



Research on Polymer Simulation, Modeling and Computation

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Message from the Guest Editors

In silico simulation, modeling, and computation have evolved into powerful tools to reveal the molecular mechanisms underlying the macroscopic phenomena and behavior of polymers, predict their physicochemical properties, and discover and design next-generation polymeric materials. A wide range of simulation methods and packages, from those based on quantum mechanics with subatomic resolutions to continuum frameworks dealing with bulk materials, are at the disposal of researchers to study polymers over a full spectrum of length, time, and energy scales under various conditions. Multiscale models are also in rapid development and validation. This Special Issue aims to serve as a platform to allow polymer researchers to exchange exciting results, recent progress, and emerging ideas on understanding polymers from the perspectives of simulation, modeling, and computation. The issue welcomes reports and reviews covering any aspect of polymer modeling, using methods including but not limited to density functional theory, molecular dynamics simulation, coarse-grained modeling, lattice Boltzmann simulation, self-consistent field theory, multiscale simulation.





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Message from the Editor-in-Chief

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