

## The role of off-plane oxygens in the enhancement of the pairing interaction in Cu-O clusters

This article has been downloaded from IOPscience. Please scroll down to see the full text article.

1990 J. Phys.: Condens. Matter 2 6337

(<http://iopscience.iop.org/0953-8984/2/29/514>)

View [the table of contents for this issue](#), or go to the [journal homepage](#) for more

Download details:

IP Address: 171.66.16.103

The article was downloaded on 11/05/2010 at 06:02

Please note that [terms and conditions apply](#).

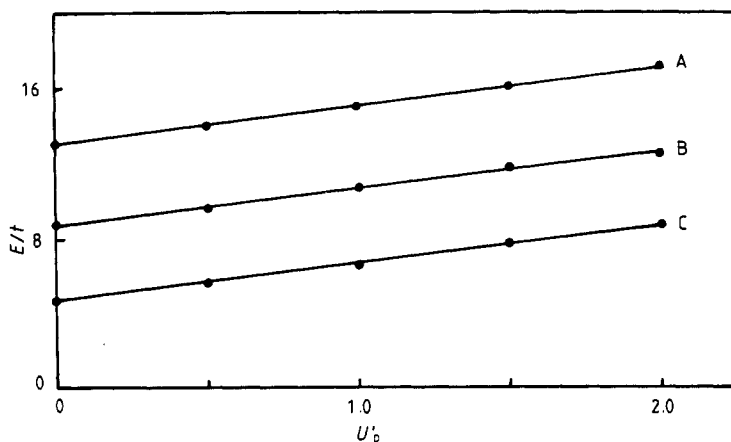
## ERRATUM

**The role of off-plane oxygens in the enhancement of the pairing interaction in Cu-O clusters** by D G Kanhere, V Sundararajan and P V Panat (*J. Phys.: Condens. Matter* 1989 1 9051–9055)

After the publication of this paper we detected an error in our program, as a result of which the main conclusion about the influence of the off-plane oxygens on the two-hole binding does not stand. In fact, no two-hole binding is observed for the range of parameters reported in the paper. The corrected versions of table 1 and figure 2 are given below.

**Table 1.** Values of  $\Delta$  for different values of  $U'_p$  for an 11-site cluster in units of  $t$ .

$V$	$U'_p$	$\Delta E(1\text{ h})$	$\Delta E(2\text{ h})$	$\Delta$
0	0.0	-4.3868	-8.3174	0.4562
	0.5	-4.3906	-8.3252	0.4560
	1.0	-4.3955	-8.3354	0.4556
	1.5	-4.4021	-8.3496	0.4546
	2.0	-4.4114	-8.3702	0.4526
2	0.0	-16.5735	-32.9086	0.2384
	0.5	-16.5737	-32.9090	0.2385
	1.0	-16.5739	-32.9093	0.2385
	1.5	-16.5741	-32.9097	0.2385
	2.0	-16.5743	-32.9101	0.2385



**Figure 2.** Total energy as a function of  $U$  for an 11-site cluster. Other parameters are:  $V = 0$ ,  $t' = 0.2$ ,  $E_d = 0$ . A, reference state; B, one-hole state; C, two-hole state.