



**Recent Developments
in Computer Simulation Studies
in Condensed Matter Physics**

February 20-24, 2017

This annual workshop series highlights recent advances in applications, algorithms, and parallel implementations of computer simulation methods for the study of condensed matter systems. Topics of interest include Monte Carlo, molecular dynamics, and other numerical studies of such physical problems as materials growth, structural and magnetic phase transitions, polymers, surfaces, nanostructures, strongly correlated electron systems and models of exotic quantum phases. Other areas of interest include interfaces, granular flow and other non-equilibrium systems, genomics, membranes and protein folding, free energy determinations, and novel simulation algorithms. Graduate student participation is encouraged.

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