



Theoretical and Computational Polymers Science: Physics, Chemistry and Biology

Guest Editor:

Dr. Hector Eduardo Roman

Department of Physics, University
of Milano-Bicocca, Piazza della
Scienza 3, 20126 Milano, Italy

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Message from the Guest Editor

The theoretical descriptions should be illustrated by discussing suitable applications to specific problems related to structural, transport and dynamical properties of linear polymers, modelled, e.g., by worm-like (reptation) and Monte Carlo-type methods. The latter are suitable for generating long and densely packed self-avoiding chains in different problems, such as the growth of thin polymeric films of nanometre size, their transport behaviour in disordered environments, and their actual packing within confined volumes such as the nucleus of a cell. The studies of polymer networks, and associated anomalous rheological properties, are welcome in view of the possible connections with fractal scaling and fractional dynamics.





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Editor-in-Chief

Prof. Dr. Alexander Böker

Lehrstuhl für Polymermaterialien
und Polymertechnologie,
University of Potsdam, 14476
Potsdam-Golm, Germany

Message from the Editor-in-Chief

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Polymers Editorial Office
MDPI, Grosspeteranlage 5
4052 Basel, Switzerland

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