

COMMENTARY

A different branch of the high T_c family?D. J. Scalapino^{a,1}

In PNAS, Li et al. (1) suggest that $\text{Ba}_2\text{CuO}_{4-y}$ is a member of a different branch of high- T_c cuprate superconducting materials. This branch is characterized as heavily overdoped with an exceptionally short Cu apical O spacing and O vacancies that are located in the CuO_2 planes. These characteristics, illustrated in Fig. 1, differ in significant ways from those of the traditional cuprate superconducting materials (2).

For example, consider $\text{La}_{2-x}\text{Sr}_x\text{CuO}_4$ (LSCO), which has the same K_2NiF_4 structure as $\text{Ba}_2\text{CuO}_{4-y}$. Using a valence count with La^{+2} , Sr^{+3} , and O^{-2} , the number of holes in the Cu 3d shell of LSCO is $1+x$. For $x=0$, LCO is antiferromagnetic. As Sr is added the antiferromagnetic Néel temperature decreases as shown in Fig. 1A and superconductivity onsets with T_c peaking at a small doping $x \sim 0.15$. At larger doping, as the system moves further away from the antiferromagnetic Mott–Hubbard phase, T_c decreases and vanishes for $x > 0.25$. For $\text{Ba}_2\text{CuO}_{4-y}$, the valence of Ba is +2 so that the hole doping $x = 2(1-y)$. Thus, the $y=0.8$ sample Li et al. (1) discuss is heavily overdoped with $x=0.4$. Nevertheless, the T_c of $\text{Ba}_2\text{CuO}_{4-y}$ is approximately 30 K higher than that of optimally doped LSCO.

Like the previously reported highly overdoped cuprate materials (3–8) $\text{Sr}_2\text{CuO}_{4-y}$, $(\text{Sr},\text{Ba})_2\text{CuO}_{4-y}$, and $\text{Cu}_{0.75}\text{Mo}_{0.25}\text{Sr}_2\text{YCu}_2\text{O}_{7.54}$, $\text{Ba}_2\text{CuO}_{4-y}$ is synthesized under high pressure and high temperature in the presence of a strong oxidizing agent. While the crystal has the K_2NiF_4 structure illustrated in the center of Fig. 1B, the CuO_6 octahedron structure is highly compressed with a much shorter Cu–O apical distance (1.86 Å) than the traditional cuprates (2.42 Å for La_2CuO_4). This is schematically illustrated at the left and right in Fig. 1B. The shortening of the Cu–O apical distance raises the energy of the $3d_{3z^2-r^2}$ orbital so that in $\text{Ba}_2\text{CuO}_{4-y}$ the Cu states near the Fermi level have both $3d_{3z^2-r^2}$ and $3d_{x^2-y^2}$ orbital character. As the authors note, a shortened apical Cu–O distance and the admixing of $3d_{3z^2-r^2}$ orbital weight in the states near the Fermi energy are found to reduce T_c in the traditional cuprate superconductors (9).

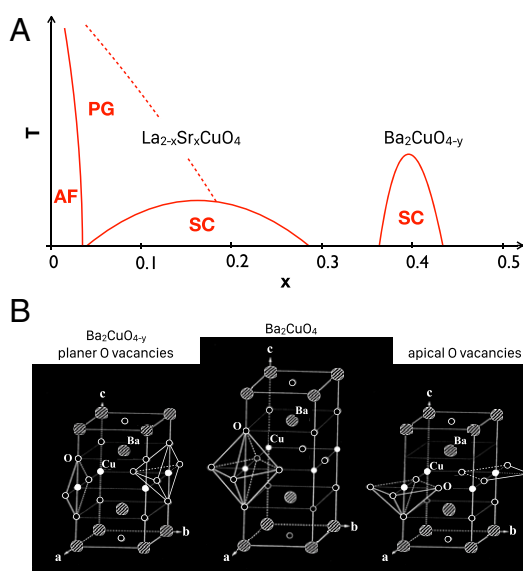


Fig. 1. (A) A schematic cuprate phase diagram as a function of hole doping x . On the left at low doping are the familiar antiferromagnetic (AF), pseudogap (PG), and d -wave superconducting (SC) regions of $\text{La}_{2-x}\text{Sr}_x\text{CuO}_4$. On the right is the heavily overdoped $x \sim 0.4$ region of superconductivity Li et al. (1) discuss for $\text{Ba}_2\text{CuO}_{4-y}$. (B) The center panel shows a fictitious Ba_2CuO_4 compound with the same K_2NiF_4 structure as La_2CuO_4 . The structures shown on the left and right have the compressed c -axis structure of $\text{Ba}_2\text{CuO}_{4-y}$. In this case the hole doping is controlled by the O vacancies. For the structure on the right, the O vacancies are on the apical sites and the CuO_4 corner-shared CuO_2 sheets remain intact. However, according to Li et al. (1) in $\text{Ba}_2\text{CuO}_{4-y}$ the O vacancies are in the plane. In this case, as shown by the structure on the left, the CuO_2 sheets are destroyed and randomly oriented Cu–O chain segments are likely formed.

Finally, the authors note at the end of the legend for figure 3 in ref. 1 that while the exact positions of the O vacancies are not known at present, they are in the CuO_2 planes. This agrees with the conclusions of Geballe and Marezio (5) regarding the O vacancies in $\text{Sr}_2\text{CuO}_{4-y}$. A number of earlier studies Li et al. (1) cite assume that the oxygen vacancies were at the apical O sites. In this

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