

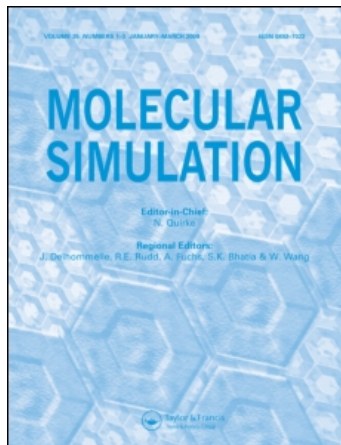
This article was downloaded by:

On: 14 January 2011

Access details: *Access Details: Free Access*

Publisher *Taylor & Francis*

Informa Ltd Registered in England and Wales Registered Number: 1072954 Registered office: Mortimer House, 37-41 Mortimer Street, London W1T 3JH, UK



Molecular Simulation

Publication details, including instructions for authors and subscription information:

<http://www.informaworld.com/smpp/title~content=t713644482>

Erreta

To cite this Article (1989) 'Erreta', *Molecular Simulation*, 4: 4, 251

To link to this Article: DOI: 10.1080/08927028908022368

URL: <http://dx.doi.org/10.1080/08927028908022368>

PLEASE SCROLL DOWN FOR ARTICLE

Full terms and conditions of use: <http://www.informaworld.com/terms-and-conditions-of-access.pdf>

This article may be used for research, teaching and private study purposes. Any substantial or systematic reproduction, re-distribution, re-selling, loan or sub-licensing, systematic supply or distribution in any form to anyone is expressly forbidden.

The publisher does not give any warranty express or implied or make any representation that the contents will be complete or accurate or up to date. The accuracy of any instructions, formulae and drug doses should be independently verified with primary sources. The publisher shall not be liable for any loss, actions, claims, proceedings, demand or costs or damages whatsoever or howsoever caused arising directly or indirectly in connection with or arising out of the use of this material.

ERRATA

Molecular Simulation, 3, pp. 27–47 (1989)

MOLECULAR DYNAMICS MODELLING OF POLYMER MATERIALS

J.H.R. CLARKE and D. BROWN

Chemistry Department U.M.I.S.T., Manchester M60 1QD, UK.

Note that the following corrections should be made to the original article:

- (i) On page 30 the coefficient C_3 in equation (2) has the value -31800 i.e. it is negative and not positive as published.
- (ii) On page 31 the 'h' in equation (3) should be bold to denote its tensor character.

Molecular Simulation, 4, pp. 121–136 (1989)

COMPUTER SIMULATIONS IN ZEOLITE CATALYSIS RESEARCH

C.J.J. DEN OUDEN¹, B. SMIT¹, A.F.H. WIELERS¹, R.A. JACKSON²,
and A.K. NOWAK³

¹ *Koninklijke/Shell Laboratorium Amsterdam (Shell Research B.V.), Badhuisweg 3,
1031 CM Amsterdam, The Netherlands*

² *Department of Chemistry, University of Keele, Keele, Staffordshire, ST5 5BG, UK*

³ *Davy Faraday Laboratory, The Royal Institution, 21 Albemarle Street,
London W1X 4BS, UK*

The caption for the figure on page 124 should read:

Figure 3 Structure of one faujasite (structure of zeolite X and Y) cage. The faujasite structure is built up by interlinking various cages. Faujasite is characterised by the presence of a pore system rather than a channel system. The entrance to the pores is 8 Å. Oxygen atoms are coloured yellow, T-sites (silicon or aluminium atoms) are tetrahedrally surrounded by oxygen atoms and are coloured red. (See colour plate VII.)