site. Hence, it transfers the nearest-neighbor superexchange interaction from one orbital to the other [36], as shown in Fig. 5(c), but with considerably reduced averaged strength because the transfer takes place only when the two orbitals on the same site are both singly occupied, while in the realistic Ni- $3d^{7.33}$ configuration,

they have only 2/3 possibility to be singly occupied [21]. Similar to the cuprates, these superexchange interactions can serve as effective pairing interactions in pressurized $\text{La}_4\text{Ni}_3\text{O}_{10}$.

Based on the above analysis, a two-orbital t - J model has been established by some of the authors here in Ref. [52]. Then, a systematic slave-boson mean-field approach is engaged to study this model. The obtained results suggest that while both orbitals are near 1/3 filling [21], the $3d_{z^2}$ orbital dominates the pairing because its weak intralayer hopping leads to weak band dispersion and hence enhances the density of state, which enhances the pairing strength. It is also shown in Ref. [52] that the lower T_c in pressurized $\text{La}_4\text{Ni}_3\text{O}_{10}$ than that in pressurized $\text{La}_3\text{Ni}_2\text{O}_7$ originates from the pairing frustration effect caused by its trilayer structure. In bilayer $\text{La}_3\text{Ni}_2\text{O}_7$ driven by the interlayer superexchange interaction, the pairing is mainly between the two layers [33,34,36], experiencing no frustration. However, in trilayer $\text{La}_4\text{Ni}_3\text{O}_{10}$, as shown in Fig. 5(d), when an inner-layer $3d_{z^2}$ electron pairs with an upper outer-layer $3d_{z^2}$ electron driven by J_{\perp}^z , the lower outer-layer $3d_{z^2}$ electron is left single. That electron has to seek pairing

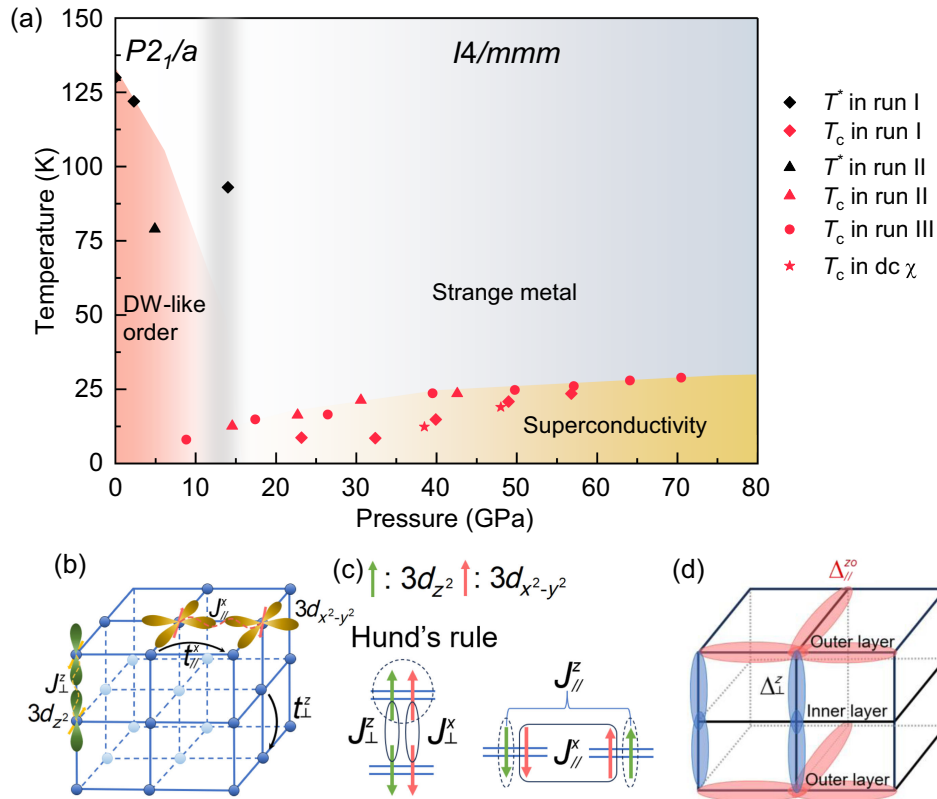


FIG. 5. The phase diagram and theoretical model of pressurized $\text{La}_4\text{Ni}_3\text{O}_{10}$. (a) Phase diagram under pressure. The black symbols represent the density-wave (DW) transition T^* measured at various pressures. The red symbols represent the superconducting transition temperatures T_c determined from the various runs. (b) The relevant orbitals, their couplings, and direct superexchange interactions J_{\perp}^z and J_{\parallel}^x . (c) An illustration of the effective superexchange interactions transferred from one orbital to the other induced by Hund's rule coupling. Note that such a transfer takes place only when both orbitals are singly occupied, while in the realistic Ni- $3d^{7.33}$ configuration, these orbitals have only a 2/3 possibility of being singly occupied. (d) The dominant superconducting pairings.

with neighboring electrons within the same layer. Since the intralayer superexchange interaction J_{\parallel}^z for the $3d_{z^2}$ electron is weak, such a pairing is fragile and easily destroyed by thermal fluctuation. This pairing frustration effect in pressurized $\text{La}_4\text{Ni}_3\text{O}_{10}$ leads to a much lower T_c than that in pressurized $\text{La}_3\text{Ni}_2\text{O}_7$, as shown in Figs. 3(a) and 5(a) in Ref. [52].

III. CONCLUSION

We have successfully grown high-quality single crystals of $\text{La}_4\text{Ni}_3\text{O}_{10}$ and performed a series of transport measurements both at ambient pressure and high pressure. We found pressure-induced superconductivity with a T_c at approximately 25 K. The emergence of superconductivity accompanied by the phase transition from $P2_1/a$ to $I4/mmm$, leads to the buckled Ni-O planes ultimately becoming flattened. Theoretical analysis displays that the AFM superexchange interactions can serve as the effective pairing interaction for the emergence of superconductivity in pressurized $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals. Our results demonstrate that high-temperature superconductivity can be achieved in other members of RP phases, from which we can explore the unconventional superconductivity in nickelates.

Note added. Recently, we became aware that there was another paper by Zhu *et al.* [53]. They reported resistance measurements on $\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals, which is consistent with our data.

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APPENDIX: MATERIALS AND METHODS

1. Material synthesis and characterization

Single crystals of $\text{La}_4\text{Ni}_3\text{O}_{10}$ were grown by the high-pressure optical floating zone technique. Raw materials of La_2O_3 (99.99%) were dried at 900 °C overnight before the reaction and then mixed with an appropriate amount of NiO. The mixture was ground thoroughly, pressed into pellets, and sintered at 1200 °C for two days with several intermediate grindings. The powders were pressed into seed and feed rods with a diameter of about 6 mm and sintered at 1300 °C for two hours. Subsequent single crystal growth was performed in a high-pressure optical floating zone furnace (HKZ-300) with a growth rate of 5 mm/h and 20 bar of oxygen. Single crystals larger than 5 mm in size could be obtained. The powder x-ray diffraction patterns were taken using a Bruker D2 phaser with Cu- K_{α} radiation ($\lambda = 1.5418 \text{ \AA}$) at room temperature. Energy-dispersive x-ray spectroscopy was employed to determine the compositions of single crystals. The high-resolution scanning transmission electron microscopy imaging was taken in JEM-ARM300F GRAND ARM.

2. STM measurements

$\text{La}_4\text{Ni}_3\text{O}_{10}$ single crystals were cleaved at 77 K with a background pressure of 2×10^{-10} Torr and then were immediately transferred into the STM head for measurements. STM experiments were performed within a Unisoku low-temperature STM. All of the STM data were obtained at 4.3 K.

3. ARPES measurements

High-resolution ARPES experiments were conducted at beamline 03U in Shanghai Synchrotron Radiation Facility. The samples were cleaved *in situ* under ultrahigh vacuum below 7×10^{-11} mbar. Data were collected with a Scienta DA30 electron analyzer. The total energy and angular resolutions were set to 20 meV and 0.2° , respectively.

4. Transport properties under ambient and high pressure

Resistivity, heat capacity, and magnetic susceptibility under ambient pressure were measured using the Physical

Property Measurement System (Dynacool, Quantum Design) and Magnetic Property Measurement System (MPMS3, Quantum Design). High-pressure electrical-transport measurements were performed in a DAC with culets of 200 and 300 μm . The cubic boron nitride and epoxy mixture was used as an insulator layer between the BeCu gasket and electrical leads. Four Pt foils were arranged according to the van der Pauw method [54,55]. We used NaCl as a pressure-transmitting medium (PTM) to improve the hydrostatic environment. Ultrasensitive magnetic susceptibility measurements were performed utilizing a bespoke beryllium-copper-alloy miniature DAC fitted with a nonmagnetic rhenium gasket. The measurements were carried out using the Magnetic Property Measurement System. The DAC comprised a pair of diamond anvils, each with a diameter of 300 μm , which confined the sample within a chamber of 240 μm in diameter. This chamber housed a single crystal sample of $\text{La}_4\text{Ni}_3\text{O}_{10}$ measuring approximately 210 μm in diameter and 33 μm in thickness. Neon gas was used as the pressure-transmitting medium to ensure an optimal hydrostatic pressure environment. The superconducting volume fraction calculation method is derived from Ref. [56] (see the Supplemental Material for details [44]). *In situ* high-pressure XRD measurements were performed at the beamline 15U in the Shanghai Synchrotron Radiation Facility ($\lambda = 0.6199 \text{ \AA}$) and mineral oil was used as PTM. The pressure was determined by the ruby luminescence method [57].

5. First-principles calculations

First-principles band structure calculations were performed using the Vienna *ab initio* simulation package [58,59] with a plane-wave basis. For the band structure, the exchange-correlation energy was considered under PBE-type generalized-gradient approximation [60] with spin-orbit coupling included. The Hubbard $U = 4.0 \text{ eV}$ was applied to describe the localized $3d$ orbitals of Ni atoms. The cutoff energy for the plane-wave basis was set to 400 eV. A Γ -centered Monkhorst-Pack k -point mesh of $19 \times 19 \times 5$ was adopted in the self-consistent calculations. For the enthalpy difference calculation, we used the projector-augmented wave [61] approach to describe the core electrons and their effects on valence orbitals and considered the $5s^25p^65d_16s_2$ for La, $3d^84s^2$ for Ni, and $2s^22p^4$ for O. We set the plane-wave kinetic-energy cutoff to 700 eV, and the Brillouin zone was sampled with the special k -mesh generated by the Monkhorst-Pack scheme with a k -point spacing of $2\pi \times 0.025 \text{ \AA}^{-1}$. The convergence tolerance was 10^{-6} eV for total energy, and all forces were converged to be less than 0.003 eV/\AA . At ambient pressure, the calculated volume was less than 1% larger than the experimental data, suggesting the reliability of the calculation.

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