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

Monday 16 May 2022

SESSION 1 - (09:00-15:00)

time	title	presenter
09:00	Intro to Python	AKIN-OJO, Omololu
10:00	Recap of Unsupervised ML	LAIO, Alessandro
11:00	Coffee break	
11:30	Unsupervised Methods in Machine Learning	LAIO, Alessandro
12:30	Break	
13:00	Machine Learning at the Atomic Scale: from Structural Representations to Chemical Insight	CERIOTTI, Michele
14:00	Lunch break	