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Numerical Coarse-Graining of Polymer Field Theories MICHAEL VILLET, GLENN FREDRICKSON, University of California, Santa Barbara — Field theoretic models of polymers are widely used to investigate polymer selfassembly, but numerical simulations of these models that include full fluctuation physics are computationally demanding and infrequently conducted. To reduce this computational cost, we propose the use of systematically coarse-grained field theories that can be simulated on a coarsely spaced lattice without truncation of important short-wavelength physics. We present a variational method for numerically executing this coarse-graining, in which fine-grained simulation data is used to parameterize trial coarse-grained models, and results from the application of this method to some model systems.

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