

Simulations of Polymer-Grafted Nanoparticle Phase Behavior

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The phase behavior of polymer-grafted nanoparticles can be computationally challenging to obtain using particle-based simulations, especially with long polymers. The theoretically-informed Langevin dynamics method enables the simulation of large collections of nanoparticles grafted with polymers much longer than those typically accessible by molecular dynamics simulations, and yet also includes fluctuations which are neglected in typical self-consistent mean field methods. The temperature-concentration phase diagram of polymer-grafted nanoparticles in cyclohexane is determined using this relatively new method, by explicitly simulating the phase boundary between particle-rich and particle-poor phases. The phase behavior depends on nanoparticle size, graft degree of polymerization and graft density.

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