# HELIOS

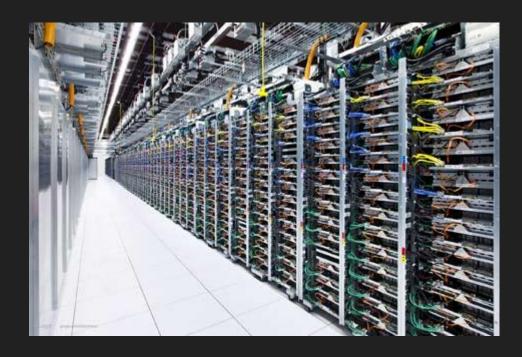
## About

- HELIOS is a computer cluster
  - computers are aggregated together to create a more powerful machine.

 User can access a specific machine (node) inside the cluster

or

 use multiple nodes at the same time for task that are computation heavy



## Connect to Helios

Should have received your logging info

If not ask for it to Mathieu Germain

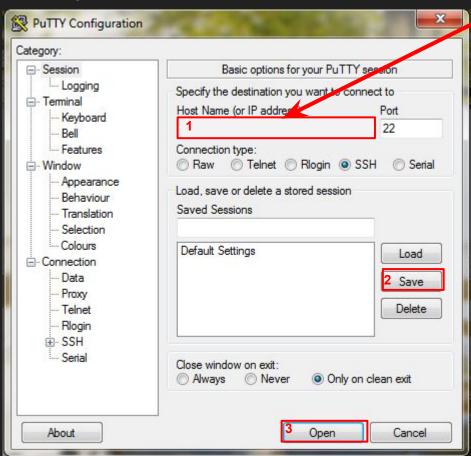
# Connect with linux

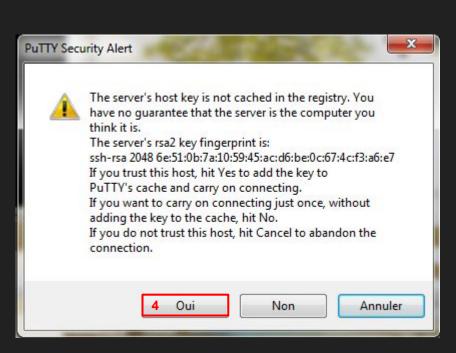
• ssh <account-name>@helios.calculquebec.ca

## Connect with windows

- Multiple solutions
  - Windows subsystem for linux
    - Makes ubuntu command line available on your windows machine
    - Can install all ubuntu packages
  - b. Install putty
    - This is a ssh client for windows
  - c. Use git bash
    - Git comes with ssh

# Putty





## HELIOS

 Once connected you should see something like this →

```
user1@mila-145D → _/helios / master • ssh user60@helios.calculquebec.ca
user60@helios.calculquebec.ca's password:
Last login: Wed Jan 9 10:52:58 2019 from 69.28.236.4
Vous êtes sur un noeud de login de Helios (Calcul Québec).
 - Nous n'effectuons pas de sauvegarde de vos fichiers.
 - N'utilisez pas le noeud de login pour executer votre code.
This is a Calcul Québec login node for Helios.
 - There is no backup of users files.
 - Do not use this node to run code.
Rapportez tout problème à / Report any problems to: helios@calculquebec.ca
Documentation: https://wiki.calculquebec.ca/
Suivre sur Twitter/Follow on Twitter: https://twitte
État des serveurs: http://serveurscq.computecanada.
Vous pouvez maintenant utiliser l'environnement logiciel de Calcul Canada
en utilisant la commande
source /admin/bin/enable cc cvmfs
ou
touch $HOME/.helios ccstac
(cette dernière sera petsis ante)
You can now use the soft are environment of Compute Canada by running
the command
source /amin/in/enable cc cvmfs
touch $HOME/.helios ccstack
(this last one will be persistent)
Due to MODULEPATH changes, the following have been reloaded:
 1) openmpi/2.1.1
[user60@helios1 ~]$
```

#### Bash - A few useful functions

mkdir: make directory

```
[delaunay@helios1 ~]$ mkdir my project
```

Is: list directories

```
[delaunay@helios1 ~]$ ls
my project script moab
```

• cd: change directory

```
[delaunay@helios1 ~]$ cd my_project/
[delaunay@helios1 my project]$
```

pwd: print working directory

```
[delaunay@helios1 my_project]$ pwd
/home/delaunay/my project
```

cat: show file content

```
[delaunay@helios1 ~]$ cat .bashrc
```

- vi: open a file with vi
  - to edit file press i
  - to quit press ESC then enter :q

```
[delaunay@helios1 ~]$ vi .bashrc
```

- Compute Resource are shared among a lot of people
- You need to demand access to those resources
- The resources will be allocated to you by the ...
  - o resource manager (torque) / job scheduler: moab
  - https://wiki.calculquebec.ca/w/Moab/en

> msub hello.pbs

- Submit a job
- hello.pbs is a bash script

```
#!/bin/bash
#PBS -A colosse-users
#PBS -l advres=MILA2019
#PBS -l feature=k80
#PBS -l nodes=1:gpus=1
#PBS -l walltime=XX:XX:XX
```

- Small jobs have priority!
- max 12h of runtime

```
[delaunay@helios1 MixedPrecisionTutorial]$ msub hello.pbs

371993

Job ID
```

```
[delaunay@helios1 MixedPrecisionTutorial]$ ls -all | grep 371993 
-rw----- 1 delaunay jvb-000-01 1406 Dec 13 14:19 371993.err 
-rw----- 1 delaunay jvb-000-01 64 Dec 13 14:19 371993.out
```

Job Output

#### Demo

```
# copy the example locally
git clone https://github.com/Delaunay/helios

# enter the example
cd helios

# Schedule the example to be run
msub hello.pbs
showq -w user=user60

# Show <job_id>.out
watch tail -n 20 $(ls -rt | grep .out | tail -n 1)
```

- Singularity
  - Sandboxed Execution
  - You have control over the container (image)
  - Pre configured container to simplify your life

```
singularity exec --nv --bind source:dest container script.sh
    exec : execute a script
    --nv : mount NVIDIA GPUs
    --bind : make host folder (source) available inside the container (dest)
    container: image of the container you want to use
    script.sh: the script you want to run
```

```
#!/bin/bash
#PBS -A colosse-users
#PBS -1 advres=MILA2019
#PBS -1 feature=k80
#PBS -l nodes=1:gpus=1
#PBS -l walltime=XX:XX:XX
PATH=$PATH:/opt/software/singularity-3.0/bin/
# set the working directory to where the job is launched
cd "${PBS O WORKDIR}"
# Singularity options
FOLDERS=$RAP,$HOME
SINGULARITY EXEC="singularity exec --nv --bind $FOLDERS $IMAGE"
# start your python script
```

```
# Schedule the example to be run
msub singularity.pbs

# Show <job_id>.out
watch tail -n 20 $(ls -rt | grep .out | tail -n 1)
```

# How to run Al stuff interactively

```
> msub -N debug -A colosse-users -l advres=MILA2019, feature=k80 nodes=1:gpus=1, walltime=15:00 -I -qtest
> module --force purge
> PATH=$PATH:/opt/software/singularity-3.0/bin/
> singularity shell --nv --bind $RAP,$HOME /rap/jvb-000-aa/COURS2019/etudiants/ift6759.simg
```

Useful for testing & debugging

• The walltime/user time allocated to them is small

## **Monitor Jobs**

- Email notification from moab
- showq -w user=<username>
  - Show the current jobs running for you
- checkjob <jobid>
  - Show details on a particular job (resource usage, status)
- mjobctl -c <jobid>
  - Cancel job

# Quality of life

Modify your ~/.bashrc to pre configure your environment

```
vi ~/.bashrc
```

Add at the end of the file:

```
source /rap/jvb-000-aa/COURS2019/etudiants/common.env
```

This will configure your environment and provide a few shortcuts

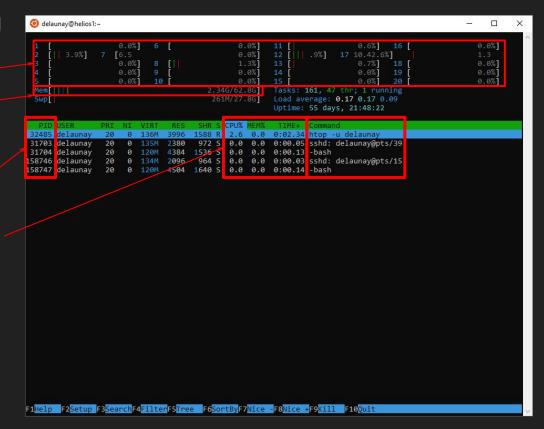
# Quality of Life

- Shortcuts:
  - mdebug
    - Start an interactive session
  - o s shell
    - Start a singularity shell
  - o s exec
    - Execute a command inside singularity
  - o show\_out
    - monitor latest output log
  - show\_err
    - monitor latest error log

## Monitor resource usage

- RAM and CPU usage monitoring
  - htop -u <username>
    - w usage per cores
    - Memory usage

- PID: Program Identifier
- Resource usage per program
- To kill a running program:
  - o kill -9 <PID>



# Monitor resource usage

GPU usage monitoring

o nvidia-smi

Software versions

[delaunay@gpu-k20-03 ~]\$ nvidia-smi Thu Dec 13 12:35:39 2018 Driver Version: 410.73 NVIDIA-SMI 410.73 CUDA Version: 10.0 Volatile Uncorr. ECC Persistence-M Bus-Id Disp.A Name Perf Pwr:Usage/Cap Memory-Usage GPU-Util Compute M. Tesla K20m 00000000:05:00.0 Off E. Process GPU Memory Usage GPU Memory Processes: PID Usage GPU Type Process name Program using GPU No running processes found

**GPU Compute Usage** 

# Monitor resource usage

• nvidia-smi --loop=1 --query-qpu=utilization.gpu,utilization.memory,memory.used,memory.total --id=0 --format=csv

```
[delaunay@gpu-k20-03 ~]$ nvidia-smi --loop=1 --query-gpu=utilization.gpu,utilization.memory,memory.used,memory.total --id=0 --format=csv utilization.gpu [%], utilization.memory [%], memory.used [MiB], memory.total [MiB]
0 %, 0 %, 0 MiB, 4743 MiB
```

- --loop=1 run nvidia-smi every second
- --query-gpu specify which statistic to print
- --id=0 show statistic only for the first gpu
- --format=csv print each iteration as a new CSV line

#### More documentation

- How to use Compute Canada Clusters
  - https://github.com/SMART-Lab/smartdispatch/wiki/How-To-Use-Compute-Canada-Clusters
  - https://docs.computecanada.ca/wiki/Python
- nvidia-smi
  - http://developer.download.nvidia.com/compute/DCGM/docs/nvidia-smi-367.38.pdf
- Singularity
  - https://www.sylabs.io/guides/3.0/user-guide/index.html