Contribution ID: **10** Type: **not specified**

Large-scale GROMACS biomolecular simulations using AWS Batch

Wednesday, December 8, 2021 3:15 PM (1 hour)

How biomolecular simulations scale on AWS CPU and GPU instances across regions using GROMACS through AWS Batch

Presenters: KNIEP, Christian (Sr. HPC/Batch Developer Advocate, AWS); Dr KUTZNER, Carsten (Max Plank Institute)