

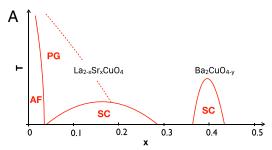
A different branch of the high T_c family?

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In PNAS, Li et al. (1) suggest that $\mathrm{Ba_2CuO_{4-y}}$ 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 $\mathrm{CuO_2}$ planes. These characteristics, illustrated in Fig. 1, differ in significant ways from those of the traditional cuprate superconducting materials (2).

For example, consider $La_{2-x}Sr_xCuO_4$ (LSCO), which has the same K_2NiF_4 structure as Ba_2CuO_{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 Ba_2CuO_{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 Ba_2CuO_{4-y} is approximately 30 K higher than that of optimally doped LSCO.

Like the previously reported highly overdoped cuprate materials (3–8) Sr_2CuO_{4-y} , $(Sr,Ba)_2CuO_{4-y}$, and $Cu_{0.75}Mo_{0.25}Sr_2YCu_2O_{7.54}$, Ba_2CuO_{4-y} is synthesized under high pressure and high temperature in the presence of a strong oxidizing agent. While the crystal has the K₂NiF₄ structure illustrated in the center of Fig. 1B, the CuO₆ octahedron structure is highly compressed with a much shorter Cu-O apical distance (1.86 Å) than the traditional cuprates (2.42 Å for La₂CuO₄). 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 Ba₂CuO_{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).



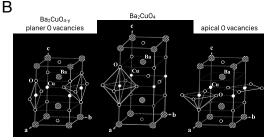


Fig. 1. (A) A schematic cuprate phase diagram as a function of hole doping x. On the left at low doping are the familiar antiferromagnetic (AF), pseudogap (PG), and d-wave superconducting (SC) regions of La2-xSrxCuO4. On the right is the heavily overdoped $x \sim 0.4$ region of superconductivity Li et al. (1) discuss for Ba₂CuO_{4-v}. (B) The center panel shows a fictitious Ba₂CuO₄ compound with the same K₂NiF₄ structure as La₂CuO₄. The structures shown on the left and right have the compressed c-axis structure of Ba_2CuO_{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 CuO₄ corner-shared CuO₂ sheets remain intact. However, according to Li et al. (1) in Ba_2CuO_{4-y} the O vacancies are in the plane. In this case, as shown by the structure on the left, the CuO₂ 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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case, while there would be CuO_5 pyramidal structures and square fourfold coordinated CuO_4 along with the CuO_6 octahedral units, the CuO_2 planes would survive intact as illustrated in the right-hand structure of Fig. 1B. However, if the oxygen vacancies are in the CuO_2 planes one will have Cu–O chain structures rather than the CuO_2 planes as indicated in the left-hand structure of Fig. 1B. The 2D CuO_2 planes, consisting of corner-shared CuO_4 units, are considered the key structural elements of the traditional cuprate superconductors. They are the defining characteristic of these superconductors and imperfections in these layers are known to reduce T_c .

As noted, Ba_2CuO_{4-y} is made under high pressure and temperature resulting in polycrystalline samples and the precise

location of the oxygen vacancies in the CuO_2 plane remains open. Nevertheless, the remarkably high T_c of this highly overdoped cuprate with its short Cu apical O separation and its O vacancies in the CuO_2 plane suggest that it is a member of a different branch of high- T_c cuprate materials, which challenges the basic tenants of many high- T_c theories.

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