and hexagonal lattices. Furthermore, the point group $\overline{3}m$ in the hexagonal description is further subdivided into two cases - one having the normal of the mirror planes in the [100] direction and the other one having the normal in the [210] direction. The Laue group $2/m$ is also subdivided into two cases - with the $b$ axis unique and with the $c$ axis unique.

Since the distribution of the X-ray intensity $I(hkl)$ in reciprocal space shows the same symmetry as a general crystal form of the corresponding Laue class, the asymmetric unit of X-ray intensity data for a crystal is the same as that for the crystal forms $\{hkl\}$ with the symmetry of its Laue group.

Table 1 shows information on the 11 Laue groups (column 3). The crystal system is given in column 1, all corresponding noncentrosymmetric point groups are listed in column 2. Column 4 describes one asymmetric unit for the intensity data of each Laue group. Alternative possibilities are separated by dashed lines. The last column shows the fraction of reciprocal space belonging to one asymmetric unit.

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References

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*International Tables for Crystallography*
Volume C: Mathematical, Physical and Chemical Tables
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Volume C is the result of many years of collaborative effort to revise and update Volumes II, III and IV of *International Tables for X-ray Crystallography*. Most of the contributions are entirely new; the remaining sections are the result of editorial condensation of material from the earlier volumes or have been revised by their original authors. Copies may be ordered directly from the publishers (Kluwer Academic Publishers, PO Box 17, 3300 AA Dordrecht, The Netherlands), from Polycrystal Book Service, Box 3439, Dayton, Ohio 45401, USA, or from any bookseller. [Price Dfl 400 (or Dfl 200 for individuals, who should give a written undertaking that the copy is for their personal use only and will not be made available to libraries etc.).]

A full review will be published in *Acta Crystallographica*, Section A.

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Book Reviews

Works intended for notice in this column should be sent direct to the Book-Review Editor (R. F. Bryan, Department of Chemistry, University of Virginia, McCormick Road, Charlottesville, Virginia 22901, USA). As far as practicable, books will be reviewed in a country different from that of publication.


Fundamentals of crystallography. By C. GIACOVAZZO,  
H. L. MONACO, D. VITERBO, F. SCORDARI, G. GILLI,  

For this book, the second of the new IUCr Texts on Crystallography, Carmelo Giacovazzo has edited, revised, expanded and brought up to date in English translation Introduzione della Cristallografia Moderna, first published in 1985. The stated aim of the book is to provide a compact comprehensive account of modern crystallographic subjects that should be useful as a text for university courses that cover crystallography, fully or only partially, but that should also be of use at the doctoral or research level. I believe that this book more than adequately meets these goals and the high quality of the translation should ensure it a significantly wider audience than that reached by the original.

Recognizing the interdisciplinary character of modern crystallography, the editor realized that a book of this kind would need to be written by several authors and that the various contributions would have to be “carefully harmonized in order to conform them to a unified plan”. The result is a text that can be used with profit in its entirety but from which instructors and students can select according to their particular interests. There are three core chapters. Chapter 1 (Giacovazzo, 60 pp.) is a concise treatment of symmetry in crystals. The first half covers symmetry elements, lattices, point groups, Laue classes, crystal systems, Bravais lattices, and plane, line and space groups. This material is fundamental for audiences at all levels. For the more advanced student, the second half of the chapter consists of appendices dealing with isometric transformations, combinations of movements, Wigner-Seitz cells, space-group rotation matrices, symmetry groups and generalized symmetry with an introduction to $G$ groups and color symmetry. Very usefully, in this last case, actual structural arrangements are used as illustrations. The second core chapter (Chapter 3, Giacovazzo, 86 pp.) covers the diffraction of X-rays by crystals. The same division between fundamental and more advanced topics is adopted here. The treatment is kinematic, with a progression from scattering by electrons to that by atoms, molecules and crystals. There are good treatments of symmetry in reciprocal space and of anomalous dispersion, and a short introduction to modulated structures. The appendices here deal with the mathematics of Fourier transforms and convolution operations, and with more subtle aspects of scattering such as Compton scattering, the anisotropic temperature factor, the Renninger effect and electron and neutron scattering. Scattering by non-crystalline materials such as gases, liquids and amorphous solids is covered, as is small-angle scattering. Two final appendices deal with electron-density mapping and a more advanced look at modulated structures and quasicrystals.

The newcomer to crystallography, the third core chapter (Chapter 5, Viterbo, 79 pp.) deals with the solution and refinement of crystal structures. After a general introduction come sections dealing with the statistical analysis of structure amplitudes, the use of the Patterson method and direct methods, and refinement by least-squares and difference Fourier methods, with a short discussion of the determination of absolute configuration. For the more experienced reader, the appendices deal with structure-factor and triplet-invariant probability distributions, Patterson vector methods, pseudotranslational symmetry, magic integers and newer multisolution methods and procedures for completing a partial model. As is the case for the other core chapters, the references here are particularly helpful and cover the literature into 1990.