

## **Announcements**

### *International Meeting* **Advances in Biomolecular Simulations**

Strasbourg, France, 12–14 March 1991

The program includes the following topics:

Potential function development  
Quantum-mechanical studies of small molecules  
Models for solvents  
Dynamical simulation methods  
Methods for searching configurational space  
Methods for studying transitions between conformers  
Protein-engineering applications

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