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

Thursday 19 May 2022

SESSION 8 - (13:30-17:30)

time	title	presenter
13:30	Neural Network Interatomic Potentials to Explore the Crystal Structure Landscape	DE GIRONCOLI, Stefano
14:30	Projects	
16:00	Coffee break	
16:30	Neural Network Interatomic Potentials with Long Range Interaction	KUCUKBENLI, Ermine