

Data collection and cell refinement: *Diffractometer Control Program for the Hilger & Watts (Y290)* (Kopf & Abeln, 1993). Data reduction: *WATSHL* (Kopf, 1987). Program(s) used to solve structure: *SHELXS86* (Sheldrick, 1990a). Program(s) used to refine structure: *SHELXL93* (Sheldrick, 1993). Molecular graphics: *SHELXTL-Plus* (Sheldrick, 1990b). Software used to prepare material for publication: *SHELXL93 CIFTAB*; *FCF2FOC* (Kopf, 1992).

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Lists of structure factors, anisotropic displacement parameters, H-atom coordinates and complete geometry have been deposited with the IUCr (Reference: AB1150). Copies may be obtained through The Managing Editor, International Union of Crystallography, 5 Abbey Square, Chester CH1 2HU, England.

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ADDENDA AND ERRATA

Acta Cryst. (1994). **C50**, 1372

The structures of alkalides and electrides. II. Structure of caesium bis(15-crown-5) electride.

Erratum. By DONALD L. WARD, RUI H. HUANG, MARK E. KUCHENMEISTER and JAMES L. DYE, *Department of Chemistry and Center for Fundamental Materials Research, Michigan State University, East Lansing, Michigan 48824, USA*

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Abstract

An error in printing is corrected. In Table 1 of the paper by Ward, Huang, Kuchenmeister & Dye [*Acta Cryst.* (1990), **C46**, 1831–1833] the *y* coordinate of atom O1 is printed

incorrectly as –0.1501 (2). The correct value of this coordinate is +0.1501 (2).

All relevant information is given in the *Abstract*.