



# Introduction to HPC Tutorials

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### **Useful Links**

- HPC Docs ぱ
- Interactive Learning site 🗗

#### Category:

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# **Tutorial 1 - Logging into the HPC**

Objective: To log into the HPC cluster using an ssh client. Prerequisites: A QMUL username and an SSH public/private keypair.

#### Instructions:

- 1. Create an SSH public/private keypair (instructions sent separately).
- 2. Follow the relevant instructions below for your installed operating system to launch a terminal within an ssh client:
  - Windows
  - Linux
  - MacOSX
- 3. Log into the HPC by executing:

ssh abc123@login.hpc.qmul.ac.uk -i /path/to/private\_key

This will attempt a login connection to the HPC cluster at address loqin.hpc.qmul.ac.uk for username abc123.

Note: Replace abc123 with your QMUL username and  $/path/to/private\_key$  with the actual path to your private SSH key.

Note: Your password will not be displayed on the screen as it is typed.

4. Once logged into the HPC, you should see the welcome message and a prompt similar to the following:

[abc123@frontend8 ~]\$

Verify you see the prompt with your username.

#### Windows Users

• Download a suitable SSH client, for example MobaXterm (https://mobaxterm.mobatek.net/download-home-edition.html). Download the portable edition to get started but if you have rights to install software on your machine, you may also select the installer edition.

*Note:* To ensure your private key does not get deleted when your machine is powered off, set a "persistent home directory" from within the MobaXterm configuration.

• Open the MobaXterm application and click the Start local terminal button.





### **Linux Users**

Launch the in-built terminal application:

- Centos7: Applications > System Tools > Terminal.
- Ubuntu: Open the dash menu (Windows key) and type terminal.

### MacOSX Users

 Launch the in-built terminal application by clicking the Finder icon in the dock > Go > Utilities > Terminal.

# **Tutorial 2 - Application Modules**

Objective:	To view, load, unload and list available application modules.
Prerequisite:	An active connection to the HPC cluster.

### Instructions:

1. List all the available software packages by executing:

module avail

2. List the available versions of R by executing:

## module avail R

If an application is not installed, you should see no output - try printing the available versions for the application "not installed".

3. Load the default version of the R application into your environment by running:

## module load R

*Note:* To load a specific version of an application, use the full module name (application name and version).

4. List the currently loaded modules by running:

### module list

The output will now show that R is loaded in your environment. Observe that some applications such as R, or samtools also load other modules as dependencies. These are also unloaded automatically during a module unload of the parent application.

- 5. Run R --version to confirm which version of R has been loaded. This is an inbuilt command within the R application, rather than a feature of modules. Other applications often display the version number in the output from the -h or --help options.
- 6. Now attempt to load the R/3.5.3 module note that a module conflict error appears. This is by design you cannot have two versions of the same application loaded at once, and one needs to be unloaded first.
- 7. Let's unload the existing R module and then load version R/3.5.3:

```
module unload R module load R/3.5.3
```





*Note:* You may instead use the module switch R/3.5.3 command to achieve the same outcome in one step.

Run the R --version command to confirm that you are now running R version 3.5.3.

8. Unload all modules at once by executing:

module purge

9. Confirm you have no modules loaded.

When requesting a new version or software package, you may be asked to test a development version before it is released. These test installations will be made available through development modules.

After executing module load use.dev, the development modules appear in the top of the module avail output and can be loaded as usual using the module load command.

## **Tutorial 3 - Batch Job Submission**

Objective: To create and submit a batch job to the HPC cluster. Prerequisite: A text editor (e.g. vim).

Many Linux users decide to use Vi or ViM (Vi IMproved) as a text editor because of its efficiency to insert, delete, find and replace text only using the keyboard. On the HPC cluster, vim is readily available on the frontend and cluster nodes. For more information on how to use Vi, click here.

There are other editors available such as Nano, which is available via a module on the HPC cluster. To use Nano, run module load nano before starting this tutorial. For more information on how to use Nano, click here.

### Instructions:

- 1. Open a text editor to create a job script.
- 2. Add the following text and save the file as example.sh:

```
#!/bin/bash
#$ -cwd
#$ -j y
#$ -pe smp 1
#$ -l h_rt=1:0:0
#$ -l h_vmem=1G
#$ -m bea
hostname
date
sleep 2m
date
```

3. Submit the job script to the HPC cluster for processing by executing:

```
qsub example.sh
```

4. Make a note of the job number shown in the output after submitting the job. This is a unique identifier which can be used to check the job status and consumed





resources.

5. Check the status of your queued / running jobs by executing:

#### qstat

Note: Running qstat -j JOBID will provide a high level of detail about job JOBID - this should be replaced with the actual job number.

- 6. Once the job has completed running, confirm that it is no longer visible in the output of the qstat command. Examine the job output file, named example.sh.oJOBID, replacing JOBID with the actual job number.
- 7. View the contents of the job completion email to find helpful information such as exit status and resources used.
- 8. Confirm your job has completed by running qstat again you should not see the job in the output.
- 9. Modify the resources in the *example.sh* script, increasing the core request to 4 and the **total** memory to 8GB. Add a job name with the #\$ -N NAME parameter and repeat steps 3-5 and observe the change of output from qstat.

*Hint:* The h\_vmem setting is per core.

If your job is queued, you may see the position in the whole cluster queue with our showqueue utility script.

10. Check your list of queued jobs with qstat and cancel any jobs submitted in this section that you no longer require.

Note: you can cancel a queued / running job by running qdel JOBID.

## **Tutorial 4 - Interactive Job Submission**

Objective:	To submit interactive jobs to the HPC cluster.
Prerequisite:	An active connection to the HPC cluster.

Jobs that request up to 1 hour will become eligible for the short queue and will often start running immediately. All other jobs may request up to 10 days runtime with the  $h_{rt=240:0:0}$  parameter.

Jobs are killed automatically when the runtime has been exceeded. Therefore, we recommend setting the runtime to the maximum of 10 days when running jobs over 1 hour. For example, if your job will take 24 hours, you should still request 10 days to cover any unexpected overheads.

#### Instructions:

1. Submit an interactive job to the HPC cluster by executing:

### qlogin

Note: By default, all jobs request 1 core, 1GB RAM and 1 hour unless otherwise specified. For example, qlogin -pe smp 2 -1 h\_vmem=2G will request 2 cores and 2GB per core, while keeping the default 1 hour maximum runtime.

2. If there are available resources, the scheduler will provide an interactive session running on a compute node. Notice the change of prompt and the node the job is





executing on, which will no longer be a frontend node. Type qstat and find your running session in the list of your jobs, named QLOGIN.

3. Load an R module and run

Rscript /data/teaching/workshop/common/4/countdown.R

to execute a trivial R script from within the interactive session. Observe the output which should show decrementing numbers from 5 to 0.

- 4. Use exit to end the interactive job session.
- 5. Notice the change of prompt which will be a frontend node. Please do not start any computational work when connected to a frontend. All jobs must be submitted via qsub or qlogin.

# **Tutorial 5 - Using Resources Effectively**

Objective: To submit jobs which use HPC resources effectively. Prerequisite: A text editor (e.g. vim).

#### Instructions:

- 1. Open a text editor to create a job script.
- 2. Add the following text and save the file as bad\_job.sh:

 $\it Note$ : The trailing backslash character breaks a single line command over multiple lines, to aid readability.

- 3. Submit the job script to the HPC cluster for processing.
- 4. Check which cluster node the job is running on with qstat.
- 5. Run ssh -t NODE top -Hu \${USER} to log into the compute node running your job and observe the process table output to see that the process *blastn* is only using up to 100% CPU (1 core) and one thread, meaning the job is using fewer cores than requested, because we requested 2 cores. Press q to return to the prompt.

Note: Replace NODE with the actual node your job is running on.

6. Add -num\_threads 4 to the end of the bad\_job.sh script. The complete blastn command should look like:

blastn -db /data/teaching/workshop/common/5/files/nt \





```
-query /data/teaching/workshop/common/5/example_seqs.fa \
-out example_seqs_blastout.txt \
-num_threads 4
```

Note: Don't forget to add a backslash character to the -out parameter.

This will instruct blast to use 4 threads during execution, which will overload a compute node, resulting in a slower performance for your job and other user's also running on the same node. Please do not do this for production jobs.

- 7. Submit the job script to the HPC cluster and run the command to observe the process table output on the node running your job, to see that *blastn* is now running on 4 threads, even though we only requested 2 cores.
- 8. Replace -num threads 4 with -num threads \${NSLOTS} in the bad job.sh script.

We are now using the NSLOTS variable which is assigned at job runtime to the number of cores requested - this is the correct method when running multi-threaded applications on the HPC cluster. Feel free to rename the script now it is no longer bad

- 9. Submit the job script to the HPC cluster and run the command to observe the process table output on the node running your job, to see that *blastn* is now correctly running on 2 threads.
- 10. When the job has completed, examine the output file example\_seqs\_blastout.txt.
- 11. Confirm the exit status of your completed job by substituting JOBID with your job number in the following command:

```
qacct -j JOBID
Note: you can use grep to only view the exit status:
qacct -j JOBID | grep exit_status
```

## References

- SSH keygen tutorial https://blog.hpc.qmul.ac.uk/ssh-tutorial/
- Logging into the HPC https://docs.hpc.qmul.ac.uk/intro/login/
- Application Modules https://docs.hpc.qmul.ac.uk/using/UsingModules/
- Submitting Jobs https://docs.hpc.qmul.ac.uk/using/
- Job Script Builder https://docs.hpc.qmul.ac.uk/using/jobscriptbuilder/
- Interactive Jobs https://docs.hpc.qmul.ac.uk/using/#interactive-jobs
- Productivity tips for HPC users https://blog.hpc.qmul.ac.uk/productivity-tips-for-apocrita-cluster-users.html#content





## **Solutions**

## **Tutorial 3**

Modify the resources in the example.sh script, increasing the core request to 4 and the total memory to 8GB. Add a job name with the #\$ -N NAME parameter and repeat steps 3-5 and observe the change of output from qstat.

- #!/bin/bash
- #\$ -cwd
- #\$ -ј у
- #\$ -pe smp 4
- #\$ -1 h\_rt=1:0:0
- #\$ -1 h\_vmem=2G
- #\$ -m bea
- #\$ -N myjobname