Advanced Workshop on High-Performance & High-Throughput Materials Simulations using Quantum ESPRESSO | (smr 3102)

Contribution ID: 1269 Type: not specified

Hands-on: Ab-initio Molecular Dynamics: computing correlations and thermodynamic properties, using thermostats

Wednesday, 18 January 2017 16:30 (2:00)

Content

Summary

Presenter(s): GIANNOZZI, P. (University of Udine)Session Classification: Day 3 -Wednesday, January 18