



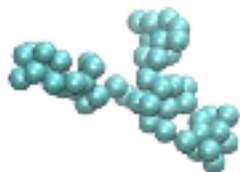


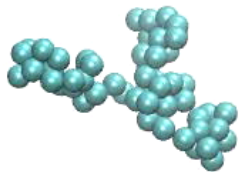
Compute different conformations
by Langevin molecular dynamic
simulation.

(see upcoming paper)

Coarse graining

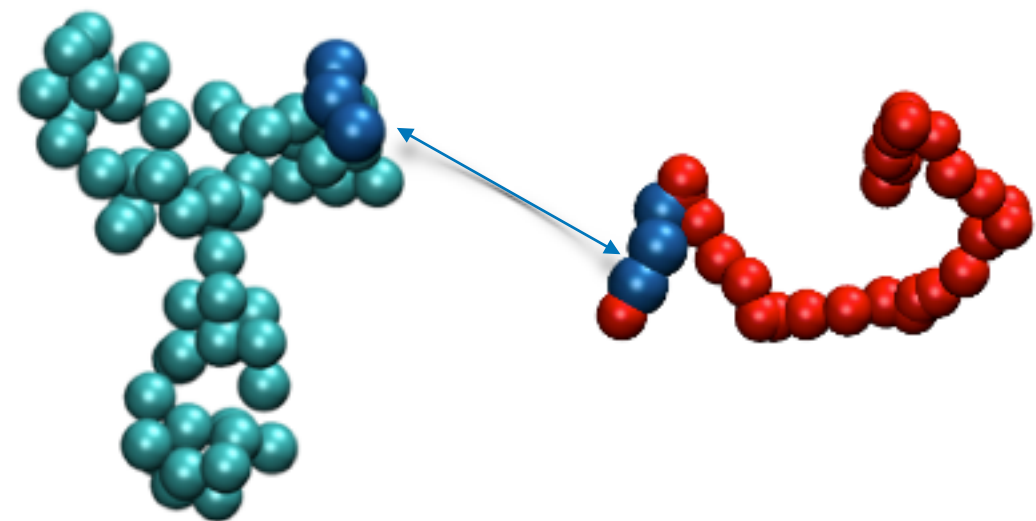
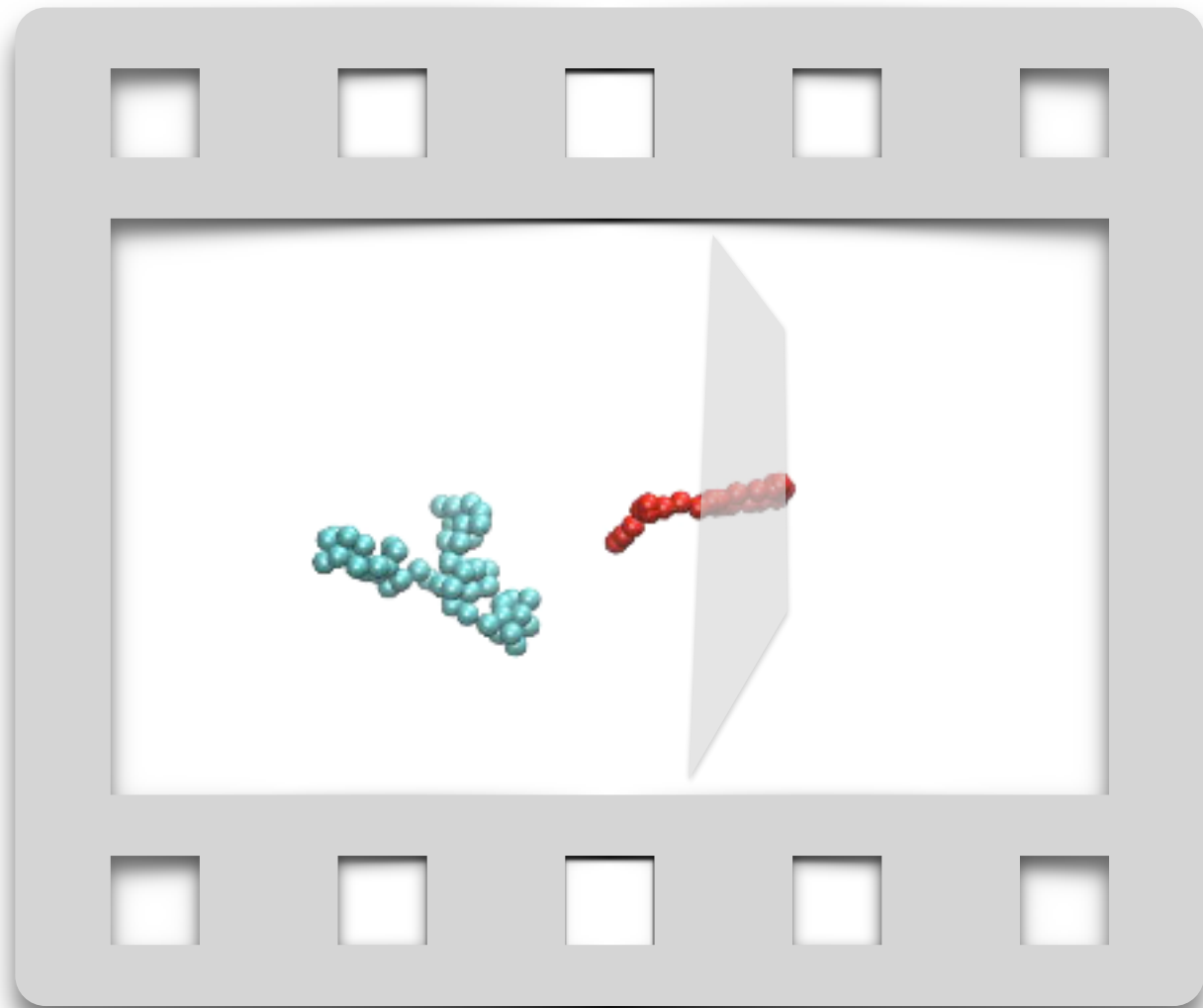
Need for speed, and the limited resolution





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