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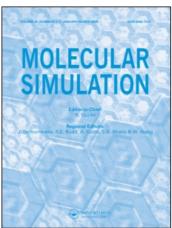
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## **Molecular Simulation**

Publication details, including instructions for authors and subscription information: http://www.informaworld.com/smpp/title~content=t713644482

## **Guest Editorial**

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To cite this Article Alba-Simionesco, Christiane , Fuchs, Alain and Vacher, René(1997) 'Guest Editorial', Molecular Simulation, 20: 1, 1

To link to this Article: DOI: 10.1080/08927029708024163 URL: http://dx.doi.org/10.1080/08927029708024163

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## **GUEST EDITORIAL**

This issue of Molecular Simulation contains invited papers presented at the meeting on "computer simulation of glasses and related materials" held under the auspices of the french cooperative research group on oxide glasses (Groupement de Recherche n°1025). This group is sponsored by CNRS (Centre National de la Recherche Scientifique). Its aim is to promote collaborative research on glassy materials. A part of its activities is the organisation of meetings on chosen themes. This meeting took place on 10th July 1996 at University Pierre et Marie Curie in Paris. The papers that follow provide an overview of the problems that were adressed in this meeting.

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