

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

Contribution ID : 1268

Type : **not specified**

**Hands-on: Ab-initio Molecular Dynamics: starting
the dynamics, computing averages**

Wednesday, 18 January 2017 14:00 (2:00)

Content

Summary

Presenter(s) : GIANNOZZI, P. (University of Udine)

Session Classification : Day 3 -Wednesday, January 18