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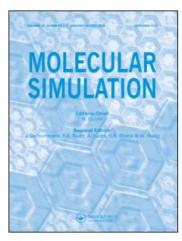
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Molecular Simulation exists to bring together the most significant papers concerned with applications of simulation methods using statistical mechanics, and original contributions to the development of simulation methodology. The aim is to provide a forum in which cross-fertilization between application areas and methodologies can take place and new developments can be encouraged.

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