Table of contents

Tuesday 17 May 2022 1

Machine Learning in Electronic Structure and Molecular Dynamics | (smr 3702)

Tuesday 17 May 2022

SESSION 4 - (13:30-17:30)

time	title	presenter
13:30	Combining the Power of High-Throughput ab initio Calculations and Machine Learning towards Materials Informatics	RIGNANESE, Gian-Marco
14:30	The NOMAD Artificial Intelligence Toolkit: Turning Materials-Science Data into Knowledge and Understanding	GHIRINGHELLI, Luca
16:00	Coffee break	
16:30	Complex Chemistry	PARRINELLO, Michele