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The number of application areas is continually increasing and *Molecular Simulation* will keep pace with events, welcoming papers on topics ranging from condensed matter physics and chemistry to biomolecules and rheological studies. Similarly, methodological papers are encouraged dealing with, for example, nonequilibrium methods or quantum simulations and new developments in languages and machines.

This journal will publish review articles and preliminary communications as well as full-length papers. Occasionally, reports of papers presented at meetings will be published.

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