Working on UC Davis Bioinformatics Core Administrated Computing Systems

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Outline

- Introduction/Questions
- Explain user space (home directories) vs share space (lab directories/shares).
- Working in a team, file/folder permissions.
- The module system, when to request software, when to install software yourself.
- Slurm

Quick Introduction to the Bioinformatics Core Introduction to the UC Davis

The **mission** of the Bioinformatics Core facility is to facilitate outstanding omics- scale research through these activities:

Data Analysis

The Bioinformatics Core promotes experimental design, advanced computation and informatics analysis of omics scale datasets that drives research forward.

Research Computing

Maintain and make available high-performance computing hardware and software necessary for todays data-intensive bioinformatic analyses.

Training

The Core helps to educate the next generation of bioinformaticians through highly acclaimed training workshops, seminars and through direct participation in research activities.

UC Davis Bioinformatics Core in the Genome Center Core Facility Manager Faculty Advisor

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Research

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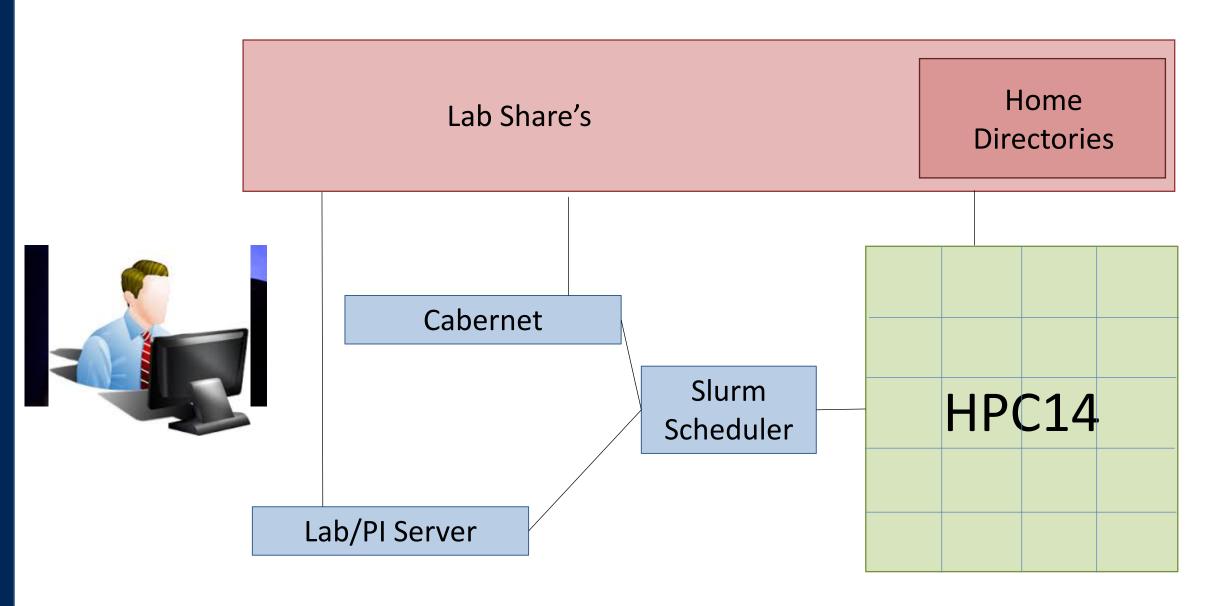
Contacts

Computing Issues, include but not limited to
 User account questions, equipment failure/malfunction, software install,
 software failures (not related to use)

helpdesk@genomecenter.ucdavis.edu

 Bioinformatics related questions, include but not limited to bioinformatic methods questions, software use, data questions

Bioinformatics.core@ucdavis.edu



User space vs shared space AND Working in a team, file/folder permissions. Groups/Labs, etc.

What is a user account?

- User account are not tied to PI/labs. Any personal who has a legitimate reason for requesting an account will be provided one.
 Users should only ever login as themselves and never as another user.
- The Bioinformatics Core provides users with a home directory that is capped at 100Gb (we are exploring reducing that to 10Gb).
- User home directories are 'personal' space, special permission must be given by university administration to explore a user's home directory.
- You now presumably can login to cabernet, but because your not yet associated with a group you can't do much. Never ever work on cabernet.

Lab Groups

- The UC Davis Bioinformatics Core manages lab group data shares, top level folder associated with a collective group. These are data storage resources purchased by a lab, or PI.
- Groups/Labs/Collaborators work together in lab shares, found under /share/group_name
 - [ex. The Bioinformatics Core share is /share/biocore]
- All lab associated analysis should be performed in lab shares. Virtually nothing should be done in your home directory, I personally have:
 - A link to my labs share
 - Software I've installed for me

Linux permissions

- Users on linux system have a user id and group(s) membership(s)
 - to see your group memberships on the command line type: groups
 - You may see an 'error' groups: cannot find name for group ID 1107119953 you can safely ignore this
- Access to additional resources [data/compute] are driven by group permissions/memberships.
- More information on linux permissions

https://www.linux.com/learn/understanding-linux-file-permissions

By default

- By default, you have sole access to your home directory by both user and group.
- By default the lab share is owned by the PI and the group may be the PI group id, or a special group id [ex. subproject].
- By default, all shares are read accessible by everyone and write accessible by the user and the group.
- By default, the group id on a share is a 'sticky' bit, this means that any files/folders created under that share folder inherit the same permission (group read/write) but are owned by the user which created them.

Working as a team

- Groups and permissions are setup to facilitate team/group interactions.
- When working on files shared with other Users you don't have to request the other user to modify permissions, the group sticky bit takes care of that for you.
- YOU DO HAVE TO BE CAREFUL of others work! Since you have group write ability you are able to modify, delete and otherwise impact others people work/project.
- Once a file/folder has been created you can modify the permission to restrict access, however should the PI request the bioinformatics core to change it back, we will.

Linux Permissions

Last thing about permissions, I like to be very permissive, so in addition to the lab share group sticky bit, I also set, in my ~/.bash_profile

umask 0002

By default all files/folders I create are group read/write

Centralized Resources

- Genomes some genomes and mapping indexes (bwa/bowtie) are available centrally to be used by everyone, they can be found under
 - /share/genomes
 - Ensembl
 - genome_reference_consortium
 - Igenomes
 - Kraken
 - ncbi
 - ncbi_genomes
 - patric

Module system

For Software

Modules

- The Environment Modules package provides for the dynamic modification of a user's environment via module files.
- Each module file contains the information needed to configure the shell for an application. Once the Modules package is initialized, the environment can be modified on a per-module basis using the module command which interprets module files. Typically module files instruct the module command to alter or set shell environment variables such as PATH, MANPATH, etc. module files may be shared by many users on a system and users may have their own collection to supplement or replace the shared module files.
- Modules are useful in managing different versions of applications. Modules can also be bundled into meta-modules that will load an entire suite of different applications, or ensure dependencies are loaded.

Modules

- Users can manage their environment (software available on the path) via modules.
- Users benefit from a familiar applications environment across UC Davis Bioinformatics Core systems (on and off-campus).
- Requests for software from any user becomes available to all users.
- Efficient system administration pooling software install/testing efforts from different projects/machines.
- Additional information <u>https://en.wikipedia.org/wiki/Environment Modules (software)</u>

Module basics

module avail, show, load, list, unload

- \$ module avail # shows the currently available module to the system # format is application/version
- \$ module show beast/1.8.3 # shows information, such as what-is, other modules loaded and path information
- \$ module load beast/1.8.3 # load the module beast version 1.8.3
- \$ module list # show currently loaded modules
- \$ module unload beast # unload the module beast

An Example

We can now use beast

```
$ module list # no modules are loaded by default
$ module avail # lets see what modules are present, lets load beast
$ which beast # currently unavailable
$ module show beast/1.8.3 # show its load parameters
$ module load beast/1.8.3 # load beast v. 1.8.3 into our environment
$ module list # show the modules currently loaded
```

\$ which beast # now we see beast is available on our path

Modules are great for

- Modules are great for keeping track of what software and specifically what version of the software is being used in an analysis.
- You can keep track by
 - Doing an analysis from a bash script, loading the modules at the beginning of a script.
 - Copy/paste recording all commands to a file (ex. Analysis.txt) to save the commands for review later
 - Use history at the end of an analysis and save all commands including module load commands
- Note of caution: There are some applications available as a part of the system (ex. samtools) we are working to identify and remove these.

When to install software or request a module

- Benefits to having software in a module:
 - System administrators install, with troublesome software or difficult dependencies they can solve issues others can't.
 - Modules are available to all users, to so request benefits many
- Difficulties with software in a module:
 - Can take time for a system administrators to install, we manage over 400 servers and 5 clusters, we do our best to be responsive.
 - If you're just trying something out that may or may not be useful can waste valuable system administrator time.

When to install software or request a module

- If you feel the software will be broadly applicable, request a module
- If it's a tough install, can't get it installed yourself, request a module
- If you just want to try something out, install it yourself
- If you need it right now, install it yourself

• If you installed it yourself AND requested a module, once the module is available delete your instance.

Slurm scheduler

SLURM

Simple Linux Utility for Resource Management

- Manages compute resources
- Schedules jobs using those resources
- Original developed at Lawrence Livermore National Labs
- Open Source
- Supports plugins extending or enhancing functionality
- Increasingly being used at academic research computing centers and national labs
 - Berkeley
 - UCSD Super Computer

Slurm Commands

sinfo, squeue, sacct, srun, sbatch, scancel

sinfo reports the state of partitions and nodes managed by Slurm. It has a wide variety of filtering, sorting, and formatting options.

squeue reports the state of jobs or job steps. It has a wide variety of filtering, sorting, and formatting options. By default, it reports the running jobs in priority order and then the pending jobs in priority order.

sacct is used to report job or job step accounting information about active or completed jobs.

srun is used to submit a job for execution or initiate job steps in real time. srun has a wide variety of options to specify resource requirements. A job can contain multiple job steps executing sequentially or in parallel on independent or shared resources within the job's node allocation.

sbatch is used to submit a job script for later execution. The script may contain one or more srun commands to launch parallel tasks.

scancel is used to cancel a pending or running job or job step. It can also be used to send an arbitrary signal to all processes associated with a running job or job step.

Slurm Commands - SINFO

- **sinfo** reports the state of partitions and nodes managed by Slurm. It has a wide variety of filtering, sorting, and formatting options. sinfo --summarize
- Node
- Nodes are computers/servers, multiple cpu with shared memory
- Partitions are logical groups of nodes (possibly overlapping)
 - gc64, gc128, gc256, gc512
 - Partition gc is the collection of gc64, gc128 and gc256 and is the default queue

Slurm Commands – SQUEUE

- **squeue** reports the state of jobs or job steps. It has a wide variety of filtering, sorting, and formatting options. By default, it reports the running jobs in priority order and then the pending jobs in priority order.
- Use squeue to view jobs currently queued and running
 - --user=username
 - --states=[running or pending]
 - --long # long format
 - --start # sometimes gives approximate start time

Slurm Commands -SBATCH

- sbatch is used to submit a job script for later execution [non-interactive].
 The script may contain one or more srun commands to launch parallel tasks.
- Login to cluster
- 2. Create a job script
 - I use a .slurm extension so that I know it's a job script
- 3. Submit the script to the cluster using sbatch
 - sbatch script.slurm
- 4. Slurm will respond with a jobid:
 - Submitted batch job 311533
- See example on github

Example Job script

```
#!/bin/bash
#SBATCH --partition=gc # partition to submit to
#SBATCH --job-name="myjob" # Job name
#SBATCH --nodes=1 # single node, anything more than 1 will not run
#SBATCH --ntasks=1 # equivalent to cpus, stick to around 20 max on gc64, or gc128 nodes
#SBATCH --mem=1G # memory pool all cores, default is 2GB per cpu
#SBATCH --time=0-01:30:00 # expected time of completion in hours, minutes, seconds, default 1-day
#SBATCH --output=my output.out # STDOUT
#SBATCH --error=my output.err # STDERR
#SBATCH --mail-user=you@gmail.com
#SBATCH --mail-type=ALL
# This will be run once for a single process
/bin/hostname
module load blat/v.35-fastq
echo "HELLO"
```

Slurm Commands — SCANCEL

- scancel is used to cancel a pending or running job or job step. It can also be used to send an arbitrary signal to all processes associated with a running job or job step.
- To delete a job
 - scancel <jobid>
 - Jobs not running yet (queued) are deleted, jobs running will be killed
- To delete many jobs
 - scancel <jobid1> <jobid2> <jobid3> ...

Slurm Command – srun and helper scripts

srun is used to submit a job for execution or initiate job steps in real time. srun has a wide variety of options to specify resource requirements. A job can contain multiple job steps executing sequentially or in parallel on independent or shared resources within the job's node allocation.

An interactive session (preferred way)

- slurmlogin
 - srun -N 1 -n 1 --pty /bin/bash
- slurmlogin-wholenode
 - srun -N1 --exclusive --pty /bin/bash
- Slurmlogin-gc512
 - srun -p gc512 --mem 500000 -N1 --exclusive --pty /bin/bash
- If using 'screen',
 - 1. 'screen' on Cabernet,
 - 2. then srun,
 - 3. then exit when done.

Task Arrays

Job arrays offer a mechanism for submitting and managing collections of similar jobs quickly and easily; job arrays with millions of tasks can be submitted in milliseconds (subject to configured size limits).

- Slurm parameter, --array=a-b, where a and b are numeric number ex. 1-5
- Sets the environment variables
 - **SLURM_ARRAY_JOB_ID** will be set to the first job ID of the array.
 - **SLURM_ARRAY_TASK_ID** will be set to the job array index value.
- The value of \$SLURM_ARRAY_TASK_ID can be used to reference rows in a file containing values to use for that job
- See example on github

Pipelines – using dependency

- sbatch returns the jobid after submission
- Can use that jobid and the sbatch argument --dependency to launch another job once the first one finishes

```
--dependency=<type:job_id[:job_id][,type:job_id[:job_id]>
```

See example on github

afterok:jobid[:jobid...]

after:jobid[:jobid...]job can begin after the specified jobs have startedafterany:jobid[:jobid...]job can begin after the specified jobs have terminatedafternotok:jobid[:jobid...]job can begin after the specified jobs have failed

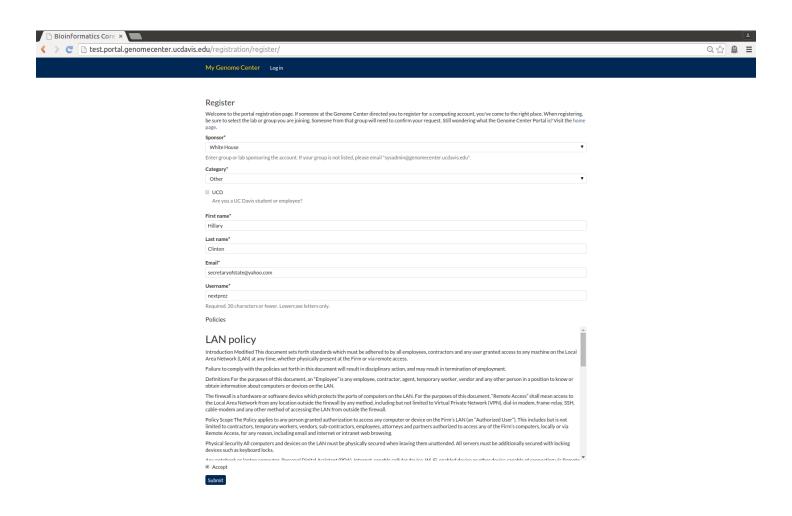
job can begin after the specified jobs have run to completion with an exit code of zero (see the <u>user guide</u> for caveats).

jobs can begin execution after all previously launched jobs with the same name and user have ended. This is useful to collate results of a swarm or to send a notification at the end of a swarm.

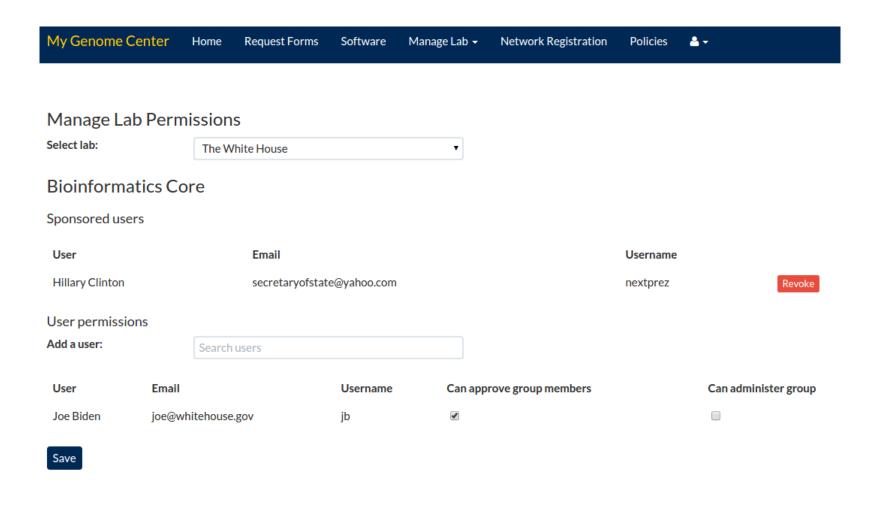
Coming soon

- HPC16
- UC Davis Bioinformatics Core Portal
 - Request users accounts
 - Manage groups/labs/access
 - View available modules and info
 - Request modules and see status of request
 - Request Sys Admin Support

Account registration



Group memberships



Request help

My Genome Center	Home	Request Forms	Software	Manage Lab →	Network Registration	Policies	4+
Submit a ticket							
Job id*							
Server*							
Command*							
Directory*							
Description*							
Submit							w w

Software Request

My Genome Center	Home	Request Forms	Software	Manage Lab ▼	Network Registration	Policies	A +
Request software	9						
Name*							
Version*							
Link*							
Description*							
Tags*							
modules data analys	sis 🗌 libr	aries 🗌 aligner 🕻	assembler	cluster dat	abases 🗌 patterns		
Submit							

Software List

My Genome Center	Home Request Forms	Software Manage Lab ▼	Network Registration	Policies	Av
Software					
Search:					
Clear search					
	gff libraries luajit n		uence U simulator U aligno	er □ assembl	ler 🗆 cluster 🗆 data analysis 🗀 epigenomic
Don't see the software you	need? Request new software t	to be installed on the cluster.			
R Versions: .3.1.0.swp, 3.0.2, l Tags:	ocal_libs, .3.1.0.swo				
abyss Versions: 1.5.2, 1.9.0		allpathslg Versions: 51298		amos Versions: 3	310
Tags:		Tags:		Tags:	.1.0
anytag Versions: 2.5.2		arachne Versions: 46223		arraymake Versions: 1	
Tags:		Tags:		Tags:	L.I
augustus Versions: 3.0.3 Tags:		bamtools Versions: 2.3.0, .2.3.0.swp, .2 Tags:	2.3.0.swo	bbcp Versions: 1 Tags:	15.02.03.01.1
bcftools		bcl2fastq		beagle	
Versions: 1.2 Tags:		Versions: 1.8.4, 2.16.0.10, 2. Tags:	17.1.14	Versions: 2 Tags:	2.1.2
beast Versions: 1.8.3, 1.8.0, 2.4. Tags:	3	bedops Versions: 2.4.11 Tags:		bedtools2 Versions: 2 Tags:	2.25.0, 2.21.0
bioutils Versions: 1.0.9 Tags:		blast Versions: 2.2.30+, 2.2.29+, 2 2.4.0+_debug Tags:	2.2.31+, 2.4.0+,	Versions: 2 Tags: modu	2.2.26
blat Versions: v.35-fastq, v.35 Tags: data analysis	;	boost Versions: 1.44.0, 1.60 Tags: libraries		bowtie Versions: 1 Tags: aligne	•
bowtie2		bsseeker2		busco	

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