

Constructing interatomic potentials from first principles using machine learning: the example of tungsten.

Wednesday, 5 November 2014 10:45 (0:25)

Content

Summary

Primary author(s) : CSÁNYI, Gábor (University of Cambridge, U.K.)

Presenter(s) : CSÁNYI, Gábor (University of Cambridge, U.K.)

Session Classification : Session