What can be done with VMD?

- Most of the “Gallery” graphics
  - the Theoretical and Computational Biophysics Group at UIUC (https://www.ks.uiuc.edu/Gallery/Science-New/)
  - on my group’s web page (http://mypages.iit.edu/~dminh/Computational_Chemical_Biology/Gallery.html)
- Scripting makes certain tasks possible

- Today we will go through the official VMD tutorial on
  - Working with a Single Molecule (http://www.ks.uiuc.edu/Training/Tutorials/vmd/tutorial-html/node2.html)
  - Working with Multiple Molecules (http://www.ks.uiuc.edu/Training/Tutorials/vmd/tutorial-html/node6.html)
  - Comparing Structures and Sequences with MultiSeq (http://www.ks.uiuc.edu/Training/Tutorials/vmd/tutorial-html/node7.html)
- The full tutorial
  - http://www.ks.uiuc.edu/Training/Tutorials/vmd/tutorial-html/index.html
  - http://www.ks.uiuc.edu/Training/Tutorials/vmd/vmd-tutorial.pdf