Debye–Waller factor was reviewed. The Commission agreed that the significance of such estimated standard deviations may be ambiguous. In the case of high anisotropy, the estimated standard deviation is necessarily large although the individual anisotropic parameters may be well determined. In future, the presence of unusual anisotropy should be referred to in the text (including the maximum and minimum amplitudes and any nonpositive-definite coefficients found), or in the table of $B_{eq}$ (by use of an asterisk), or illustrated by a plot of the atomic vibrational ellipsoids.

**Calculated hydrogen-atom coordinates:** Calculated hydrogen-atom coordinates will be published in future only at the Co-editor’s discretion and if they are necessary to the understanding of the paper. They will otherwise be deposited.

**Graphical chemical formulae:** A graphical structural formula should always be given in the report of a structure determination of an organic or organometallic compound. The figure showing the atomic positions is not adequate for this purpose. However, a plot of the atomic vibrational ellipsoids can be used to illustrate atomic positions. Authors should ensure that such plots are of good contrast and quality. The numbering of atoms should be consistent throughout a paper and, as far as possible, correspond to the systematic name [Acta Cryst. (1982). B38, 700].

**Least-squares planes:** Least-squares planes and the deviations from them will only be published if they are referred to in the text of the paper and are significant in the consideration of the structure. They will otherwise be deposited.

**Absorption correction:** In a structural paper the absorption correction, if any, should always be described and the maximum and minimum corrections stated.

**Melting point:** In a structural paper the melting point of a compound should always be given if it is known.

**Stereofigures:** The requirement on stereofigures is one per structure unless the Co-editor and referees feel that more are necessary for the understanding of the structure described in the paper. These stereo views must fit into a single column (80 mm wide). Authors are reminded that a nonstereo view (half a pair) is often an acceptable alternative to the stereo pair. In stereo pairs the relative sizes of the molecule and the whole figure should be such that when the figure is printed in a column of 80 mm, the individual atoms are easily distinguishable. The center-to-center separation in stereo figures must not exceed 55 mm. In a charge density paper only one or two figures are required to illustrate the techniques or results described: any others will be deposited. The text should be adequate to give the remaining information.

**Figures**

**Size:** Illustrations should normally present information so that each figure or part of a figure can be printed in one column (80 mm width). Co-editors will need to be satisfied that the information density is high enough, if authors wish figures to be printed larger than this.

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