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Avoided Quantum Criticality and Magnetoelastic Coupling in BaFe$_{2-x}$Ni$_x$As$_2$

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We study the structural and magnetic orders in electron-doped BaFe$_{2-x}$Ni$_x$As$_2$ by high-resolution synchrotron x-ray and neutron scatterings. Upon Ni doping $x$, the nearly simultaneous tetragonal-to-orthorhombic structural ($T_s$) and antiferromagnetic ($T_N$) phase transitions in BaFe$_2$As$_2$ are gradually suppressed and separated, resulting in $T_s > T_N$ with increasing $x$, as was previously observed. However, the temperature separation between $T_s$ and $T_N$ decreases with increasing $x$ for $x \geq 0.065$, tending toward a quantum bicritical point near optimal superconductivity at $x = 0.1$. The zero-temperature transition is preempted by the formation of a secondary incommensurate magnetic phase in the region $0.088 \leq x \leq 0.104$, resulting in a finite value of $T_N = T_s + 10$ K above the superconducting dome around $x = 0.1$. Our results imply an avoided quantum critical point, which is expected to strongly influence the properties of both the normal and superconducting states.

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A determination of the structural and magnetic phase diagram in correlated electron materials is important for understanding their underlying electronic excitations. In the iron pnictides, superconductivity arises at the border of both antiferromagnetic (AF) and structural orders [1–5]. This motivates the exploration of quantum critical points, where the transition temperatures for such orders are continuously suppressed to zero by a nonthermal control parameter. For the iron pnictide superconductors derived from electron or hole doping of their parent compounds, the most heavily studied materials are probably the electron-doped BaFe$_{2-x}$Co$_x$As$_2$ (where $T = \text{Co, Ni}$) because of the availability of high-quality single crystals [6–18]. In the undoped state, BaFe$_2$As$_2$ exhibits a tetragonal-to-orthorhombic structural transition at temperature $T_s$, and an AF phase transitions below nearly the same temperature $T_N = T_s \approx 138$ K [3,4]. Upon electron doping of BaFe$_2$As$_2$ via partially replacing Fe by Co or Ni, various experiments, including transport [8,9], neutron [11–16], and high-resolution x-ray scattering [4,18], reveal that the structural ($T_s$) and magnetic ($T_N$) phase transition temperatures in BaFe$_{2-x}$Co$_x$As$_2$ gradually decrease and separate with increasing $x$ and have $T_s > T_N$ for all doping levels. In the initial x-ray [10] and neutron [11] scattering work on BaFe$_{2-x}$Co$_x$As$_2$, it was suggested that the separated $T_s$ and $T_N$ smoothly extend into the superconducting dome, resulting in distinct structural and magnetic quantum critical points at different $x$. Subsequent x-ray [18] and neutron [12–14] scattering experiments on superconducting BaFe$_{2-x}$Co$_x$As$_2$ samples with coexisting AF order revealed that superconductivity actually competes with the static AF order and lattice orthorhombicity. As a consequence, the smoothly decreasing $T_s$ and $T_N$ are reported to bend back below $T_s$, and the orthorhombic structure above $T_s$ for an optimally doped sample evolves back to a tetragonal structure well below $T_c$ (termed the “reentrant” tetragonal phase) [18].

Although previous neutron [11–13] and x-ray diffraction [18] experiments have established the magnetic and structural phase transitions in BaFe$_{2-x}$Co$_x$As$_2$, similar measurements have not been carried out on BaFe$_{2-x}$Ni$_x$As$_2$. In this Letter, we describe neutron and x-ray scattering studies of structural and magnetic phase transitions in BaFe$_{2-x}$Ni$_x$As$_2$, focusing on materials near optimal superconductivity [Fig. 1(a)]. While neutron scattering experiments on BaFe$_{2-x}$Co$_x$As$_2$ revealed a commensurate-to-incommensurate AF phase transition near optimal superconductivity [15–17], much remains unknown about the temperature and doping evolution of the orthorhombic lattice distortion for samples with an incommensurate AF order. Here, we find that $T_s > T_N$ for samples with commensurate AF order ($x \approx 0.065$), similar to the earlier results on BaFe$_{2-x}$Co$_x$As$_2$ [11–13,18]. However, $T_s$ and $T_N$ tend to reconverge for larger values of $x$: $T_s - T_N$...
pnictides that was predicted by theory [23] and observed quantum critical point of the isoelectronically tuned iron [22] measurements, as well as its connection with the BaFe$_2$As$_2$ by extensive experiments [24,25].

from the NMR [19,20], thermoelectric [21], and ultrasonic critical point in the carrier-doped iron pnictides, as inferred important to clarifying the nature of the purported quantum interest is illustrated schematically in Fig. 1(b). Our results are ordered moment [16]. The resulting overall phase diagram short-range incommensurate AF order with a very small critical point at $T_N$ is indicated by a secondary quantum bicritical point.

and temperature dependence of the orthorhombic lattice distortion order parameter $\delta = (a_0 - b_0)/(a_0 + b_0)$. The inset shows the expanded view of $T_N$, $T_N$, and $T_c$ and temperature dependence of the orthorhombic phase are clearly marked. The inset shows the presence of a single Gaussian structural peak. (b) Schematic theoretical phase diagram for an avoided quantum bicritical point.

decreases for $x > 0.065$. This implicates a quantum bicritical point at $T = 0$, which is interrupted by a secondary short-range incommensurate AF order with a very small ordered moment [16]. The resulting overall phase diagram is illustrated schematically in Fig. 1(b). Our results are important to clarifying the nature of the purported quantum critical point in the carrier-doped iron pnictides, as inferred from the NMR [19,20], thermoelectric [21], and ultrasonic [22] measurements, as well as its connection with the quantum critical point of the isoelectronically tuned iron pnictides that was predicted by theory [23] and observed by extensive experiments [24,25].

We have carried out neutron scattering experiments on BaFe$_{2-x}$Ni$_x$As$_2$ with $x = 0.085$, 0.092, 0.096, 0.1, 0.104 and 0.108 using the RITA-II cold neutron triple-axis spectrometer at the Paul-Scherrer Institute; the HB-1A thermal triple-axis spectrometer at the High-Flux Isotope Reactor, Oak Ridge National Laboratory; and the C5 triple-axis spectrometer at the Canadian Neutron Beam Centre, Chalk River Laboratories [26]. We have also performed high-resolution synchrotron x-ray diffraction experiments on identical BaFe$_{2-x}$Ni$_x$As$_2$ samples using beam line X22C at the National Synchrotron Light Source, Brookhaven National Laboratory. The details of the experimental procedure are given in the Supplemental Material [27]. Although neutron scattering probes the bulk sample, whereas the length scale for x-ray diffraction is typically about $\sim 5$ microns [28], both techniques are measuring the intrinsic properties of these materials.

We first describe the determination of the Ne`el temperatures for BaFe$_{2-x}$Ni$_x$As$_2$ using neutron scattering. Figure 2(a) shows transverse scans along the [1, K, 3] direction at different temperatures for the $x = 0.085$ sample. Consistent with earlier results [16], a well-defined commensurate AF order appears below 44 K. Figure 2(b) shows temperature dependence of the magnetic order parameter. Again, consistent with earlier results [15–17], the AF order appears approximately below $T_N = 44 \pm 5$ K and is suppressed from the onset of $T_c$. Figure 2(c) plots similar data for $x = 0.092$ and 0.096, showing $T_N = 40 \pm 5$ and 32 $\pm 5$ K, respectively [16]. In the previous work on optimally

FIG. 1 (color online). (a) Electronic phase diagram of BaFe$_{2-x}$Ni$_x$As$_2$ as a function of Ni doping $x$ as determined from our neutron and x-ray scattering experiments. The PM Tet, PM Ort, AF Ort, and IC Ort are paramagnetic tetragonal, paramagnetic orthorhombic, commensurate AF orthorhombic, and incommensurate AF orthorhombic phases, respectively. The AF Ort, IC Ort, and PM Tet structures in the superconducting (SC) phase are clearly marked. The inset shows the expanded view of $T_N$, $T_N$, and $T_c$ and temperature dependence of the orthorhombic

FIG. 2 (color online). (a) Transverse scans along the [1, K, 3] direction at different temperatures for BaFe$_{2-x}$Ni$_x$As$_2$ with $x = 0.085$. The magnetic scattering of each temperature was obtained for an avoided quantum bicritical point.

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electron-doped BaFe$_{1.9}$Ni$_{0.1}$As$_2$ [29], it was suggested, based on cold neutron data on mosaic crystals ($\sim 0.6$ g) counting 1 min/point, that there is no measurable static AF order. Our new measurements on the $x = 0.1$ sample ($\sim 0.34$ g) with much longer counting time (30 min/point on HB-1A) reveal a weak static AF order with magnetic scattering 5 times smaller than that of $x = 0.096$ [Figs. 2(c) and 2(d)]. Similar measurements on $x = 0.104$ also show the presence of a weak static AF order, which is 50% smaller than that of the $x = 0.1$ sample. In spite of their small moments, the temperature dependence of the magnetic order parameters for both samples indicates that their Néel temperatures are essentially unchanged at $T_N = 30 \pm 5$ K [Fig. 2(d)]. Finally, we find no evidence of static AF order for a $x = 0.108$ sample ($\sim 0.5$ g) by counting 40 min/point on CS [Fig. 2(d)].

In order to compare the onset of orthorhombicity with antiferromagnetism, high-resolution x-ray scattering measurements were performed on the samples identical to those used for neutron scattering. In all cases, we carried out longitudinal scans along the $[H, 0, 12]$ direction. Figure 3(a) shows the outcome for $x = 0.085$, which has a superconducting $T_c = 16.5$ K. At $T = 58$ K, a temperature well above $T_c$, we see a single instrumentation resolution-limited peak, consistent with a tetragonal lattice. On cooling to $T = 45, 30$, and 17 K, the single peak splits into two peaks with increasing peak separations as temperature decreases down to $T_c$. Upon further cooling below $T_c$, the peak separations become smaller, as if the system turns back toward the tetragonal structure [18]. Figure 3(b) shows similar temperature-dependent scans for $x = 0.096$. Although the split peaks appear to become a single peak at $T = 10.5$ K, its width is still larger than that in the tetragonal phase ($T = 34$ K), suggesting that the nearly optimal superconductor has an orthorhombic lattice distortion at $T = 10.5$ K. To see how such orthorhombic lattice distortion evolves at lower temperatures, we carried out additional measurements using a cryostat capable of going down to 2 K. The longitudinal $[H, 0, 12]$ scans in Fig. 3(c) show broad peaks at temperatures below 10 K, suggesting the presence of an orthorhombic lattice structure even at 2 K.

To quantitatively analyze the temperature dependence of the orthorhombic lattice distortion, we define lattice orthorhombicity $\delta = (a_o - b_o)/(a_o + b_o)$, where $a_o$ and $b_o$ are lattice parameters of the orthorhombic unit cell [18]. Figure 4(a) shows the temperature dependence of $\delta$ for BaFe$_{2-x}$Ni$_x$As$_2$ with $x = 0.075, 0.085, 0.092, 0.096$, and 0.1. Figures 4(b) and 4(c) compare the ordered moment squared $M^2$ with the lattice orthorhombicity $\delta$, and their similar temperature dependence suggests a strong magnetoelastic coupling.

The optimally doped $x = 0.1$ sample ($T_c = 20.2$ K) deserves special attention. Its temperature-dependent $[H, 0, 12]$ scans are shown in Fig. 3(d). Although we can no longer see the double peaks, we observe a peak broadening that does not disappear at low temperatures. We therefore used the FWHM of the peak in order to determine the lattice orthorhombicity $\delta$, similar to the analysis of BaFe$_{2-x}$Co$_x$As$_2$ by Nandi et al. [18]. The deduced temperature dependence of $\delta$ is shown in Fig. 4(a) with red squares and appears to have a sharp cusp near the superconducting $T_c$. We conjecture that this cusp occurs because the electron-lattice coupling results in a lattice response to the superconducting fluctuations near $T_c$. At the lowest temperature measured $T = 11$ K, the value of $\delta$ is too small ($2 \times 10^{-5}$) in order to unambiguously claim the orthorhombicity. However, taken together with magnetization squared for incommensurate AF order [see Fig. 4(c), which has a similar temperature dependence], we conclude that a weak static AF order likely coexists with orthorhombic lattice distortion in the optimally superconducting BaFe$_{2-x}$Ni$_x$As$_2$, different from the reentrant tetragonal transition seen in BaFe$_{2-x}$Co$_x$As$_2$ [18].

Figure 4(d) shows the Ni-doping dependence of $\delta$ and the ordered moment squared $M^2$, while Fig. 4(e) compares the doping dependence of $\delta$ in BaFe$_{2-x}$Ni$_x$As$_2$ and in the previously reported BaFe$_{2-x}$Co$_x$As$_2$ [18]. The essentially continuous suppression of both $M^2$ and $\delta$ near $x = 0.1$ provides further evidence for an extrapolated quantum critical point. For the magnetic ordering, this represents
except that here the lattice quantum field $\phi$ is endowed with its own dynamics and undergoes Landau damping $\Gamma_{\phi}$, making it inherently quantum critical with the dynamic exponent $z = 3$. In two spatial dimensions, $d + z = 5$ for the $\phi$ field and $d + z = 4$ for the $M$ fields. Because they are above or at the upper critical dimension, a quantum bi-critical point for both orders is expected in the presence of the magnetoelastic coupling $\eta$. This is similar to the result of the $O(3) \times \mathbb{Z}_2$ model [23] and is indicated schematically in Fig. 1(b). Indeed, as noted above, our measurements find that $T_N$ and $T_s$ get closer to each other as the quantum critical point is approached [see Figs. 1(a) and 4(f)] and the two order parameters disappear at the same point [Fig. 4(d)]. However, the appearance of an emergent incommensurate magnetism at $x = 0.088$ severely reduces the scattering rate $\gamma$ and $\Gamma_{\phi}$ (in addition to modifying other parameters of the effective theory), thereby eliminating the quantum critical point. A quantum critical point preempted by an emergent order is often referred to as “avoided” quantum criticality [31–33].

From direct measurements of the order parameters for both the AF and structural transitions, our results provide a solid basis for quantum criticality in carrier-doped iron pnictides, which has so far been indirectly deduced from the temperature dependences of magnetic, transport, or acoustic properties [19–22]. In addition, because the primary AF order in the electron-doped iron pnictides discussed here is commensurate, our results suggest that the quantum critical point arising under the carrier doping is surprisingly similar to that induced by isoelectronic doping [23–25]; the main distinction of the carrier doping is to introduce a secondary incommensurate order. This reveals an important universality of the underlying physics for the iron pnictides under carrier and isoelectronic dopings.

Summarizing the results presented in Figs. 2–4, we show in Fig. 1(a) the refined phase diagram of BaFe$_{2-x}$Ni$_x$As$_2$, in agreement with the theoretically expected one [Fig. 1(b)]. While the phase diagram is mostly consistent with the earlier work on BaFe$_{2-x}$Co$_x$As$_2$ at low electron-doping levels [18], our key new finding is that when $x$ approaches optimal doping, the magnetic and structural transition temperatures converge to the purported quantum bicritical point, with both order parameters disappearing near $x = 0.1$ [Fig. 4(d)] as a result of magnetoelastic coupling. However, the emergent short-range incommensurate magnetism helps the system avoid the quantum critical fate, resulting in an apparent saturation of $T_s \sim T_N \approx 30$K above the superconducting $T_c$ near optimal doping $x = 0.1$, as shown in Fig. 1(a). These results elucidate the quantum criticality in the carrier-doped iron pnictides and its connection with that of the isoelectronically doped counterparts, and reveal a rich theoretical picture that should be further explored in future work.

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