SHORT COMMUNICATIONS

Contributions intended for publication under this heading should be expressly so marked; they should not exceed about 1000 words; they should be forwarded in the usual way to the appropriate Co-editor; they will be published as speedily as possible.

Acta Cryst. (1989). C45, 838

Structure of α-trans-cinnamic acid. Addendum. By Derk A. Wierda, Timothy L. Feng and Andrew R. Barron, Department of Chemistry, Harvard University, Cambridge MA 02138, USA

(Received 20 March 1989)

Abstract

The Related literature section of the paper by Wierda, Feng & Barron [Acta Cryst. (1989), C45, 338-339] should include the following additional reference: R. F. Bryan & D. P.

Freyburg (1975). J. Chem. Soc. Perkin Trans. 2, pp. 1835-1840.

All relevant information is contained in the Abstract.

Acta Cryst. (1989). C45, 838

Structure of 3β-dimethylamino-21-norcon-5-enine-20-one dihydrate. Erratum. By R. RADHAKRISHNAN and M. A. VISWAMITRA, Department of Physics and ICMR Centre on Genetics and Cell Biology, Indian Institute of Science, Bangalore-12, India and K. K. BHUTANI and R. M. VAID, Regional Research Laboratory, Jammu-1, India

(Received 28 March 1989)

Abstract

In the original publication [Acta Cryst. (1989), C45, 463-465] the final author was erroneously given as M. Ali. The synthesis of the title compound by an unusual oxidative

demethylation of the alkaloid conessine has recently been described [K. K. Bhutani & R. M. Vaid (1988). *Tetrahedron Lett.* 29, 359-360].

All relevant information is contained in the Abstract.

Acta Cryst. (1989). C45, 838

Structure of 1-(4-chlorophenacyl)-2-methyl-5-morpholino-4-nitroimidazole. Erratum. By Teresa Borowiak, Irena Wolska and Mariola Baryla, Faculty of Chemistry, Adam Mickiewicz University, Grunwaldzka 6, 60–780 Poznań, Poland and Stanislaw Sobiak, Department of Chemical Technology of Drugs, Karol Marcinkowski Academy of Medicine, Grunwaldzka 6, 60–780 Poznań, Poland

(Received 5 April 1989)

Abstract

An error in technical editing is corrected. The sentence beginning on line 5 of page 451 of the paper by Borowiak, Wolska, Baryla & Sobiak [Acta Cryst. (1989), C45, 448–451] should read: According to these bond distances, the inactivity of N(16) towards protonation can result from the structure

$$\begin{cases} N(16)^{+} = C(5) - C(4) = N(15)^{+} \\ O(151)^{-} \\ O(152)^{-} \end{cases}$$

that contributes to a resonance hybrid.

All relevant information is contained in the *Abstract*.

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