2/11/2020 Week 5 Module 1
Interactive Exercise for Molecular Dynamics

- In this module, you will learn how to
  - prepare a python script to run OpenMM using the OpenMM script builder (http://builder.openmm.org)
  - navigate the UNIX terminal
  - execute programs with a command line interface (CLI)
  - run a short MD simulation on your desktop machine
  - visualize the simulation with VMD

Preparing a script to run MD