## **UF HPC Training**

## NGS Mapping and Assembly February 20, 2014

- 1. Log into UF HPC's Galaxy instance: http://galaxy.hpc.ufl.edu/
- 2. Get some data:
  - a. Shared Data: Data Libraries: Training datasets: wine yeast.100K.fg
- 3. NGS: QC and manipulation: FASTQ Groomer
  - a. Input FASTQ quality scores type: Sanger
- 4. NGS: QC and manipulation:FastQC: FastQC:Read QC
  - a. Use the defaults or add a title for easier reference later
  - b. Notice poor quality at ends of reads
- 5. NGS:QC and manipulation:FASTQ Quality Trimmer
  - a. Window size: 5
  - b. Quality score: 20
  - c. Rerun FastQC on trimmed dataset
- 6. NGS: Mapping: Map with Bowtie for Illumina
  - a. Use a built-in index:
  - b. Select: S. cerevisiae (CGD) 2011
- 7. NGS:SAM Tools:**SAM-BAM** 
  - a. Convert your SAM file to BAM using the defaults
- 8. Click on the View in Trackster icon in the BAM results window
- 9. Select a chromosome and see where reads mapped

## From the command line:

- 1. Login to HiPerGator: ssh <user>@gator.rc.ufl.edu
- 2. Go to your scratch space, make a directory called bowtie test and cd into it
  - a. cd /scratch/lfs/\$USER
  - b. mkdir bowtie test
  - c. cd bowtie test
- 3. Copy/scratch/lfs/bio/training/2014-02-20/bowtie.pbs to bowtie test
  - a.cp /scratch/lfs/bio/training/2014-02-20/bowtie.pbs .
- 4. Edit the bowtie.pbs file to have your e-mail
  - a. nano bowtie.pbs
- 5. Submit the Bowtie run
  - a.qsub bowtie.pbs
- 6. Now let's look at Velvet: make a directory called velvet test and cd into it
  - a. cd .. (moves you up one directory)
  - b. mkdir velvet test
  - c. cd velvet test
- 7. Copy /scratch/lfs/bio/training/2014-02-20/velet.pbs to velvet\_test
  - a.cp /scratch/lfs/bio/training/2014-02-20/velvet.pbs .
- 8. Edit the velvet.pbs file to have your e-mail.

- a. nano velvet.pbs
- 9. Submit the velvet run
  - a. qsub velvet.pbs
- 10. Compare the resulting contig files

## Contents of velvet.pbs:

```
Here are the PBS directives, the information for the
#!/bin/bash
                                scheduler:
#PBS -N velvet
#PBS -M <your e-mail>
                                In addition to CPUs, RAM and walltime, this has
#PBS -m abe
                                information for log files, and e-mail notification.
#PBS -o velvet.test.out
#PBS -e velvet.test.err
                                PBS directive lines start with #PBS and should be at the
#PBS -l nodes=1:ppn=4
#PBS -1 pmem=900mb
                                top of the file
#PBS -1 walltime=00:05:00
                           Remember to run out of scratch space—the command, cd
cd $PBS O WORKDIR
                           $PBS O WORKDIR, changes from home to where you typed
                           gsub. This should be part of most scripts you make.
module load velvet
                           Note loading of the module for the application we are running,
# Make and output
                           use the module system to save headaches!
# directory for Velvet
mkdir test_run
#Run Velvet with kmer of 21
                                                                  Run velvet once
velveth test run/ 21 -fastq -short \
                                                                  using kmer of 21
   /scratch/lfs/bio/training/2014-02-20/wine yeast.100k.fq
velvetg test run/ -min contig lgth 500
#Get things ready to use threaded (OMP) version of Velvet
                                                                Note that for the
#Set OMP THREAD LIMIT--should be the same as ppn above
                                                                threaded version of
export OMP THREAD LIMIT=$PBS NP
                                                                Velvet, you need to set
#Set OMP_NUM_THREADS--should be 1 lower than ppn
                                                                some environment
NUM THREADS=$((PBS NP-1))
                                                                variables
export OMP NUM THREADS=$NUM THREADS
echo Limiting Velvet to $PBS_NP threads total with $NUM_THREADS slave threads.
# Rerun Velvet using a kmer of 51, and the threaded version
# Note there isn't a flag to tell Velvet how many threads to use
# It will use all the cores on a node unless you tell it not to with
# $OMP THREAD LIMIT and $OMP NUM THREADS
                                                                Run velvet again,
mkdir test_run_kmer51
                                                                this time using
velveth max99 OMP test run kmer51/ 51 -fastq -short \
                                                                multiple CPUs,
  /scratch/lfs/bio/training/2014-02-20/wine yeast.100k.fq
```

**Note that this script runs Velvet twice as an example**. You would not typically want to do this... *Either* run on a single core, like the first time through, and adjust resource requests to nodes=1:ppn=1, *or* run on multiple cores, and set OMP\_THREAD\_LIMIT and OMP\_NUM\_THREADS as in the example.

and kmer of 51

velvetg\_max99\_OMP test\_run\_kmer51/ -min\_contig\_lgth 500