Steps for cluster generation

- 1. Make folder for project. For example, make a folder names cluster_MSA_afv2
 - Mkdir cluster_MSA_afv2
- 2. Upload fasta file to "cluster_MSA_afv2"
 - 1. Vi test models.fasta
- 3. Upload the python scripts (pipeline_CFB) to current folder
 - 1. Scp -r Downloads/pipelineCFB vmischley@login.expanse.sdsc.edu:/expanse/lustre/projects/was136/vmischley/cluster_MSA_afv2
- 4. Move the submit script for MSA generation (CFS_submit_script.slurm) to folder that you want the MSA folder to be created
 - Cp pipeline_CFB/CFS_submit_script.slurm ./
- 5. Edit CFS_submit_script.slurm. Change the name of the fasta file and change the name of the MSA folder. For example, change fasta file to test_models.fasta
 - 1. Vi CFS_submit_script.slurm
- 6. Run CFS_submit_script.slurm. This generates the MSAs. It will take between 24-48 hours. This will generate an MSA folder.
 - Sbatch CFS_submit_script.slurm
- 7. Once the MSA folder is generated, run run_batch_pipeline.py . This will rename the MSAs and will split the MSAs into jobs that will take 48 hours to complete. It will also create the submit scripts for you. Three arguments: location of MSA folder (use full path) number of recycles and version of AF.
 - 1. Python pipeline_CFB/run_batch_pipeline.py /expanse/lustre/projects/was136/vmischley/cluster_MSA_afv2/test_models_MSA 12 alphafold2_multimer_v2
- 8. Change into the model folder where all of the submit scripts are
 - 1. Cd models
- 9. Submit all of the submit scripts
 - 1. Sbatch 1_CFB_test_models_MSA_1_submit.slurm
- 10. After completion, download folder
- 11. Run move af.py to copy only pdb files and .json files into one folder (optional)