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addenda and errata

Order–disorder transition in monoclinic sulfur: a precise structural study by high-resolution neutron powder diffraction. Corrigendum

W. I. F. David,[a] R. M. Ibberson,[a]* S. F. J. Cox[a] and P. T. Wood[b]‡

[a]ISIS Facility, CCLRC, Rutherford Appleton Laboratory, Chilton, Didcot, Oxfordshire OX11 0QX, England, and [b]School of Chemical Sciences, University of East Anglia, Norwich, Norfolk NR4 7TJ, England

‡ Present address: The University Chemical Laboratory, University of Cambridge, Lensfield Road, Cambridge CB2 1EW, England.
Correspondence e-mail: r.m.ibberson@rl.ac.uk

Revised lattice parameters for Table 1 of the paper by David *et al.* (2006), *Acta Cryst.* B**62**, 953–959, are given.

The lattice constants for β-sulfur at 100 K given in Table 1 of the paper by David *et al.* (2006) are incorrect. Correct values are: a = 10.8027 (1), b = 10.6911 (1), c = 10.6689 (1) Å; β = 95.7124 (10)°; V = 1226.06 (2) Å^3. Also, the a and c axis labels shown in Fig. 2(a) should be reversed.

References

David, W. I. F., Ibberson, R. M., Cox, S. F. J. & Wood, P. T. (2006). *Acta Cryst.* B**62**, 953–959.

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