

Acta Cryst. (1980). B36, 2858

The structure of $K_2(K_{0.41}, H_2O_{0.59})_6 Na_{3.98} H_3 O_{0.78} X_{0.68}^{+} (Fe_{0.05}^{2+}, \square_{0.95}) Fe_6^{3+} O_2 (SO_4)_{12} \cdot 11 \cdot 91 H_2 O$: erratum.
By FERNANDO SCORDARI, *Istituto di Mineralogia e Petrografia, Piazza Umberto I°, 70121 Bari, Italy*

(Received 22 September 1980)

Abstract

A printer's error is corrected. In the paper by Scordari [*Acta Cryst.* (1980), B36, 1733–1738] the sentence beginning on the nineteenth line of the left-hand column of p. 1735 should read: After several attempts we preferred to fix the

0567-7408/80/112858-01\$01.00

coordinates of the atoms refined in $P\bar{3}$ and then to proceed to refine in $P3$ the occupancy coefficients of all the other atoms deduced from the Fourier map and steric considerations.

All the information is given in the *Abstract*.

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Acta Cryst. (1980). B36, 2858–2859

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