

Corrections and Additions

Acta Cryst. (1984), **C40**, 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 $a = 9.54$ (1), $b = 5.34$ (1), $c = 5.02$ (1) Å, $\alpha = 86.95$ (5), $\beta = 134.65$ (5), $\gamma = 104.90$ (5)°. Also for Form II the correct atomic coordinates are

	<i>x</i>	<i>y</i>	<i>z</i>
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-diacetato-copper(II) methanol solvate, [Cu(C₁₆H₂₈N₂O₈)]·CuCl₂·CH₃OH: erratum. By P. GLUZIŃSKI, J. W. KRAJEWSKI and Z. URBAŃCZYK-LIPKOWSKA, *Institute of Organic Chemistry, Polish Academy of Sciences, 01-224 Warszawa, Poland*, and G. D. ANDRETTI and G. BOCELLI, *Centro di Studio per la Strutturistica Diffraattometrica del CNR, Istituto di Strutturistica Chimica dell'Università di Parma, 43100 Parma, Italy*

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A printer's error is corrected. The title of the paper by Gluziński, Krajewski, Urbańczyk-Lipkowska, Andreotti & Bocelli [*Acta Cryst.* (1984), **C40**, 778–781] is incorrect. The correct wording is given in the title of this erratum.

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