Corrections and Additions

Acta Cryst. (1984). C**40**, 1762

The crystal structure of terephthalic acid: errata.

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Errors in the paper by Bailey & Brown [*Acta Cryst.* (1967), **22**, 387–391] are corrected. For Form II the correct unit-cell parameters are $a = 9{\cdot}54$ (1), $b = 5{\cdot}34$ (1), $c = 5{\cdot}02$ (1) Å, $\alpha = 86{\cdot}95$ (5), $\beta = 134{\cdot}65$ (5), $\gamma = 104{\cdot}90$ (5)°. Also for Form II the correct atomic coordinates are

	x	*y*	*z*
C(1)	0·147	−0·002	0·000
C(2)	0·175	0·217	0·184
C(3)	−0·031	−0·224	−0·186
C(4)	0·301	−0·007	0·003
O(1)	0·461	0·196	0·181
O(2)	0·275	−0·209	−0·158.

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Structure of the copper(II) dichloride complex of 1,4,10,13-tetraoxa-7,16-diazacyclooctadecane-7,16-diacetatocopper(II) methanol solvate, $[Cu(C_{16}H_{28}N_2O_8)].CuCl_2.CH_3OH$: erratum.

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A printer's error is corrected. The title of the paper by Gluziński, Krajewski, Urbańczyk-Lipkowska, Andreetti & Bocelli [*Acta Cryst.* (1984), C**40**, 778–781] is incorrect. The correct wording is given in the title of this erratum.