- Simulation Group

01. Molecular dynamics study of transport coeffients for hard-chain fluid
02. Computational physics from Imperial College of Science London
03. Monte Carlo simulation Syracuse university
04. Project MD group
05. Center for material simulation
06. Molecular dynamics
07. CASINO(monte CArlo Simulation of electroN trajectory in sOlids)
08. Northup computational group
09. Material and process simulation center
10. Molecular simulation of liquid water
11. Network science(molecular modeling and simulation of surface)
12. Molecular simulation laboratory
13. Computational material Science
14. Molecular dynamics simulation
15. Theoretical chemistry of SNU
16. Classical molecular dynamics simulation
17. Raymond Kapral-chemical physics theory group
18. University of Durham, science Lab.
19. Computer chemistry center
20. The geometry calculation center