# **CURC Cheat Sheet**

**Guide to CURC Resources** 

## **Accessing CURC**

#### **Get a Research Computing account**

https://bit.ly/3LR8XYi

#### Logging in from a terminal

ssh <username>@login.rc.colorado.edu

#### Windows/Mac Clients:

PuTTY: SSH client for Windows WinSCP: SCP client for Windows

FileZilla: FTP client for Linux, Windows, Mac Xming/XQuartz: X11 server for Windows/Mac

#### Logging in from the CURC Web Portal, OnDemand

- Visit: ondemand.rc.colorado.edu
- Web portal to view, edit, down/upload files, manage and create job templates, and access interactive applications

#### **Accessing different CURC clusters**

module load slurm/summit # Summit module load slurm/blanca # Blanca module load slurm/alpine # Alpine

#### **Accessing Compile Nodes/Jobs**

ssh scompile # Summit compile node
acompile # Start Alpine compile job

### **CURC Resources**

#### **High Performance Computing (HPC) Clusters**

Summit2nd-Gen CURC ClusterBlancaCURC Condo ClusterAlpineNext-Gen CURC cluster

#### Storage

• Cluster Directories

/home (2GB) | Backed up daily
/projects (250GB) | Backed up daily
/scratch/summit (10TB) | No back up, 90 day purge
/rc\_scratch (10TB) | No back up, 90 day purge

- PetaLibrary
  - Paid storage service that supports storage, archive, and sharing of research data
  - Backed up with ZFS Snapshots, snapshots do count against your storage quota

#### Cloud

- Δ\Λ/ς
  - Research Computing Cloud Team managed Amazon Web Servies (AWS) accounts
- CUmulus
  - CURC on-premise cloud service

# **Monitoring Tools**

**Slurmtools**: A module that contains a collection of functions to assess recent usage statistics.

#### module load slurmtools

**XDMOD**: A web portal for viewing metrics at the system, partition, and user-levels.

• Visit: xdmod.rc.colorado.edu

**Curc-Quota**: CLI utility to check the storage capacity of your directories on the CURC system.

curc-quota

## **Data Transfers**

Transferring files between CURC and your system

scp source <uname>@login.rc.colorado.edu:destination scp <uname>@login.rc.colorado.edu:source destination

#### Other transfer options

- Globus: Browser application (Recommended)
- Rsync: CLI sync utility
- Sftp: CLI interactive utility
- · Rclone: CLI cloud transfer utility

### **SLURM**

SLURM is an open-source cluster management and job scheduling system for Linux clusters

#### **SLURM Scheduling**

sbatch <file> | Submits a job script <file> **sinteractive** | Submits interactive job Show job queue for <user>. squeue -u <user> scancel <jobid> Deletes the job with <jobid> scontrol hold <jobid> | Hold job with <jobid> scontrol release <jobid> Release job with <jobid> Cluster status salloc Request new resource allocation srun Launch parallel job step **sacct** | Display job accounting information

#### **#SBATCH Directives**

nodes= <count></count>	Number of nodes
tasks-per-node= <count></count>	Processes per node
ntasks= <count></count>	Total processes
cpus-per-task= <count></count>	CPU cores per process
nodelist= <nodes></nodes>	Preferred nodes
exclude= <nodes></nodes>	Nodes to avoid
time= <min></min>	Time limit; either min or
	dd-hh:mm:ss
mem= <count></count>	RAM per node; e.g. 10G
output= <file></file>	Standard output; defaults to
	slurm-jobid.out if omitted
error= <file></file>	Write standard error to file
array= <arrayspec></arrayspec>	Define job array
gres=gpu: <type>:<count></count></type>	Number of GPUs
mail-user= <email></email>	Email for job alerts
mail-type= <type></type>	Email alert types: BEGIN,
	END, FAIL, REQUEUE, ALL
account= <account></account>	Account to charge
depend= <state>:<jobid></jobid></state>	Job dependency. state =
	after,
job-name= <name></name>	afterok, afterany, afternotok
constrain= <attribute></attribute>	Job name
	Request CPU type: e.g.,
	westmere-ex, sandybridge,
partition= <name></name>	ivybridge, haswell
	Submit to partition: e.g.,
	shas, smem, sgpu, amilan-
	ucb

## **SLURM (Cont.)**

#### **SLURM Environment Variables**

SLURM_JOBID	Job ID
SLURM_SUBMIT_DIR	Job submission directory
SLURM_SUBMIT_HOST	Name of host from which
	job was submitted
SLURM_JOB_NODELIST	Names of nodes allocated
	to job
SLURM_ARRAY_TASK_ID	Task id within job array
SLURM_JOB_CPUS_PER_	CPU cores per node
NODE	allocated to job
SLURM_NNODES	Number of nodes allocated
	to job

### **Software**

Research Computing uses a module system called LMOD to load most software into a user's environment. Most software is not accessible by default and must be loaded in.

#### **LMOD Commands**

module avail	Shows all available module
module load <module></module>	Loads <module> in the</module>
	environment, specify version
	with <module>/version**</module>
module spider <string></string>	Searches for module names
	matching <string></string>
module keyword <string></string>	Searches for modules
	containing <string> in name</string>
	or description
module list	List currently loaded modules
module unload <module></module>	Removes <module> from</module>
	environment
module purge	Removes all modules from
	environment
module save <collection></collection>	Saves currently loaded
	modules to collections
module savelist	Returns all saved module
	collections
module describe	Get modules in a saved
	collection

#### Important notes:

- Modules should be loaded in job scripts, interactive jobs, or on compile nodes only, *not* on one of the login nodes
- \*\*It is highly recommended that modules are loaded using their versions, e.g., module load cmake/3.14.1
- A module's dependencies must be loaded before the module can be loaded.