Comment on "Electronic Raman Scattering in High- T_c Superconductors: A Probe of $d_{x^2-y^2}$ Pairing"

Devereaux et al. [1] calculated electronic Raman scattering in a superconductor [2,3] for a gap function of $d_{x^2-y^2}$ symmetry. Strong anisotropies in the spectra of the superconducting gap excitations were predicted. "Quantitative agreement" with Raman data for Bi-2212 in B_{1g} and B_{2g} polarization configurations is claimed. Since the data are taken as evidence for $d_{x^2-y^2}$ pairing we point out flaws in [1] which invalidate its conclusions.

While the data presented for Bi-2212 in Figs. 2 and 3 of [1] agree with the calculation, no data for the A_{1g} component are presented. This is attributed in Ref. [13] of [1] to the impossibility of obtaining A_{1q} data from experiments involving a single polarization configuration. However, provided A_{1g} is large, its extraction from measurements for two polarization configurations is standard procedure of Raman spectroscopy. A_{1g} is large in literature data for Bi-2122 [4], Y-123 [5], and Tl-2223 [6]. In Fig. 1 we show that the A_{1g} spectrum of a single domain Y-123 crystal, obtained by us from the xx , yy , and $x'y'$ (i.e., $x^2 - y^2$) polarization configurations is stronger than (x, y)) the B_{1q} and B_{2q} spectra. Its shape is the same as that of the measured xx and yy spectra.

Within the standard theory of scattering by normal carriers [7] in the range of Fig. 1 only unscreened scattering, related to mass fluctuations around and among the various sheets of the Fermi surface, should be observable well below the plasma frequency [8]. It has been shown [9] that the mass fiuctuations obtained from the band structure [10] without resonance factors, account well for the normal state data in single domain Y-123. The A_{1g} mass fluctuations are actually several times larger than the average mass since the latter reverses sign around the Fermi surface (see Fig. 30 of $[11]$). Since the mass fluctuations also determine the γ of Eq. (4) of [1] we expect the A_{1g} scattering to be mostly unscreened while the curve in Fig. 1 of [1] corresponds to screened scattering.

The unscreened A_{1g} scattering, in particular that part due to fluctuations among various sheets of the Fermi surface, should have a maximum at $\omega = 2\Delta_0$, since it samples equally all values of $\Delta(\mathbf{k}, T)$ which have a stationary point for $\Delta = \Delta_0$. This has been confirmed in Fig. 3 of [12] and contradicts the experiments of Fig. 1: A_{1g} peaks well below $2\Delta_0$ if, as required in [1], $2\Delta_0$ is given by the peak in B_{1g} . This results from the fact that the only spectrum which should peak well below $2\Delta_0$ is that for xy polarization, which has a well defined orthorhombic symmetry $(\neq A_g)$ and samples the Fermi surface regions around zero gap.

Last, but not least, in Raman spectroscopy, like in many measurements which are not sensitive to the phase

FIG. 1. A_{1g} , B_{1g} , and B_{2g} Raman response of a single domain Y-123 crystal determined from the xx , yy, and $x'y'$ (i.e., x^2-y^2) spectra displayed on the same scale. To compare with A_{1g} calculations of [1], which assume tetragonal symmetry, we have averaged xx and yy polarizations. This is not essential to our conclusions [9]. Phonons have been removed from the data.

of the order parameter, there is no way to distinguish between $d_{x^2-y^2}$ and a strongly anisotropic s-wave gap (e.g., $|x^2-y^2|^2$).

In conclusion, the electronic scattering observed in high- T_c superconductors below T_c disagrees with the theoretical expectations based on $d_{x^2-y^2}$ pairing.

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