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**The experimental charge density in sulfur-containing molecules: a study of the deformation electron density in sulfamic acid at 78 K by X-ray and neutron diffraction. Corrigendum.** BY J. W. BATS and P. COPPENS, *Department of Chemistry, State University of New York at Buffalo, Buffalo, New York 14214, USA* and T. F. KOETZLE, *Department of Chemistry, Brookhaven National Laboratory, Upton, New York 11973, USA*

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An improved reproduction of Fig. 1 of the paper by Bats, Coppens & Koetzle [*Acta Cryst.* (1977), B33, 37–45], giving deformation density in sections through the sulfamic acid molecule, is shown.

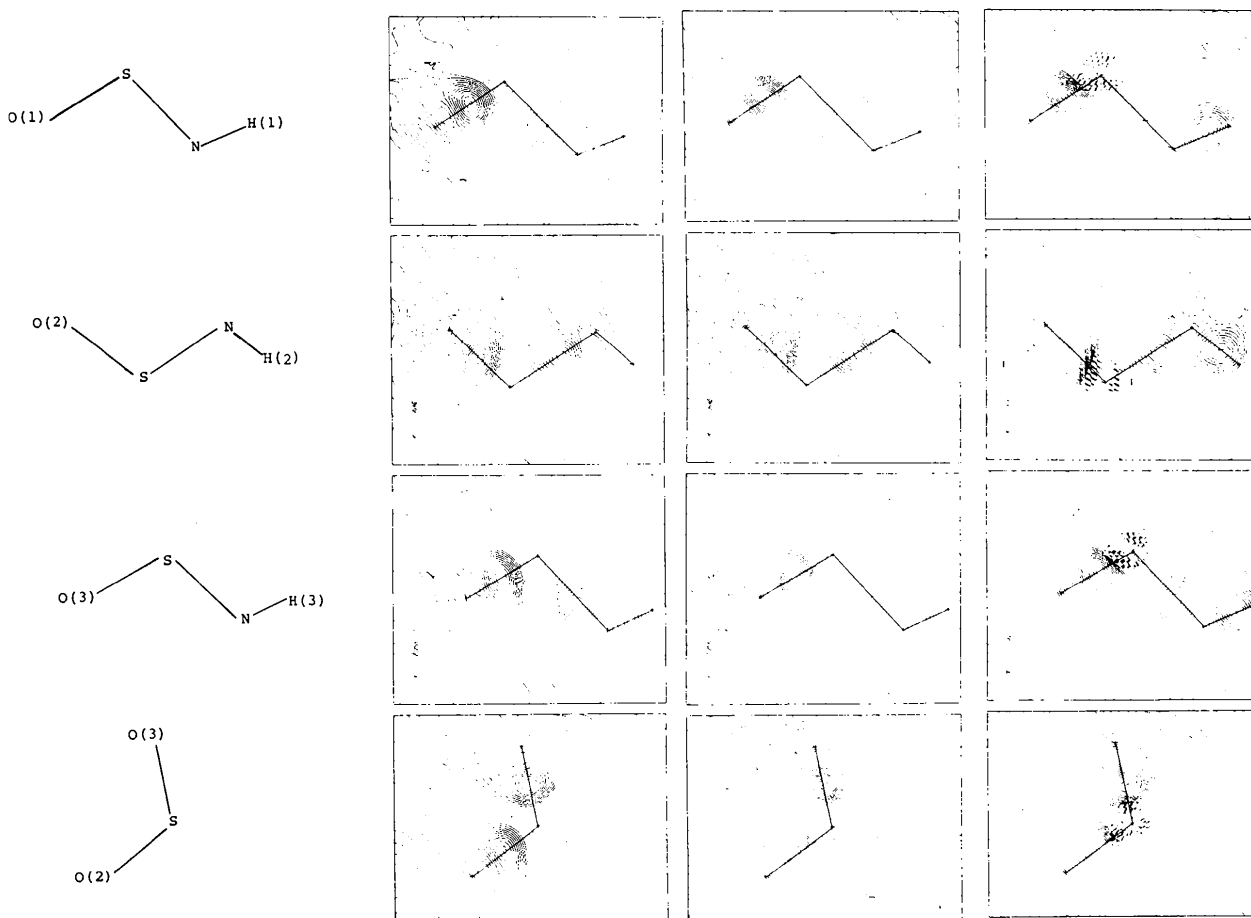


Fig. 1. Deformation density in sections through the molecule. Column 1: X–X (conventional), column 2: X–X (high order,  $\sin \theta/\lambda > 1.0 \text{ \AA}^{-1}$ ), column 3: rescaled X–N maps. Data cut-off in Fourier synthesis:  $(\sin \theta/\lambda)_{\max} = 0.85 \text{ \AA}^{-1}$ . The contour interval is  $0.05 \text{ e \AA}^{-3}$ . Negative contours are shown as broken lines.