Research Computing New User Seminar

Shelley Knuth shelley.knuth@colorado.edu

www.rc.colorado.edu

Slides:

https://github.com/ResearchComputing/Final_Tutorials/tree/master/RC_Access_Specific_Topics/New_User_Seminar

Outline

- What is RC?
- Steps to get access to our systems
 - Accounts
 - Two-factor authentication
 - Allocations
 - Logging in
- Navigating our systems
 - Storage spaces
 - Data transfer Globus
 - Software
- Running jobs

What is Research Computing?

- Provide services for researchers that include:
 - Large scale computing
 - Data storage
 - High speed data transfer
 - Data management support
 - Consulting
 - Training
- We are likely best known for:
 - Janus/Summit
 - PetaLibrary

What Would I Use Summit For?

- Research Computing is more than just Summit
- But it is what we are most known for
- So what would you use Summit For?
 - Solving large problems that require more:
 - Memory than you have on your personal computer
 - Cores/nodes/power than you have on your personal computer
 - Large visualization jobs
 - High memory jobs
- Not a place for:
 - Large data storage

Hardware - Summit Supercomputer

- 475 compute nodes (Intel Xeon Haswell)
- 24 cores per node
- 11,400 total cores
- Omni-Path network
- 1.2 PB scratch storage
- GPFS File system



Additional Types of Summit Compute Nodes

- 10 Graphics Processing Unit (GPU) Nodes
 - NVIDIA Tesla K80 (2/node)
- 5 High Memory Nodes
 - 2 TB of memory/node, 48 cores/node
- Phi Nodes (planned summer 2017)
 - 20 nodes
 - Intel Xeon Phi

How To Access RC Resources?

- 1. Get an account
- 2. Set up two factor authentication
- 3. Set up an allocation
 - Don't need for Blanca or PetaLibrary
- 4. Log in
- 5. Create greatness

 After you login, you will need to do many additional things that we will discuss today

Getting an RC Account

- CU Boulder users and affiliates:
- Request an account through the RC Account request portal
 - https://portals.rc.colorado.edu/accounts/account-request/create
- CSU Users:
 - Request an account through CSU
 - Talk to Richard Casey
 - Then get an RC user account
- RMACC Users:
 - We are working on a login system through XSEDE
 - Not operational yet

Setting up Two-Factor Authentication

- Two factor authentication is required to access our system
- Require this to provide an extra level of authentication
- Two methods for achieving this:
 - Duo
 - Access through a smart phone app
 - Vasco OTP (one time password)
 - Physical device

Duo Authentication

- Once you get an account, contact <u>rc-help@colorado.edu</u> to request a Duo invitation
- Once you get the invitation, you'll get a series of steps to complete Duo enrollment
- RC supports Duo "push" and "phone call" for authentication
- Greatly prefer "push"

Vasco Authentication

- Physical device that generates a new password every 30 seconds
- To get the device, go to the IT Service Center or request a time to pick one up from Research Computing
- You will need to show an ID to get a device
- First, register the device
 - Otp.colorado.edu
- You will set a four digit password

Allocations

- You will need an allocation to use any of our resources
 - The exception is if you will ONLY be using the PetaLibrary or Blanca
- Currently, to request an allocation please email rc-help@colorado.edu and ask for a General account
 - Need to provide a few sentences on your project
- In the future, we will have a place on our website to submit a more formal request
- Once you have some benchmarks, you will want to move to a project allocation

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Why Do I Need An Allocation?

- I have an account why do I need an allocation?
 - An account validates you are eligible to use RC resources
 - An allocation allows us to keep track of your use of the system
 - This is important because:
 - We need to make sure we have enough resources to accommodate all of our users
 - Helps for reporting to NSF and the CU Research & Innovation Office
 - Applying for an allocation beyond a general allocation:
 - Gives you higher priority in the system

What is Fair Share?

- Fair share scheduling uses a complex formula to determine priority in queue
- Looks at load for each user and each QOS and balances utilization to fairly share resources
 - Involves historical use by user plus how long job has been in the queue
- System will first look at weighted average utilization of user over last 3-4 weeks
- Then compare it to the fair share target percentage of a user

Fair Share Target Percentage

- The target percentage depends on your priority based on your project proposal
- Everyone not associated with a project shares a target percentage of 13% (20% of the CU fraction)
 - No guaranteed level per user
- If you are under (over) your target percentage (based on a 3-4 week average) your priority is increased (decreased)
- Reminder this all only impacts pending jobs
- If no other pending jobs and enough resources are available then your job will run regardless of your previous usage

Allocations

- Need an allocation? Plan to run on Summit?
- Make a request now!
- Include 2-3 sentences describing your proposed usage
- Email <u>rc-help@colorado.edu</u>

Logging In

- It's important to note that you are NOT logging into any specific resource
 - Summit, etc
- When you log in, you land on our login nodes
- From there, you can access our other resources

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RC Resource Access

- To login to an RC login node: ssh username@login.rc.colorado.edu
- If logging in with Duo, your enter your password as: duo:identikey password
- If logging in with Vasco, your enter your password as:

 Pin+six-digit number on Vasco

Navigating our Systems

- Now that you've logged in, now what?
 - What are the different node types we have?
 - What are the different storage spaces?
 - What should I be putting in these storage spaces?
 - How do I transfer data around?
 - How do I deal with software?

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Different Node Types

- Login nodes
 - Four virtual machines
 - This is where you are when you log in
 - No heavy computation, interactive jobs, or long running processes
 - Script or code editing, minor compiling
 - Job submission
- Compile nodes
 - Where you compile code
- Compute/batch nodes
 - This is where jobs that are submitted through the scheduler run
 - Intended for heavy computation

Storage Spaces

Home Directories

- /home/\$USER
- Not for direct computation
- Small quota (2 GB)
- Backed up

\$PROJECT Space

- /projects/\$USER
- Mid level quota (250 GB)
- Large file storage
- Backed up

Scratch Directory

- /scratch/summit/\$USER
- Much larger depends on system
- Files purged around 90 days

What Belongs Where?

- /home
 - Scripts
 - Code
 - Very small files
 - Inappropriate for sharing files with others
 - Inappropriate for job output
- /projects
 - Code/files/libraries relevant for any software you are installing (if you want to share files with others)
 - Mid-level size input files
 - Appropriate for sharing files with others
 - Inappropriate for job output
- /scratch/summit
 - Output from running jobs
 - Large files
 - Appropriate for sharing files with others
 - THIS IS NOT APPROPRIATE FOR LONG TERM STORAGE

Transferring Data

- Globus is Research Computing's preferred method of data transfer
- Designed with researchers in mind
- End points between computers make for efficient data transfer with an easy to use interface
 - Endpoints are different locations that data can be moved to/from
 - Personal or multi-user
- Rsync and sftp through the login nodes is good for small transfers

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Setting Up Globus

- Create an account at Globus.org
- Make your personal computer an endpoint
- Transfer data
- www.globus.org

We will come back to this at the end if we have time

Software

- Common software is available to everyone on the systems
- Can install your own software
 - But you are responsible for support
 - We are happy to assist
- Research Computing uses modules to manage software
 - You can load modules to prepare your environment for using software
 - Set any environment variables
 - Set environment so application can find appropriate libraries, etc.

Important Things to Know About Modules

- Some modules might require a specific hierarchy to load
 - For some modules, you may need to specify a specific version
 - For example, ml R/3.3.0
 - For other modules, you may be able to be more generic
 - For example, ml matlab
- Some modules may require you to first load other modules that they depend on
- To find dependencies for a module, type module spider <package>
- To find out what software is available, you can type ml avail
- To set up your environment to use a software package, type ml
 <package>/<version>

Job Submission

Running Jobs

- What is a "job"?
- Interactive jobs
 - Work interactively at the command line of a compute node
- Batch jobs
 - Submit job that will be executed when resources are available
 - Create a text file containing information about the job
 - Submit the job file to a queue

Job Scheduling

- On a supercomputer, jobs are scheduled rather than just run instantly at the command line
 - Shared system
 - Jobs are put in a queue until resources are available
- Need software that will distribute the jobs appropriately and manage the resources
 - Simple Linux Utility for Resource Management (Slurm)
 - Keeps track of what nodes are busy/available, and what jobs are queued or running
 - Tells the resource manager when to run which job on the available resources

Partitions and 'Quality of Services'

- There are several ways to define where your job will run
- Partitions (basically a queue):
 - Resources/hardware
- QoS:
 - Tells what the limits or characteristics of a job should be
 - Maximum wall time
 - Number of nodes
- One partition might have multiple QoS
- A QoS might exist on multiple partitions

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Partitions

- By default, jobs will run on the general compute (Haswell) nodes
- Recommend specifying a partition
- Do this using the -p partition flag

Available Partitions

Partition	Description	# of nodes	cores/node	GPUs/node
shas	General Compute (Haswell)	380	24	0
sgpu	GPU- enabled nodes	10	24	effectively 4
smem	High-memory nodes	5	48	0
sknl	Phi (Knights Landing) nodes - [not currently available]	20	68	0

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Quality of Service

QoS	Description	Maxwall	Max jobs/user	Max nodes/user
normal	Default QoS	Derived from partition	n/a	256
debug	For quick turnaround when testing	1 H	1	32
long	For jobs needing longer wall times	7 D	n/a	20
condo	For groups who have purchased Summit nodes	7 D	n/a	n/a

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Useful Slurm Commands - sbatch

- sbatch: submit a batch script to slurm
- You can use a bunch of flag options in a batch script or on the command line
- Useful to put in script so have for future use

Example:

```
sbatch test.sh
```

OR

```
sbatch --partition=shas test.sh
```

http://slurm.schedmd.com/sbatch.html

SBATCH Options

```
#SBATCH <options>
                         sbatch <options>
Allocation:
                 -A=<account no>
Partition:
                 --partition=<partition name>

    Sending emails:

                 --mail-type=<type>
Email address:
                --mail-user=<user>

    Number of nodes: --nodes=<nodes>

    Quality of service: --gos=<gos>

Reservation:
                 --reservation=<name>
Wall time:
                 --time=<wall time>
Job Name:
                 -job-name=<jobname>
```

 FYI: You do NOT actually type <> above – this designates something specific you as a user must enter about your job

Working on Summit

- Right now we are in a transition phase between Summit and Janus
- You should be running all your jobs on Summit
- To ensure this, make sure you load the slurm module on Summit

ml slurm/summit

After you run this command you can run sbatch to submit jobs

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Blanca

- If you are a Blanca user, you need an RC account, but not an allocation
- To run jobs as a Blanca user, once you've logged into a login node, load the Blanca slurm module

ml slurm/blanca

- Only certain users have access to Blanca paid service
- If you are unsure, you can ask your advisor or RC
 - But likely if you are unsure you don't have access

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PetaLibrary

- To access the PetaLibrary, you login in to one our RC's login nodes as normal
- Then you cd to either /work/<groupname> or /archive/<groupname>,
 depending on your PetaLibrary service
 - <groupname> is the name set for your group when you set up the PetaLibrary service
 - You do not include the <>
- Only certain users have access to PetaLibrary paid service
- If you are unsure, you can ask your advisor or RC
 - But likely if you are unsure you don't have access

Practice Examples

Submit Your First Job!

- Submit a slurm job with the following instructions:
- 1. The job should run the Unix "hostname" command
- 2. The job will be submitted from a bash script named hostname_summit.sh
- 3. The job will run on 1 node
- 4. We will request 1 minute wall time
- 5. Run from the debug QOS
- 6. Run on the shas partition
- 7. User the new_user reservation
 - This is only for this workshop

Hostname_summit.sh

```
#!/bin/bash
#SBATCH --nodes=1
                                           # Number of requested nodes
#SBATCH --time=0:01:00
                                           # Max wall time
                                           # Specify debug QOS
#SBATCH --qos=debug
#SBATCH --partition=shas
                                           # Specify Summit haswell nodes
#SBATCH --output=hostname %j.out
                                           # Rename standard output file
#SBATCH --reservation=new user
                                           # Reservation name
# purge all existing modules
module purge
# The directory where you want the job to run
cd /projects/$USER/tutorials/parallelization workshop/new
```

hostname

Running the script

 Load up the slurm module module load slurm/summit

• Submit the job: sbatch hostname_summit.sh

Check output

Another slurm command

squeue

View information about jobs located in the slurm scheduling queue

OPTIONS:

- User: -u <user list>
- Queues: --qos=<qos_list>

EXAMPLE:

http://slurm.schedmd.com/squeue.html

Your turn

- Submit a slurm job with the following instructions:
- 1. The job should run first the whoami command, then the Unix "sleep" command for 30 seconds, then the hostname command
 - Syntax for these Unix commands are below:

whoami sleep 30 hostname

http://slurm.schedmd.com/squeue.html

Your turn

- Submit a slurm job with the following instructions:
- 1. The job will be submitted from a bash script named sleep.sh
- 2. The job will run on 1 node
- 3. Request a 1 minute wall time
- 4. Run the job from the normal QOS
- 5. Run the job from the Summit GPU partition
- 6. Name your job sleep
- 7. Email yourself the results at the end of the job run
 - Hint: Requires two SBATCH options to do this see link at top of this slide
- 8. Use the new_user reservation

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Sleep.sh

```
#!/bin/bash
#SBATCH --nodes=1
                                                # Number of requested nodes
#SBATCH --time=0:01:00
                                                # Max walltime
#SBATCH --qos=normal
                                                # Specify normal QOS
                                                # Specify Summit GPU nodes
#SBATCH --partition=shas
#SBATCH --output=sleep %j.out
                                                # Rename standard output file
#SBATCH --job-name=sleep
                                                # Job submission name
                                                # Email you when the job ends
#SBATCH --mail-type=end
###SBATCH --mail-user=<user>@colorado.edu
                                                # Email address to send to
                                                # Reservation name
#SBATCH --reservation=new user
# purge all existing modules
module purge
# The directory where you want the job to run
cd /projects/$USER/tutorials/parallelization workshop/new
whoami
sleep 30
hostname
```

Running an external script

- Let's run a Matlab program
- We will run the batch script matlab.sh
- This script calls and runs matlab_tic.m

Running the script

Submit the job:

sbatch matlab.sh

Check output

Matlab.sh

```
#!/bin/bash
#SBATCH --nodes=1
                                         # Number of requested nodes
#SBATCH --time=0:02:00
                                         # Max walltime
#SBATCH --qos=debug
                                         # Specify debug QOS
#SBATCH --partition=shas
                                         # Specify Summit haswell nodes
#SBATCH --output=matlab %j.out
                                         # Output file name
#SBATCH --reservation=new user
                                        # Reservation name
# purge all existing modules
module purge
# Load Matlab module
module load matlab
# Run matlab without a GUI
matlab -nodisplay -nodesktop -r "clear; matlab tic;"
```

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Your turn

- Submit a slurm job with the following instructions:
- Create an R program called R program. R that creates a vector called "planets" and then list the planets in the vector
 - Syntax: planets -> planets <- c("Mercury", "Venus", "Earth", "Mars", "Jupiter", "Saturn", "Uranus", "Neptune", "Pluto")
- 2. Print off the vector
 - Syntax: planets
- 3. Create a bash script called R code.sh that runs the R script
 - Syntax: Rscript R code.sh
- 4. The job will run on 1 node
- We will request a 1 minute wall time
- Specify the debug QOS
- 7. Specify the shas partition
- The output will be put in a file called R_code_%j.out
- 9. Use the new_user reservation
- 10. Don't forget to load the R module!

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Solution - R_code.sh

```
#!/bin/bash
#SBATCH --nodes=1
#SBATCH --time=0:01:00
#SBATCH --qos=debug
#SBATCH --partition=shas
#SBATCH --output=R code %j.out
###SBATCH --reservation=new user
# purge all existing modules
module purge
# Load the R module
module load R/3.3.0
# Run R Script
Rscript R program.R
```

```
# Number of requested nodes
# Max walltime
# Specify debug QOS
# Specify Summit haswell nodes
# Output file name
# Reservation name
```

Solution – R_program.R

```
#Simple R code example by Shelley Knuth (shelley.knuth@colorado.edu)
# Create vector
planets <- c("Mercury", "Venus", "Earth", "Mars", "Jupiter", "Saturn", "Uranus",
"Neptune", "Pluto")
# Print off vector
planets
```

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Interactive jobs!

- Sometimes we want our job to run in the background
- Sometimes we want to work in program in real time
- For example, Matlab
- Let's run an interactive Matlab job

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Interactive job

- NEVER EVER EVER EVER EVER
- EVER EVER EVER EVER EVER EVER EVER
- Run Matlab from the command line without running an interactive job first
- BAD USER!

Interactive job

- To do this, we are going to log out and log back in
 - Only necessary for demo
 - Need to add something to the sign in process
- For Mac Users:

```
ssh -X username@login.rc.colorado.edu
```

- For Windows Users, must set up X-forwarding through your SSH client program
- Also must have an X-server package on your laptop
 - Xming for Windows or XQuartz for Mac

Interactive job

- To work with Matlab interactively, we're going to request some time from the supercomputer
- When the resources become available then we will start up Matlab
- Commands to run:

```
module load slurm/summit
sinteractive --reservation=new_user
```

• Once we receive a prompt, then:

```
module load matlab
matlab
```

Once we finish we must exit!

Setting Up Globus

- Create an account at Globus.org
- Make your personal computer an endpoint
- Transfer data
- www.globus.org

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Questions?

- Email rc-help@colorado.edu
- Twitter: @CUBoulderRC
- Link to survey on this topic:

http://tinyurl.com/curc-survey16

Slides:

https://github.com/ResearchComputing/Final_Tutorials/tree/master/RC_Access_Specific_Topics/New_User_Seminar