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Common Crystalline and Magnetic Structure of superconducting A_2 Fe₄Se₅ (A=K, Rb, Cs, Tl) single crystals measured using neutron diffraction

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Single crystal neutron diffraction studies on superconductors A_2 Fe₄Se₅, where A = Cs, K, (Tl,Rb) and (Tl,K) ($T_c \sim 30$ K), uncover the same Fe vacancy ordered crystal structure and the same block antiferromagnetic order as in K₂Fe₄Se₅. The Fe order-disorder transition occurs at $T_S = 500-578$ K, and the antiferromagnetic transition at $T_N = 471-559$ K with ordered magnetic moment ~ 3.3 μ_B /Fe at 10 K. Thus, all recently discovered A intercalated iron selenide superconductors share the common crystalline and magnetic structure, which are very different from previous families of Fe-based superconductors, and constitute a distinct new family.

The iron pnictide and chalcogenide superconductors, as exemplified by the ReFeAs(O,F) [1–4], (Ba,K)Fe₂As₂ [5], Li_{1-x} FeAs [6] and Fe_{1+x} Se [7] families of materials, share the Fe square lattice as the common structure feature, which determines electronic states at the Fermi surfaces responsible for high- T_c superconductivity [8]. The $Fe_{1+x}(Se, Te)$ and Li_{1-x} FeAs have the same crystal structure [9–11], in which the excess Fe and Li deficiency at the same crystallographic site serve to maintain close to +2 oxidation state of the Fe ions. Recently, $K_{0.8}Fe_2Se_2$, regarded as K intercalated FeSe in the BaFe₂As₂ structure, has been discovered as a superconductor with T_c above 30 K [12]. Subsequent works replacing K by Rb or Cs [13, 14] have led to more new superconductors of T_c up to ~ 32 K. The heavy departure of Fe from the +2 valency implied by the nominal chemical formulas seem supported by angle-resolved photoemission (ARPES) measurements [15], which would upset theoretical foundation of the prevailing mechanism of previously discovered Fe-based superconductors [16–19].

However, x-ray and neutron diffraction structure refinement studies have shown the correct compositions for the A=K or Cs intercalated iron selenide superconductors as close to $A_{0.8}$ Fe_{1.6}Se₂ [20, 21]. Thus, the new superconductors are not heavily electron doped as previously thought, and Fe has a formal oxidation state close to +2 as in the $Fe_{1+x}(Se,Te)$ superconductors [9, 10]. The $\sim 20\%$ Fe vacancies in $K_{0.8}$ Fe_{1.6}Se₂ are not randomly distributed in the BaFe₂As₂ structure below an orderdisordered transition at $T_S \approx 578$ K [21]. Instead, they form an ordered structure with the Fe1 site almost empty and the Fe2 site fully occupied in the $\sqrt{5} \times \sqrt{5} \times 1$ supercell as shown in Fig. 1 (b)-(c) [21]. At a slightly lower temperature $T_N \approx 559$ K, magnetic moments carried by Fe ions at the Fe2 site develop a block checkerboard antiferromagnetic order. The Fe vacancy and antiferromagnetic order also drastically alter the topology of the Fermi surface according to band structure calculations [22, 23], which would have important physics ramifications and affect the interpretation of the ARPES [15, 24, 25] and optic [26] data.

Also reported are $(Tl,K)Fe_uSe_2$ and $(Tl,Rb)Fe_{1.72}Se_2$ superconductors of similarly high $T_c \sim 30$ K [27, 28]. What distinguishes these superconductors from the Tlless ones is the appearance of the superconducting state at the quantum critical point of an antiferromagnetic insulator [27]. It is imperative then to find out whether the K, Cs or Rb superconductors and the (Tl,K) or (Tl,Rb) superconductors are two different types of new Fe selenide superconductors. We report here singlecrystal neutron diffraction investigation on the Rb, Cs, (Tl,K) and (Tl,Rb) superconductors from about 10 to 580 K that shows the same Fe vacancy and antiferromagnetic order as in the $K_{0.8}Fe_{1.6}Se_2$ superconductor, and the identification of a low T_N in the previous bulk study [27] is incorrect. Hence, superconductivity in all five $A_{0.8}$ Fe_{1.6}Se₂ [A=K, Cs, Rb, (Tl,K) and (Tl,Rb)] superconductors develops in a common crystalline and magnetic structure framework, which is distinctly different from the previous Fe square lattice-based families of iron superconductors [1, 5–7]. Per chemistry convention, $A_{0.8}$ Fe_{1.6}Se₂ is better written as A_2 Fe₄Se₅ (245).

Single crystals of $Rb_2Fe_4Se_5$ ($T_c \approx 32$ K) [14], $Cs_2Fe_4Se_5$ ($T_c \approx 29$ K) [29], (Tl,K)₂Fe₄Se₅ ($T_c \approx 28$ K) [30] and $(Tl,Rb)_2Fe_4Se_5$ ($T_c \approx 32$ K) [27] superconductors were grown using the Bridgeman method at USTC and ZU, and bulk superconductivity is indicated by nearly 100% diamagnetic response reported in previous studies using samples grown with the same recipe[14, 27, 29, 30] and shown in Fig. 1(a). A small piece of Cs₂Fe₄Se₅ single crystal from the same batch was used in the previous x-ray crystal structure refinement study at 295 K [20]. Single crystal neutron diffraction experiments were performed at the High Flux Isotope Reactor (HFIR) of the Oak Ridge National Laboratory. The Wide Angle Neutron Diffractometer (WAND) was used to collect Bragg peaks in the $(h0\ell)$, $(2h, h, \ell)$ and $(3h, h, \ell)$ reciprocal planes at various temperatures. Vertically fo-



FIG. 1. (a) Magnetic susceptibility of A_2 Fe₄Se₅ [A=Rb, Cs, (Tl,K) and (Tl,Rb)] showing the superconducting transition. (b) Common crystal and magnetic structure in the tetragonal I4/m unit cell. The cyan sphere represents Se, yellow sphere the intercalating A, the orange cube the Fe1 vacancy site, and the purple sphere the occupied Fe2 site with +(-)sign indicating up(down) magnetic moment along the *c*-axis. There are 4 formula units per cell. (c) The top Fe-Se layer. The dark and light cyan plaquettes emphasize the four-spin blocks of the up and down spins respectively, which forms a checkerboard nearest-neighbor antiferromagnetic pattern. From one Fe layer to another, magnetic moments are antiferromagnetically coupled. The solid line denotes the I4/m unit cell, one of the two twins, which breaks the high-temperature symmetry of the I4/mmm unit cell (dashed line) at the Fe order-disorder transition at T_S .

cused Ge(113) monochromator was used to produce neutron beam of wavelength $\lambda = 1.460$ Å. The curved, onedimensional ³He position-sensitive detector of 624 anodes covers 125° of the scattering angle. The samples were also investigated using the Four-Circle Diffractometer (FCD) with neutrons of $\lambda = 1.536$ Å at selected temperatures. The order parameters were measured at WAND for the (Tl,Rb), Cs and Rb samples, and at the fixed $E_i = 14.7$ meV triple-axis spectrometer HB1A for the (Tl,K) one. Sample temperature was regulated by high temperature Displex close cycle refrigerator.

In term of the high temperature I4/mmm unit cell of the BaFe₂As₂ structure, the Fe vacancy order is represented by the appearance of the structural superlattice peaks characterized by $\mathbf{Q}_S = (\frac{1}{5}, \frac{3}{5}, 0)$ [20], and the antiferromagnetic order by $\mathbf{Q}_M = (\frac{2}{5}, \frac{1}{5}, 1)$ [21]. Both types of superlattice peaks are accommodated by the $\sqrt{5} \times \sqrt{5} \times 1$ tetragonal I4/m unit cell [21, 31] which we use to label reciprocal space in this paper and in which $\mathbf{Q}_S = (110)$ and $\mathbf{Q}_M = (101)$. Fig. 2 shows single-crystal diffraction pattern in the $(h0\ell)$ plane at 295 K. Structural Bragg peaks $\{h0\ell\}$ (h = 2, 4, 6, ... and $\ell = 0,$ 2, 4, ...) due to the Fe vacancy order as well as magnetic Bragg peaks $\{h0\ell\}$ (h = 1, 3, 7, ... and $\ell = 1, 3, 5,$...) caused by the block checkerboard antiferromagnetic order [21] are clearly visible. These "superlattice" peaks

TABLE I. Physics properties of the A_2 Fe₄Se₅ superconductors. Lattice parameters at 295 K, ordered magnetic moment of Fe at 295 K (*m*) and at ~10 K (*M*) are provided together with T_c , T_N and Fe vacancy order-disorder transition T_S .

Α	Κ	Rb	\mathbf{Cs}	Tl,K	Tl,Rb
a (\mathring{A})	8.7306(1)	8.788(5)	8.865(5)	8.645(6)	8.683(5)
c (Å)	14.1128(4)	14.597(2)	15.289(3)	14.061(3)	14.388(5)
$m~(\mu_B)$	2.76(8)	2.95(9)	2.9(1)	2.61(9)	2.7(1)
$M(\mu_B)$	3.31(2)	3.3(1)	3.4(2)	3.2(1)	3.2(1)
T_c (K)	32	32	29	28	32
T_N (K)	559(2)	502(2)	471(4)	506(1)	511(1)
T_S (K)	578(2)	515(2)	500(1)	533(2)	512(4)

are forbidden in the $BaFe_2As_2$ structure.

Like in previous x-ray and electron single crystal diffraction studies [20, 31, 32], twins with the same caxis exist in all crystals examined in this neutron single crystal diffraction study. They show up as those "superlattice" peaks at non-integer h indices in Fig. 2. For example, the $(-1.41, 0, \ell)$ with ℓ even in Fig. 2(a) can be identified as the Fe vacancy peaks (11ℓ) from a twinned domain. During the data reduction, care was paid to resolve the twin distribution. Data collected using WAND and FCD were consistent with each other. Data analysis was conducted using the Fullprof program suite. The relative occupancy of Tl and Rb (Tl and K) was fixed by that from the energy dispersive x-ray (EDX) spectroscopic measurements since our neutron diffraction refinements did not have enough sensibility for an independent determination. Fixing or floating the ratio does not affect appreciably the value of refined magnetic moment. The total number of independent Bragg reflections used in the refinements were 155, 131, 190 and 132 for the A=Rb, Cs, (Tl,K) and (Tl,Rb) samples, respectively, and the RF-factors were 4.8, 5.6, 4.5 and 7.7%, respectively. A summary of results of the four superconductors investigated in this study is listed in Table I, together with those of $K_2Fe_4Se_5$ from previous study [21].

Two separated phase transitions due to antiferromagnetic order and the iron vacancy order have been discovered in neutron diffraction study on $K_2Fe_4Se_5$ [21], but only magnetic transition has been reported previously for the Cs compound in a combined μ SR and thermometry study [33]. In Fig. 3(a), integrated intensity of magnetic Bragg peak (101) and structural Bragg peak (118) of $Cs_2Fe_4Se_5$, simultaneously measured on WAND, are shown as a function of temperature. Similar to the case of the K intercalated superconductor, there are two separated magnetic and structural transitions at $T_N = 471(4)$ K and $T_S = 500(1)$ K, respectively. The order parameters at \mathbf{Q}_S and \mathbf{Q}_M look very similar to those of K_2 Fe₄Se₅ [21], albeit now that magnetic and structural transition temperatures of T_N and T_S are lower [Table I and Fig. 3]. The T_N reported here agrees with $T_N \approx 477$ K from



FIG. 2. Single crystal neutron diffraction pattern of (a) $Cs_2Fe_4Se_5$, (b) $Rb_2Fe_4Se_5$, (c) $(Tl,Rb)_2Fe_4Se_5$ and (d) $(Tl,K)_2Fe_4Se_5$ measured in the (h0 ℓ) plane of the I4/m unit cell at T=295 K. In addition to Bragg peaks of the high temperature I4/mmmstructure, two new types of contributions to Bragg reflection appears: 1) structural {200}, {202}, {400}, {402}, etc. caused by the Fe vacancy order, and 2) magnetic {103}, {301}, {303}, {701}, etc. caused by antiferromagnetic order of the Fe moments.

the previous μ SR study [33], which also reports a microscopic coexistence of superconductivity and an unidentified antiferromagnetic order in a nominal Cs_{0.8}Fe₂Se_{1.96} ($T_c \approx 28.5$ K) superconductor. The T_S agrees with that from an x-ray study appearing after we finished the measurement [34]. Different from previous Fe-based superconductors where antiferromagnetic order coexisting with superconductivity can be regarded as relic from the parent antiferromagnet [35], antiferromagnetic order in 245 superconductors is very strong. In addition to the high T_N , all the magnetic moments listed in Table I at ~10 K are higher than previous record 2 μ_B /Fe found in nonsuperconducting *parent* compounds [9].

The structural peak at \mathbf{Q}_S , which represents the breaking of the high temperature I4/mmm symmetry when the vacant (orange cube in Fig. 1) and occupied (purple sphere in Fig. 1) Fe sites segregate below T_S , saturates rapidly to its low temperature value [Fig. 3(a)]. Therefore, below the room temperature where relevant physics processes relating to the superconductivity occur, we are dealing with a well ordered crystal structure in the I4/msymmetry. Structurally, and consequently electronically, using the BaFe₂As₂ family of superconductor materials as a model is too remote to the reality. The modification of electronic structure by the new I4/m crystal symmetry and by the block antiferromagnetic order has been vividly and dramatically demonstrated in first-principle band structure calculations [22, 23]. Future investigations on the 245 superconductors should be based on the correct I4/m crystal structure.

Comparing to the temperature dependence of the structural peak, the magnetic Bragg peak of $Cs_2Fe_4Se_5$ show a gradual rise below T_N [Fig. 3(a)], as previously found for $K_2Fe_4Se_5$ [21]. The T_N is lower than T_S as ex-



FIG. 3. (a) Magnetic (103) and structural (118) Bragg peaks vs. temperature, serving as order parameters for the antiferromagnetic and Fe vacancy order-disorder transitions, respectively, in Cs₂Fe₄Se₅. (b) Normalized magnetic Bragg intensity representing the squared magnetic moment as a function of temperature for A_2 Fe₄Se₅ superconductors. Inset: Magnetic (101) peak of (Tl,Rb)₂Fe₄Se₅ showing that its intensity saturates when T_c is approached.

pected, since the long-range magnetic order builds on the ordered Fe pattern. This is reminiscent of the antiferromagnetic transition in V_2O_3 albeit the underlying transi-

tion is to an orbital occupation order [37]. Fig. 3(b) illustrates the magnetic order parameters of all five A₂Fe₂Se₅ samples with T_N s listed on Table I. Like in the K compound, the linear dependence of magnetic Bragg intensity is interrupted by a flat plateau when T_c is approached. For example, the inset of Fig. 3(b) shows a temperature scan measured with much better statistics using tripleaxis spectrometer for the (Tl, Rb)₂Fe₄Se₅ sample, suggesting interaction between the antiferromagnetic order and superconductivity as previously discovered in the case of heavy fermion superconductors [36].

Since the discovery of the nominal $K_{0.8}Fe_2Se_2$ superconductor [12], there had been chaotic confusion about sample composition of the new iron selenide superconductors due to the lack of reliable structure determination work. The Fe-Se square lattice was initially regarded as intact as in the previous families of iron-based superconductors. Analogue to the heavily electron doped iron pnictide superconductors was advocated. Theories based on both the AFe_2Se_2 of the $BaFe_2As_2$ structure and $AFe_{1.5}Se_2$ of a $\sqrt{2} \times 2\sqrt{2} \times 1$ supercell structure framework were advanced for the new superconductors [16, 38]. The superconductivity in the (Tl.K) system appeared to occur at a magnetic quantum critical point [27] and supported a doped Mott insulator scenario [39, 40] similar to the case of cuprate superconductors. While electronic properties are indeed very different from previous ironbased superconductors as probed in ARPES [15, 24, 25], optic [26], NMR [41–43], μ SR [33] and Raman scattering [44] studies, an appropriate understanding cannot be achieved without knowing the superconducting composition, crystal structure and magnetic order determined in neutron and x-ray diffraction structural studies.

In summary, our systematic neutron diffraction works show that Cs, Rb, (Tl,Rb) and (Tl,K) intercalated Fe selenide superconductors are members of the common family of the A_2 Fe₄Se₅ superconductors which share the same Fe vacancy- and antiferromagnetic order. Not only the crystal structure of the A_2 Fe₄Se₅ superconductors in the $\sqrt{5} \times \sqrt{5} \times 1 I4/m$ unit cell is very different from that in previous iron-based high- T_c superconductors, leading to drastically different Fermi surface topology, the coexistence of superconductivity with the very high T_N and large moment antiferromagnetic order is unprecedented. Although the same Fe ions are involved, given the drastically different crystalline and magnetic structures, the A_2 Fe₄Se₅ superconductors are unlikely to share the same physics mechanism as in previously discovered families of iron-based superconductors. Future search for new high- T_c superconductors has a wider playground including vacancy decorated lattice.

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