JOURNAL OF SOLID STATE CHEMISTRY **51,** 133 (1984)

Erratum

Volume **49,** Number 2 (1983), in the article "A New Delafossite-Type Compound $CuYO_2$," by T. Ishiguro, N. Ishizawa, N. Mizutani, and M. Kato, pages 232–236, the atomic coordinates in Table IV (page 235) are incorrect. The correct coordinates are given below.

Atom	x	y	z
Cu	$\frac{1}{3}$	$\frac{2}{3}$	$\frac{1}{4}$
Y	0	0	0
O	$\frac{1}{3}$	$\frac{2}{3}$	0.0893(11)