

CRUK cluster practical sessions (SLURM)

Part I – processes & scripts



login

Log in to the head node, clust I-headnode, using **ssh** and your usual user name & password.

```
SSH Secure Shell 3.2.9 (Build 283)
Copyright (c) 2000-2003 SSH Communications Security Corp - http://www.ssh.com/
This copy of SSH Secure Shell is a non-commercial version.
This version does not include PKI and PKCS #11 functionality.

Last login: Mon Sep 19 10:44:07 2016 from bp7r25j.cri.camres.org
[user@cluster ~]$
```

You're ready to start.

navigate

Find out where you are using pwd.

Make a directory (**mkdir**) and move into it (**cd**)

```
[user@cluster ~]$ pwd
/home/user
[user@cluster ~]$ mkdir training
[user@cluster ~]$ cd training/
[user@cluster training]$
```

processes

You can see your current processes using **ps**.

You can see what else *this* computer is doing using **top**

[user@cluster training]\$ top

top output

top uses the whole screen. Type 'q' to get your screen back.

```
top - 16:26:38 up 58 days, 22:33, 36 users, load average: 0.12, 0.14,
0.12
Tasks: 618 total, 1 running, 617 sleeping, 0 stopped,
Cpu(s): 0.1%us, 0.2%sy, 0.0%ni, 99.5%id, 0.2%wa, 0.0%hi, 0.0%si,
0.0%st
Mem: 16437908k total, 10473016k used, 5964892k free, 2611564k
buffers
Swap: 16779852k total, 162896k used, 16616956k free, 2158536k cached
 PID USER
             PR NI VIRT RES SHR S %CPU %MEM
                                               TIME+ COMMAND
 975 root 0 -20 22712 3832 2196 S
                                      1 0.0 28:44.67 lim
4686 root
            15 0
                       0
                            0
                                0 S 0 0.0
                                              3:11.36 nfsd
           15 0 11048 1592 864 R 0 0.0
                                              0:00.14 top
19175 user
                                              0:12.04 init
   1 root 15 0 10364 600 564 S 0 0.0
```

The 'sleep' command

The **sleep** command doesn't do much – but you can control how many seconds it does it for, and it doesn't use much CPU or I/O

```
[user@cluster training]$ sleep 10
[user@cluster training]$
```

Stop and suspend

If we get bored, change our mind, or think something is wrong we can interrupt jobs. To stop a job, type '^C' at the command line (that's [Ctrl]+[C] together).

```
[user@cluster training]$ sleep 100
[user@cluster training]$
```

If you don't want to stop the job, you can suspend it. Type '^Z' (that's [Ctrl]+[Z]). Type 'fg' to bring the job back to the foreground.

backgrounding

When we have suspended a job (which will never finish). To get it to carry on, we can put it in the 'background' using **bg**

You can put a job in the background deliberately using the '&' character at the end of the command.

Killing processes

If you don't want to wait for it to finish, or think it is broken in some way, you can terminate it using the **kill** command.

Kill has a variety of gentle options to allow the process to exit gracefully. If these fail one – signal -9, or –KILL will normally remove the process.

A simple example

Sleep is a good example, but it doesn't produce any output. We want to wrap it up with messages – in unix you use **echo** to do this.

The colon here allows us to put multiple commands on a single line.

```
[user@cluster training]$ echo start; sleep 1; echo finish
start
finish
[user@cluster training]$
```

Creating a script

Cluster programming makes use of scripts, so we'll turn this list of commands into a script.

You can type directly into a file using **cat** if you know that the end of file character is a '^D'.

```
[user@cluster training]$ cat > script.sh
#!/usr/bin/bash
echo start
sleep 10
echo finish
[user@cluster training]$
```

You can run a script by executing **bash <scriptname>** or by making it directly executable with **chmod**. The './' is important – the shell only looks for executables in certain places – the '**PATH**'.

```
[user@cluster training]$ chmod u+x script.sh
[user@cluster training]$ ./script.sh
start
finish
```

Running the script

Now we are ready to start running our script, or sending it as a cluster job.



Cluster practical sessions

Part II – cluster job submission



Submitting a job

Now we know enough to run our script on the cluster.

Simply submit the job using **sbatch**.

- the output file is written to a Lustre file system directory
- Create directory with username if it doesn't exist – e.g. mkdir /scratcha/stlab/garret0 l
- /home is writeable from cluster nodes, but won't perform as well.

[user@cluster training]\$ sbatch --output=/scratcha/stlab/garret01/%N-%j.out script.sh Submitted batch job 200875

 All read and write operations from within jobs running on nodes should use either /scratchb or /scratcha directories.

Look at running jobs

While the job is running, you can see it with squeue.

```
[user@cluster training]$ squeue
           JOBID PARTITION
                              NAME
                                      USER ST
                                                   TIME NODES NODELIST (REASON)
          200876
                                                   0:02
                  general script.s
                                                             1 clust1-node-3
                                      user R
                                                4:01:05
          200867 general
                              bash sawle01 R
                                                            1 clust1-node-2
                  general MB99.6.v eldrid01 R 9-00:28:46
          175393
                                                            1 clust1-node-30
                 general vardict eldrid01 R 9-01:40:03
          175330
                                                            1 clust1-node-1
```

Once it's finished, you can see the output.

```
[user@cluster training]$ squeue
            JOBID PARTITION
                               NAME
                                                     TIME NODES NODELIST (REASON)
                                        USER ST
           200867 general
                               bash sawle01 R
                                                  4:01:33
                                                               1 clust1-node-2
                   general MB99.6.v eldrid01 R 9-00:29:14
           175393
                                                              1 clust1-node-30
           175330
                   general vardict eldrid01 R 9-01:40:31
                                                              1 clust1-node-1
[user@cluster training]$ ls /scratcha/group/user/
clust1-node-3-200877.out
```

What happened?

The output went into the file as expected:

```
[user@cluster training]$ cat /scratcha/group/user/clust1-node-3-200877.out
start
finish
```

Other information is stored, and available via sacct:

```
[user@cluster training]$ sacct -j 200877

JobID JobName Partition Account AllocCPUS State ExitCode

200877 script.sh general group 1 COMPLETED 0:0
200877.batch batch group 1 COMPLETED 0:0

[user@cluster training]$ sacct -j 200877 --format JobID, MaxRSS, State, AllocCPUS

JobID MaxRSS State AllocCPUS

200877 COMPLETED 1
200877.batch 2012K COMPLETED 1
```

An alternative way to submit

You can submit a job directly to SLURM with **srun**. This still requires resources – it's more commonly used as part of an existing job.

```
[user@cluster training]$ srun /usr/bin/bash script.sh
start
finish
```

You can also use this to generate an interactive session:

```
[user@cluster training]$ srun --pty /usr/bin/bash
[user@clust1-node-3 training]$
```

Killing a job

Just as for processes, but using scancel

```
[user@cluster training] $ sbatch --output=/scratcha/group/user/%N-%j.out script.sh
Submitted batch job 200889
[user@cluster training]$ squeue
                                                          NODES NODELIST (REASON)
            JOBID PARTITION
                                NAME
                                        USER ST
                                                      TIME
           200889
                    general script.s
                                                      0:02
                                                                1 clust1-node-3
                                        user R
           175393 general MB99.6.v eldrid01 R 9-00:56:02
                                                                1 clust1-node-2
                                                                1 clust1-node-30
                    general vardict eldrid01 R 9-02:07:19
           175330
                                                                1 clust1-node-1
[user@cluster training]$ scancel 200889
[user@cluster training]$ squeue
            JOBID PARTITION
                                NAME
                                        USER ST
                                                      TIME NODES NODELIST (REASON)
           200867
                    general
                                bash sawle01 R
                                                   4:28:21
                                                                1 clust1-node-2
                   general MB99.6.v eldrid01 R 9-00:56:02
           175393
                                                                1 clust1-node-30
           175330 general vardict_ eldrid01 R 9-02:07:19
                                                                1 clust1-node-1
```

NOTE: Do not use **skill** it is **NOT** a SLURM command!

Killing isn't bad...

The scheduler manages the shutdown and still records details of the job.

200889 script.sh general group 1 CANCELLED+ 0:
200889.batch batch group 1 CANCELLED 0:1

Basic parallelism

Now we're ready to use the cluster at full power!

One way to do this is with a job array. You can create one of these using the

--array=I-N syntax in sbatch

```
[user@cluster training]$ sbatch --array=1-10 --output=/scratcha/group/user/%N-%j.out script.sh
Submitted batch job 200900
[user@cluster training]$ ls / scratcha/group/user
clust1-node-10-200908.out clust1-node-12-200900.out clust1-node-4-200902.out
clust1-node-10-200904.out clust1-node-12-200906.out clust1-node-11-200909.out
clust1-node-3-200901.out clust1-node-5-200903.out clust1-node-7-200905.out
clust1-node-9-200907.out
```

Or using the **srun** with the **-n** or **-N** parameters.

```
[user@cluster training]$ srun -n srun -n 41 hostname
clust1-node-9.cri.camres.org
...
clust1-node-13.cri.camres.org
[user@cluster training]$
[user@cluster training]$ srun -N 3 hostname
clust1-node-19.cri.camres.org
clust1-node-25.cri.camres.org
clust1-node-8.cri.camres.org
[user@cluster training]$
```



Ende

