

the triennial International Congresses of Crystallography. The first prize will be presented during the XIV Congress at Perth, Australia, in 1987. This year will be the seventy-fifth anniversary of the discovery of X-ray diffraction in 1912.

Any scientist who has made contributions of exceptional distinction to the science of crystallography is eligible for the Ewald Prize, irrespective of nationality, age or experience. No restrictions are placed on the time or the means of publication of his or her contributions. The prize may be shared by several contributors to the same scientific achievement.

Nominations for the Ewald Prize are invited. They should be submitted in writing, accompanied by supporting documentation, to the Executive Secretary of the International Union of Crystallography, 5 Abbey Square, Chester CH1 2HU, United Kingdom. The closing date for nominations is 30 September 1986.

TH. HAHN  
President

K. V. J. KURKI-SUONIO  
General Secretary

*Acta Cryst.* (1986). B42, 414

#### Commission on Journals Atomic Labelling

The Commission on Journals recently voted to simplify the form of atomic labelling. Atoms of the same chemical species within an asymmetric unit should be distinguished in future by an appended arabic numeral, with parentheses around the numeral required only in the event of ambiguity, such as C11 *versus* C1(1) or C(11).

*Acta Cryst.* (1986). B42, 414

#### Appointments in the IUCr Office

A notice concerning recent appointments in the IUCr office appeared in the July 1986 issue of *Acta Crystallographica*, section A, p. 288.

### Book Reviews

*Works intended for notice in this column should be sent direct to the Book-Review Editor (J. H. Robertson, School of Chemistry, University of Leeds, Leeds LS2 9JT, England). As far as practicable books will be reviewed in a country different from that of publication.*

*Acta Cryst.* (1986). B42, 414-415

**International tables for crystallography: Brief teaching edition of Volume A: Space group symmetry.** Edited by T. HAHN. Pp. viii + 119. Dordrecht: Reidel, 1985. Price Dfl27.50, \$8.50, £7.60.

This is a slim volume consisting of 24 selected space-group descriptions and several pages of the basic text sections from Volume A of *International tables for crystallography*. The declared purpose of the *Teaching edition* is:

(i) to provide a handy and inexpensive tool for researchers and students to become familiar with the use of the space-group tables in Volume A;

(ii) to be used in classroom teaching;

(iii) to serve as a laboratory handbook, because the 24 examples include most of the frequently occurring space groups, for both organic and inorganic crystals.

The basic text section consists of material found in chapters 1, 2, 3 and 5 of Volume A: namely *Symbols and terms used*, *Guide to the use of the space-group tables*, *Space-group determination and diffraction symbols*, and *Transformations in crystallography*.

I approve strongly of the idea of bringing out an introductory book on the excellent, but detailed, Volume A, as there is a serious need for material which can be read and understood by beginners and non-specialists. I was therefore delighted to hear that the International Union of Crystallography had, at last, brought out such a text. However, I am sorry to say that, when I opened my review copy, I felt

somewhat surprised and, even, disappointed. The title had suggested a digestible and specially written account, which could be recommended to someone meeting, for the first time, the beauty and mysteries of crystallographic symmetry. Instead of the elementary treatment I had anticipated, it seems that the basic text has been lifted entirely without change from Volume A.

As a result, I certainly could not expect a novice to take the text away and learn from it by him/herself. Indeed, because no attempt has been made to adapt the text to suit its declared aims, on almost every page peculiar inconsistencies can be found which ensure that it cannot stand alone as a teaching text. It is particularly irritating to encounter so many references to sections that are missing entirely. As typical examples, in the *Guide to the use of the space-group tables*, page 10, the reader is referred to sections 8.3.1 and 9.1 for further details on lattice centring and to section 9.3 for reduced bases. On page 30 the index of a maximal subgroup requires the footnote in section 8.1.5 to be consulted. Again, picking a page at random, I find on page 39 reference to a missing Table 4.3.1, and so on. This occurs so frequently throughout that one might as well use Volume A in the first place.

The main problem with the *Teaching edition* is that it falls between two stools. On the one hand, it is too difficult to act simply as a text for learning about the space-group tables and, on the other, it is not complete enough for the experienced researcher to use as a handbook. In any case, those who already know about space groups and the *International tables* are unlikely to want to consult the introduc-

tory texts in this book. I believe that it would have been better either to have rewritten the introductory text to suit the level of reader expected or to have published just the space-group pages, together with the useful key on the inside covers (incidentally, the book could then have been even less expensive!). I wish I could be more positive about the *Teaching edition* but, frankly, I am not sure to whom it can be recommended.

A. M. GLAZER

Clarendon Laboratory  
University of Oxford  
Parks Road  
Oxford OX1 3PU  
England

*Acta Cryst.* (1986). B42, 415

**X-ray crystallography.** By E. M. UYGUR. Pp. x+318. Published by METU, Ankara, Turkey, 1983 (available from the Metallurgical Engineering Department). Price US \$5.00.

This book, which is based on 'class notes developed over many years of teaching of X-ray diffraction and crystallography courses' by the author, would be better titled *Elements of X-ray diffraction*. The rather general title of *X-ray crystallography* would demand a more comprehensive coverage, which the book lacks. I would particularly like to stress the fact that there is not even a cursory treatment of the determination of crystal structure and the phase problem, which is a vital area in X-ray crystallography.

The geometry of crystals, symmetry mapping and projections are routinely discussed in the first three chapters. This is followed by a treatment of generation and properties of X-rays in Chapter 4. Chapters 5 and 6 deal with X-ray diffraction proper and the experimental techniques.

Except for the drawback mentioned earlier the sections covered have been treated with clarity with good examples and illustrations. The book should find use as a supplement to standard texts in X-ray diffraction and crystallography.

R. SRINIVASAN

Department of Crystallography & Biophysics  
University of Madras  
Guindy Campus  
Madras-600025  
India

*Acta Cryst.* (1986). B42, 415-416

**Structure and statistics in crystallography.** Edited by A. J. C. WILSON. Pp. vii+225. New York: Adenine Press, 1985. Price US \$65.00.

This book describes the proceedings of the symposium on Crystallographic Statistics held in Hamburg, West Germany in August 1984 in the course of the Thirteenth International Congress of the International Union of Crystallography. It also includes a few papers presented in the main Congress

but which were considered as closely linked with the symposium topic.

The first and last contributions in the book stand out on their own, the first by J. Karle on the statistical basis underpinning direct methods and the last on 'expert systems' of data acquisition. The direct-methods contribution does not contain any new material but it is a splendid review of the theoretical developments which led to direct methods having their present pre-eminence and in which Karle and Hauptman played such a leading role. The 'expert systems' paper by H. J. Milledge and her collaborators presents the principles by which the data-acquisition process by a diffractometer may be optimized by analysing the data during collection.

The remainder, the bulk, of the book divides into two roughly equal parts - the first concerned with intensity statistics and the second with refinement processes. A most interesting paper by Weiss *et al.* deals with the representation of probability density functions by Fourier series, which is much better than previous methods using the central-limit theorem or based on the Edgeworth or Gram-Charlier series. There follow three papers dealing respectively with the effects of heavy atoms, non-crystallographic centres of symmetry and non-crystallographic translational symmetry on the normal or cumulative intensity or  $|E|$  distributions. Since many crystal structures contain heavy atoms or a great deal of symmetry to do with the chemistry of molecules rather than the requirements of space groups, it is clear that departures from idealized random distributions of almost-equal atoms must be common. The papers presented here show much success in predicting distributions from known structural features; it is not quite so clear that the inverse problem has been solved.

After a paper by Parthasarathy & Elango on the best way of testing for symmetry elements from intensity statistics the section is rounded off by a contribution from Wilson on fluctuations and errors in intensity distributions. He concludes, regretfully, that there is no obvious easy way of representing distributions modified by random or systematic errors.

The second section, on refinement, starts with a paper by Prince commenting on the precision and accuracy which may be obtained in structure refinement by the Rietveld method. He concludes that while the calculated standard deviations may give a general indication of the precision of the parameters found they are not an accurate assessment of the r.m.s. errors. Clearly this paper was controversial; the following paper by Rollett is a discussion of Prince's paper and he comes to a contrary conclusion.

The next two papers are concerned with the application of information theory to refinement. The first, by Collins, is on the very topical subject of parameter estimation by entropy maximization. This is a good paper to read; firstly it confirms that there is a certain arbitrariness in the entropy function which is maximized and secondly it demystifies a topic which for many crystallographers has taken on the characteristics of a deity - all powerful and incomprehensible. The following paper by Wilkins *et al.* is similarly to be commended especially in providing a practical procedure for incorporating prior knowledge into information-theory procedures.

The three papers which follow, on the modification of weights in least-squares analysis, variance of intensities in the Bond method and the use of maximum likelihood and