

13.751 (3) Å and thus the periodicity of the spiral of strong hydrogen bonding is much greater. The observed differences in the solid-state CD spectra of the morphine methyl iodide and sulfate salts from that of morphine free base must arise from this difference in periodicity or from the superposition of a second chiral hydrogen-bonding network on the primary chiral hydrogen-bonding network in the solid-state structure of morphine free base.

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Corrections and Additions

Acta Cryst. (1984). **C40**, 1490

Structure of methyl 8-isopropyl-3,3a,8,8a-tetrahydroindeno[2,1-c]pyrazole-8a-carboxylate, $C_{15}H_{18}N_2O_2$: erratum. By L. TOUPET and J. C. MESSAGER, *Groupe de Physique Cristalline, ERA au CNRS n° 070015, Université de Rennes, Campus de Beaulieu, 35042 Rennes CEDEX, France*

(Received 22 May 1984)

In the abstract of the paper by Toupet & Messenger [*Acta Cryst.* (1984). **C40**, 330–331], the cell parameters are incorrect. The correct values are: $a = 5.791$ (4), $b = 15.503$ (4), $c = 15.954$ (5) Å, $\alpha = 82.24$ (5), $\beta = 79.35$ (6), $\gamma = 79.13$ (5)°, $V = 1375$ (4) Å³.

0108-2701/84/081490-01\$01.50

Acta Cryst. (1984). **C40**, 1490

Structure of 6-chloro-4-phenyl-1,2,3-benzoxathiazine 2,2-dioxide, $C_{13}H_8ClNO_3S$: erratum. By SAFIA MEHDI and B. RAMA RAO, *X-ray Section, Regional Research Laboratory, Hyderabad-500 007, AP India*

(Received 15 June 1984)

A printer's error is corrected. In the paper by Safia Mehdi & Rama Rao [*Acta Cryst.* (1984). **C40**, 1057–1059] two of the atom labels in Table I are incorrect. The correct labels with their corresponding fractional coordinates are given below:

	x	y	z
C(5)	9604 (9)	4384 (5)	3791 (4)
C(6)	9878 (9)	5304 (5)	3782 (4)
Cl	10979 (3)	5813 (2)	4392 (1)

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