

this in turn will tend to make the Hessian more strongly convex, thus aiding the structure-determination process [see (14)]. By this means, the information contained in a molecular envelope (solvent flattening) and molecular fragments could, for example, be incorporated in the map.

### References

- BRICOGNE, G. (1993). *Acta Cryst. D* **49**, 37–60.  
 GULL, S. F. (1989). *Maximum Entropy and Bayesian Methods*, edited by J. SKILLING, pp. 53–71. Dordrecht: Kluwer Academic Publishers.
- GULL, S. F. & SKILLING, J. (1984). *Indirect Imaging*, edited by J. A. ROBERTS, pp. 267–279. Cambridge Univ. Press.  
 JAYNES, E. T. (1986). *Maximum Entropy and Bayesian Methods in Applied Statistics*, edited by J. JUSTICE, pp. 1–58. Cambridge Univ. Press.  
 MÜLLER, J. J. & HANSEN, S. (1994). *J. Appl. Cryst.* In the press.  
 SKILLING, J. (1989). *Maximum Entropy and Bayesian Methods*, edited by J. SKILLING, pp. 42–52. Dordrecht: Kluwer Academic Publishers.  
 TITTERINGTON, D. M. (1985). *Astron. Astrophys.* **144**, 381–387.  
 WILKINS, S. W., STEENSTRUP, S. & VARGHESE, J. N. (1985). *Structure and Statistics in Crystallography*, edited by A. J. C. WILSON, pp. 113–123. Adenine Press.

*Acta Cryst.* (1994). **A50**, 550

**Determination of quasicrystalline structures: a refinement program using symmetry-adapted parameters. Erratum.** By L. ELCORO, J. M. PEREZ-MATO and G. MADARIAGA, Departamento de Física de la Materia Condensada, Facultad de Ciencias, Universidad del País Vasco, Apartado 644, Bilbao, Spain

(Received 1 May 1994)

### Abstract

A typesetting error in equation (28) of Elcoro, Perez-Mato & Madariaga [*Acta Cryst.* (1994), **A50**, 182–193] is corrected. The correct equation is

$$\begin{aligned} F(\mathbf{H}) = & [|A|/V(\mathbf{a}_i)] \sum_{\mu, m} p_m(\mu) f_m(\mathbf{H}) \\ & \times \sum_R \exp(-\widetilde{\mathbf{R}}\mathbf{H}\beta_\mu^m \mathbf{R}\mathbf{H}) \\ & \times \exp[2\pi i \hat{\mathbf{h}} \cdot (\hat{\mathbf{R}}\hat{\theta}_\mu + \hat{\mathbf{t}})] \int d\varphi_1 \dots d\varphi_{n-4} \end{aligned}$$

$$\times \int_a^b dr J(r, \varphi_1, \dots, \varphi_{n-4}) \exp[2\pi i (\tilde{\mathbf{T}}_\mu^{-1} \tilde{\mathbf{R}}_i \mathbf{h}) \cdot \mathbf{x}_i], \quad (28)$$

where

$$a = \sum_i a_i^{\mu, \text{in}} Z_i(\varphi_1, \dots, \varphi_{n-4})$$

and

$$b = \sum_i a_i^{\mu, \text{ex}} Z_i(\varphi_1, \dots, \varphi_{n-4}).$$

All information is given in the *Abstract*.

*Acta Cryst.* (1994). **A50**, 550

**Dynamical X-ray diffraction from imperfect crystals: a solution based on the Fokker–Planck equation. Erratum.** By T. J. DAVIS, CSIRO Division of Materials Science and Technology, Private Bag 33, Rosebank MDC, Clayton, Victoria 3169, Australia

(Received 25 May 1994)

### Abstract

The following mathematical expressions were incorrectly printed in the paper by Davis [*Acta Cryst.* (1994), **A50**, 224–231].

Page 225: the correct expression for  $\chi_{-h}$  is

$$\chi_{-h} = -C(\gamma_h/\gamma_o)\chi'_{-h}.$$

Page 228: equations (22) and (23) should read

$$\begin{aligned} R(t) = R_o &+ [R(t') - R_o] \exp[2i\alpha\omega(t-t')] \\ &\times (1 + [R(t') - R_o](\chi_{-h}/2\omega)) \\ &\times \{1 - \exp[2i\alpha\omega(t-t')]\}^{-1}, \quad (22) \end{aligned}$$

$$\omega = \pm(\beta^2 - \chi_h \chi_{-h})^{1/2}. \quad (23)$$

All information is given in the *Abstract*.