

# Notes for Authors

## International School on Mathematical and Theoretical Crystallography

### 1. Submission

Authors should submit their paper *via the web* at <http://journals.iucr.org/a/services/conferenceproceedings.html>. Papers are expected to be submitted by **1 September 2005**.

### 2. Manuscript preparation

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The manuscript should be prepared using Word,  $\text{\TeX}$  or  $\text{\LaTeX}$ . Authors are encouraged to use the templates available from the Editorial Office by e-mail ([med@iucr.org](mailto:med@iucr.org)), by ftp (from the 'templates' directory at <ftp://ftp.iucr.org>) or on the WWW at <http://journals.iucr.org/a/services/authorservices.html> where full instructions are given.

Figures should be prepared in PostScript, encapsulated PostScript or TIFF formats. The resolution of bitmap graphics should be a minimum of 600 d.p.i.

For submission, the text and figures should be provided as a single Word/RTF, PostScript or PDF file. Once the paper is accepted, separate source files for the text and figures should be uploaded at the URL provided by the Co-editor.

#### 2.2. General

Contributions will be published as full-length papers and should not exceed six journal pages (~6000 words). The paper should conform to the general editorial style of *Acta Crystallographica Section A*.

#### 2.3. Title and authors

The title should be short and informative. The contact author should provide an e-mail address and a fax number. The e-mail address will be used for editorial communications and despatch of electronic proofs and reprints.

#### 2.4. Abstract and synopsis

All contributions must be preceded by an *Abstract* in English. The *Abstract* should be no more than 200 words and it should make no reference to tables, diagrams or formulae contained in the body of the paper. Literature references in the *Abstract* are discouraged but if a reference is unavoidable it

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Authors should supply a synopsis of no more than two sentences for inclusion in the contents listing of the journal.

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The International System of Units (SI) is used except that the ångström (symbol Å, defined as  $10^{-10}$  m) is generally preferred to the nanometre (nm) or picometre (pm) as the appropriate unit of length. Recommended prefixes of decimal multiples should be used rather than ' $\times 10^n$ '.

#### 2.6. Figures

Figures should be prepared using one of the file formats listed in §2.1.

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Tables which are produced in Word should be prepared using the Word table editor.

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Displayed mathematical equations which are produced in Word should be prepared using the Word equation editor.

The Editor may require that Appendices giving details of mathematical derivations given only in outline in the main text be deposited in the IUCr electronic archive.

#### 2.9. References

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Andrews, M. D., Wright, H. E. & Clarke, S. A. (2005). In preparation.  
Brünger, A. T. (1992). *X-PLOR. Version 3.1. A System for X-ray Crystallography and NMR*. Yale University, Connecticut, USA.

Collaborative Computational Project, Number 4 (1994). *Acta Cryst. D* **50**, 760–763.

Crowther, R. A. (1972). *The Molecular Replacement Method*, edited by M. G. Rossmann, pp. 173–178. New York: Gordon and Breach.

International Union of Crystallography (2001). (*IUCr*) *Structure Reports Online*, <http://journals.iucr.org/e/journalhomepage.html>.

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