

293 K, $F(000) = 840$, $wR = 0.042$ for 1781 observations. Methyl 1-methyl-2-nitrophenyl-4,5-diphenylpyrrole-3-carboxylate (III), $C_{25}H_{20}N_2O_4$, $M_r = 412.5$, monoclinic, $P2_1/n$, $a = 14.675$ (7), $b = 8.278$ (6), $c = 17.397$ (3) Å, $V = 2095$ (2) Å³, $Z = 4$, $D_x = 1.31$ g cm⁻³, $\mu = 0.84$ cm⁻¹, $T =$

293 K, $F(000) = 864$, $wR = 0.039$ for 2038 observations. These studies confirm the orientation of the reaction predicted by theoretical calculations.

Le résumé contient tous les détails pertinents.

Acta Cryst. (1992). C48, 220

Crystal structure and photochemistry of a methyl phenyl diester derivative of dibenzobarrelene.

Erratum. By JAMES TROTTER and FRED C. WIREKO, *Department of Chemistry, University of British Columbia, Vancouver, BC, Canada V6T 1Y6*

(Received 29 July 1991)

Abstract

An error in technical editing is corrected. In the paper by Trotter & Wireko [*Acta Cryst.* (1991), C47, 793–797] the title compound is incorrectly named as a 2,3-diester on

pages 793, 794 and 796; its correct name is methyl phenyl dibenzobarrelene-11,12-dicarboxylate.

All relevant information is given in the *Abstract*.

Acta Cryst. (1992). C48, 220

Structure of the modified nucleoside 2',3'-dideoxy-3'-fluorocytidine. Erratum. By H. L. DE WINTER, N. M.

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(Received 10 October 1991)

Abstract

In the paper by De Winter, Blaton, Peeters, De Ranter, Van Aerschot & Herdewijn [*Acta Cryst.* (1991), C47, 832–835] the fractional y coordinates of atoms C1'A, C3'A and

F3'A should be 0.5916 (3), 0.3130 (3) and 0.3924, respectively.

All relevant information is given in the *Abstract*.

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