



Library Indian Institute of Science Education and Research Mohali



DSpace@IISERMohali / Thesis & Dissertation / Master of Science / MS-18

Please use this identifier to cite or link to this item: <http://hdl.handle.net/123456789/5559>

Title:	Understanding Granular System Dynamics using Molecular Dynamic Simulation Techniques
Authors:	Awasthi, Rohan
Keywords:	Dynamic Simulation
Issue Date:	May-2023
Publisher:	IISER Mohali
Abstract:	This thesis aims to understand the dynamics of granular systems using molecular dynamic simulation techniques. Granular systems are collections of particles with sizes ranging from a few micrometers to a few centimeters, and they interact via dissipative mechanisms such as inelastic collisions and frictional forces. The first part of this study focuses on the decay of kinetic energy, the velocity distribution of grains, and the density and velocity fields as a function of friction coefficient for unperturbed systems. In the second part, the systems are perturbed by applying homogeneous and inhomogeneous driving, and their steady-state properties are analyzed using analytical tools. Through this investigation, a deeper understanding of the dynamics of granular systems is gained, which can have applications in fields such as materials science, engineering, and geophysics.
Description:	Embargo Period
URI:	http://hdl.handle.net/123456789/5559
Appears in Collections:	MS-18

Files in This Item:

File	Description	Size	Format	
Need To Add...Full Text_PDF		15.36 kB	Unknown	View/Open

Show full item record



Items in DSpace are protected by copyright, with all rights reserved, unless otherwise indicated.