Lammps - Molecular Dynamics Simulations

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(Dated: January 1, 2020)

Abstract

Working with lammps!

I. INTRODUCTION

II. GENERAL INFO

- Project website:
- Language reference:
  - http://lammps.sandia.gov/doc
  - github: https://github.com/lammps/lammps/
- Tutorials
  - Simple tutorial: lammps for dummies
  - Tutorial: tutorial 1
  - Examples: tutorial 2
  - Good for input file: slide 40 on!
- making videos: see linux-docs.pdf

Acknowledgements: We thank .... PVS acknowledge financial support from the German DAAD (grant 57314018) and DFG (grant 4056192179), respectively.

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