

# HPCs at OU

Fabia U. Battistuzzi

battistu@oakland.edu

Tomas Hajek hajek@oakland.edu

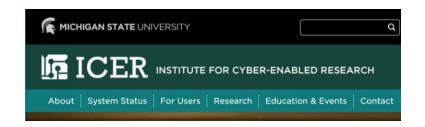
Stephen Villagonzalo <a href="mailto:svillagonzalo@oakland.edu">svillagonzalo@oakland.edu</a>

John Johnston <a href="mailto:jbjohnston@oakland.edu">jbjohnston@oakland.edu</a>



# Oakland University... StartPage RecentChanges FindPage HelpContents HPCMatilda

MSU ICER icer.msu.edu





# For general information:

https://kb.oakland.edu/uts/ResearchComputingH PC#Introduction to HPC clusters

### Oakland University... FindPage ResearchComputingHPC

Getting Started FAQ

### Public Documents

**OU Community** Banner Unlock/Reset Help

OU Forms and Tech Depot

Argos Reporting Banner Document Management Degree Works FAQ

Degree Works GoAnywhere Google Groups

MDUU Oracle Java SE Changes

### Common Issues

NetID Information Peer-to-Peer File Sharing

University Technology Services (UTS)

### RESEARCH COMPUTING AND HPC

### Contents

- · Research Computing and HPC
  - · Introduction to HPC clusters
  - Using the Cluster
  - Data Management
  - · Application Specific Help
  - System Status
  - MSU iCER HPCC Buy-in

### INTRODUCTION TO HPC CLUSTERS

- · What is an HPC Cluster
- · HPC cluster at Oakland

### **USING THE CLUSTER**

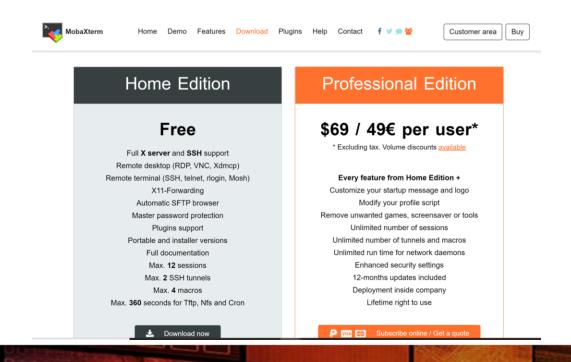
- Requesting a Matilda Account
- · Logging in to the Matilda HPC Cluster
- Logging into Matilda with MobaXterm
- Logging into Matilda with PuTTY
- Cluster Software Modules System
- Available HPC Software
- Launching and Controlling Jobs
- SBATCH Options
- Job Script Examples

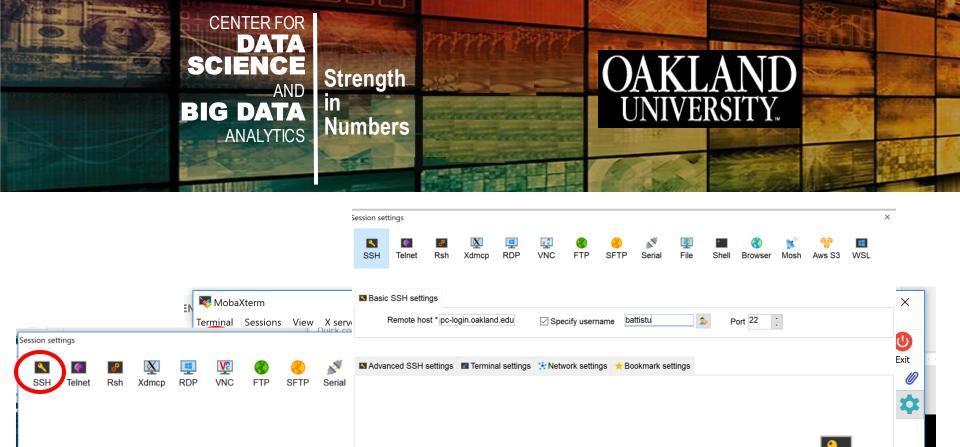


To access the cluster from Windows:

MobaXterm download (home edition)

https://mobaxterm.mobatek.net/download.html





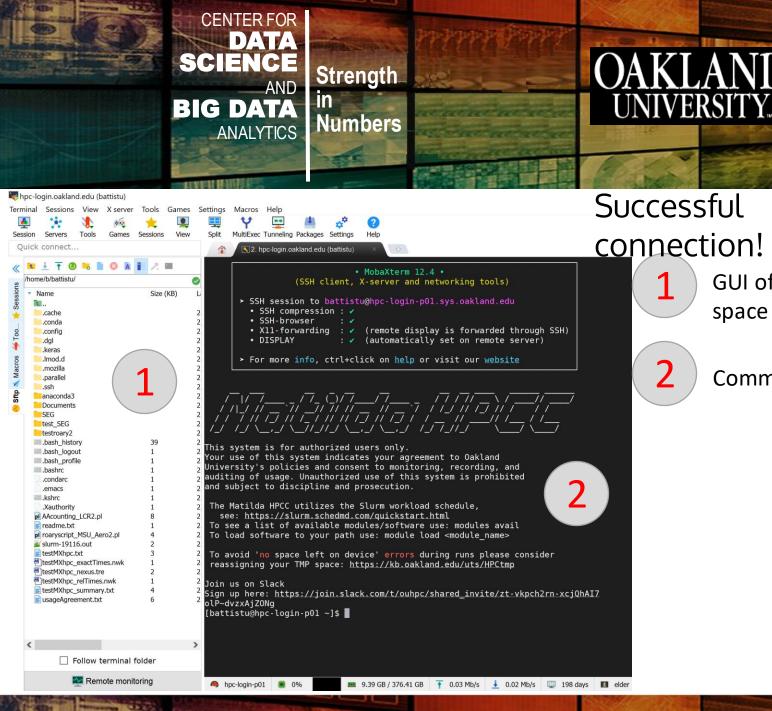
Remote host: hpc-login.oakland.edu



Username: netid





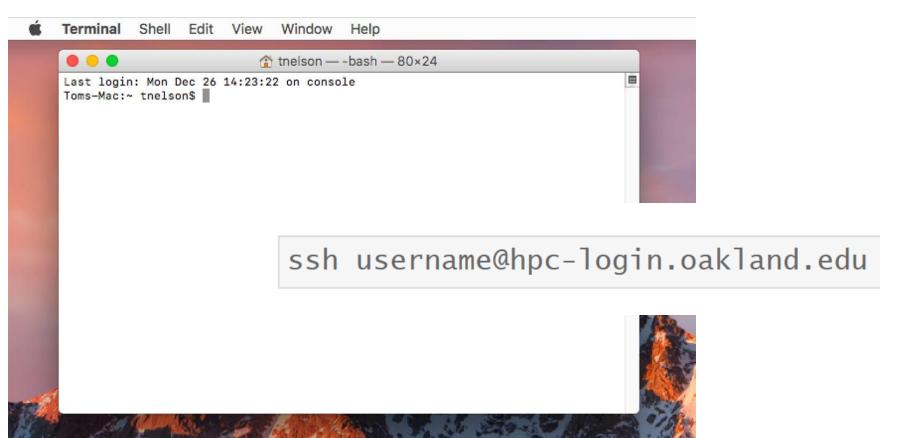


Successful connection!

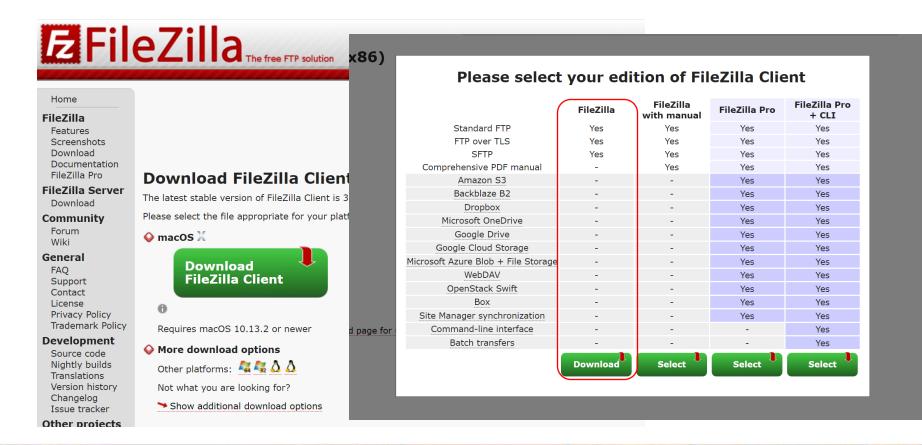
- GUI of your home space on the cluster
  - Command line space



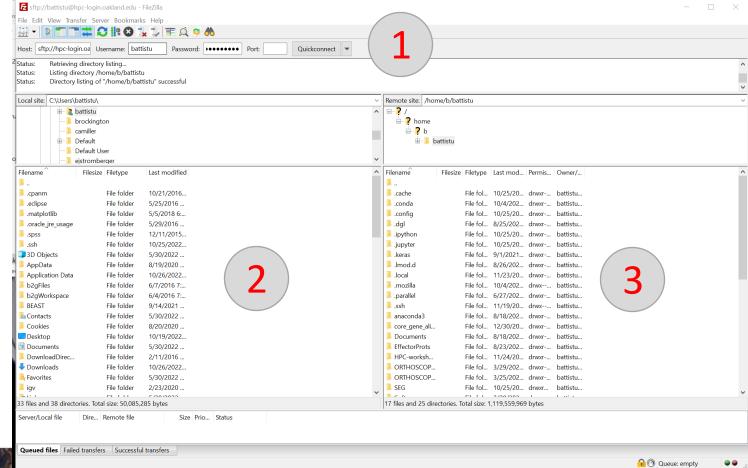
### To access the cluster on Mac:



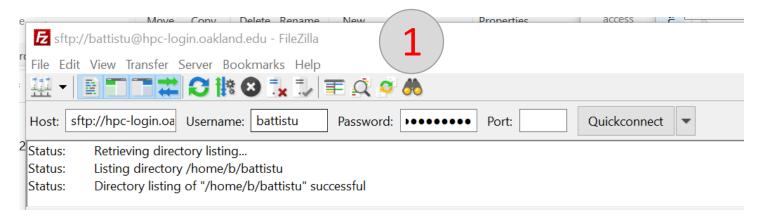












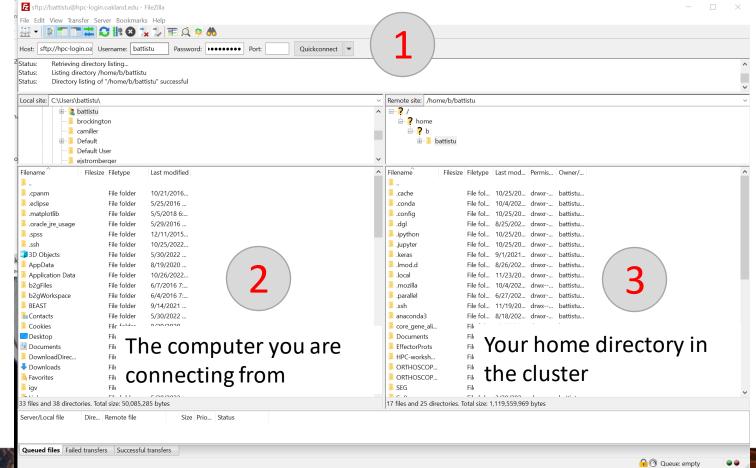
Host: hpc-login.oakland.edu

Username: netID

Password: netID password

Port: 22

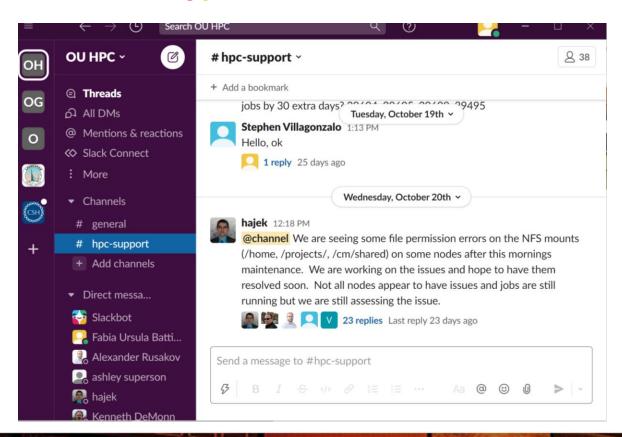


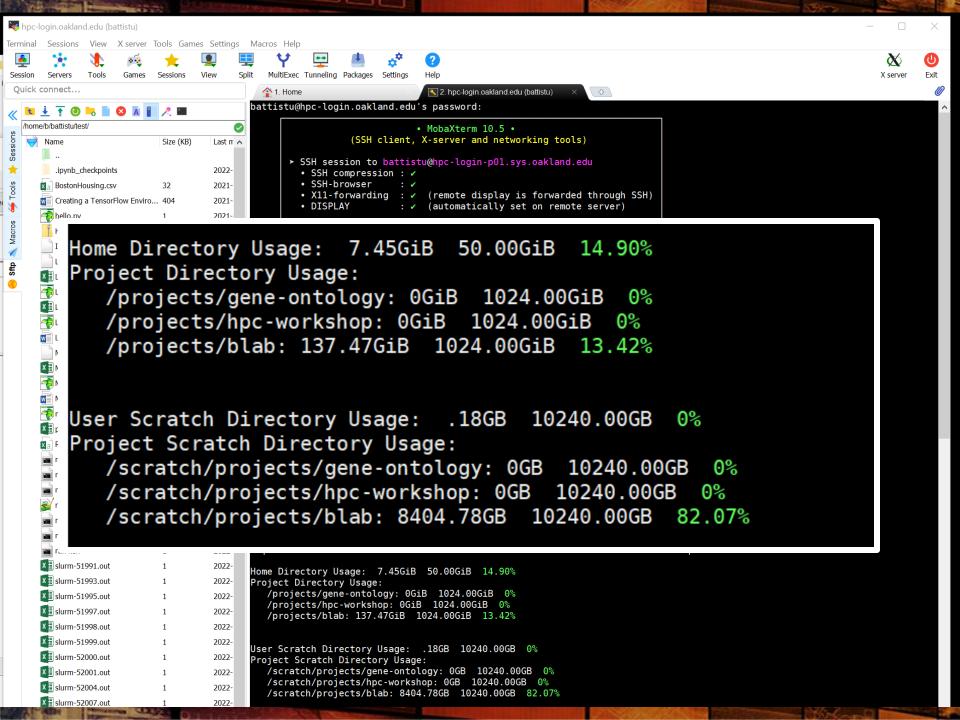




# Technical support:









To get a Matilda account: forms.oakland.edu

If you had an account created just for this workshop, it will be deactivated Dec. 6. To request a long-term account follow the instructions

M

**MarketPlace Access Request** 

MarketPlace Account Unlock/Password Reset Request

**MarketPlace Product Request** 

Matilda HPC Cluster Access Request

Miscellaneous Pay - Staff Employees Only

**Modify an Approved Undergraduate Program** 

MSDNAA Student Account Request (No Login)

My PUB Alumni Profile and Impact Statement

N

**NetID Guest Account Request** 

**NetID Shared Account Request** 

0

Oakland University On-Campus Accident Form (No Logic



To access Matilda from outside of campus: forms.oakland.edu

Request a VPN account, set up DUO authentication with your phone (or another device)



Software & Hosted Solution Purchasing Checklist (No Login)

Special Account Request (pdf)

**Special Credit Offering Request** 

Student Employee/Intern Confidentiality Agreement

Т

**Technology Control Plan Template** 

Touchnet Unlock Account / Password Reset

**Travel Authorization Request (UTS)** 

**Tuition Assistance Application** 

**Tuition Assistance Application for Faculty** 

۷

Vendor/Consultant Account Access Request

Virtual Labs Change Request

**VPN Access Request** 

Υ

Youth protection Approval Request

Z

**Zoom HIPAA Request** 



To execute an analysis:

Use SLURM to schedule your job in the cluster. This will make the job go in a queue and allow to request resources (# nodes, # of CPUs, types of nodes, etc)

For testing purpose, you can skip the cue and run the job on the login node (only for testing purposes!)



Let's run "Hello world"





### To run you will need:

- 1. The script
- 2. Python loaded on the cluster

### To write the script on windows:

- 1. Open Notepad++ on Windows (do not use Word) or BioEdit on Mac on your computer (not the cluster)
- 2. Type: print("Hello world")
- 3. Save as "helloworld.py"
- 4. Upload (or drag and drop) the file into your home directory in the cluster

### To write the script in Linux:

- 1. In your home directory type: nano (this is a text editor in linux)
- Type: print("Hello world")
- 3. Ctrl+x, give a file name and follow instructions to save it

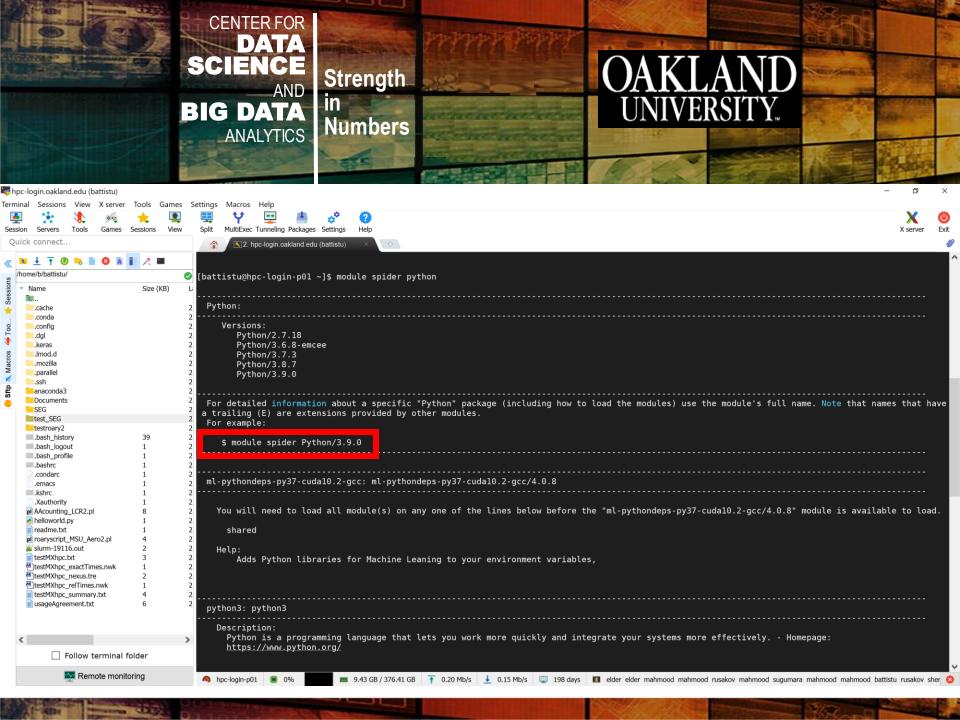


# To run you will need:

- 1. The script
- 2. Python loaded on the cluster (python is automatically loaded when you log in but for practice let's go through the process)

# In the cluster type:

module spider python





# To run you will need:

- 1. The script
- 2. Python loaded on the cluster



# To run you will need:

- 1. The script
- 2. Python loaded on the cluster

```
battistu@hpc-login-p01 ~] module list

urrently Loaded Modules:

1) shared 2) DefaultModules 3) dot 4) slurm/slurm/19.05.8 5) default-environment 6) gcc/9.2.0

battistu@hpc-login-p01 ~]$
```



# To run you will need:

- 1. The script
- 2. Python loaded on the cluster

```
battistu@hpc-login-p01 ~] module load Python/3.9.0
battistu@hpc-login-p01 ~] module load Python/3.9.0
battistu@hpc-login-p01 ~] module list

Currently Loaded Modules:

1) shared 2) DefaultModules 3) dot 4) slurm/slurm/19.05.8 5) default-environment 6) gcc/9.2.0 7) Python/3.9.0

battistu@hpc-login-p01 ~]$
```



### To run you will need:

- 1. The script
- 2. Python loaded on the cluster

# To run your script type:

python3 helloworld.py

Case sensitive
Do not use spaces
Meaningful names



[battistu@hpc-login-p01 ~]\$ python3 helloworld.py Hello world [battistu@hpc-login-p01 ~]\$ ■



To submit your job you will need:

- 1. The script
- 2. The SLURM file
- 3. Python loaded on the cluster



#!/bin/bash --login



python3 helloworld.py > output.txt ### call your executable

scontrol show job \$SLURM\_JOB\_ID ### write job info to output file



```
1 ₽#!/bin/bash --login
   ######## SBATCH Lines for Resource Request #########
   #SBATCH --time=00:01:00
                                       # limit of wall clock time - how long the job will run in hrs:min:sec (same
   as -t). Max 168:00:00 (7 days)
   #SBATCH --nodes=1
                                     # number of different nodes - could be an exact number or a range of nodes
    (same as -N)
   #SBATCH --ntasks=1
                                       # number of tasks - how many tasks (nodes) that you require (same as -n)
   #SBATCH --cpus-per-task=40
                                        # number of CPUs (or cores) per task (same as -c). Max for most nodes: 40
                                       # memory required per allocated CPU (or core) - amount of memory (in bytes)
   #SBATCH --mem-per-cpu=2G
   #SBATCH --job-name HelloWorld # you can give your job a name for easier identification (same as -J)
10
11
12
   ######### Command Lines to Run #########
13
   module load Python/3.9.0
14
15
   cd /home/b/battistu
                                ### change to the directory where your code is located
16
17
   python3 helloworld.py > outut.txt
                                            ### call your executable
18
   scontrol show job $SLURM JOB ID ### write job information to output file
19
20
```



Upload the SLURM file into your home directory Run the job:

sbatch MatildaSLURM\_example.sb

```
battistu@hpc-login-p01 ~]$
battistu@hpc-login-p01 ~]$ sbatch MatildaSLURM_example.sb
batch: error: Batch script contains DOS line breaks (\r\n)
batch: error: instead of expected UNIX line breaks (\n).
battistu@hpc-login-p01 ~]$
```



module spider dos2unix
module load shared
module load dos2unix/7.4.2
dos2unix MatildaSLURM example.sb

Repeat sbatch command



```
battistu@npc-login-p01 ~]$
battistu@hpc-login-p01 ~]$
battistu@hpc-login-p01 ~]$
battistu@hpc-login-p01 ~]$ sbatch MatildaSLURM_example.sb
bubmitted batch job 31287
battistu@hpc-login-p01 ~]$ squeue -u battistu

JOBID PARTITION NAME USER ST TIME NODES NODELIST(REASON)
battistu@hpc-login-p01 ~]$
```

- pl AAcounting LCR2.pl
- helloworld.py
- MatildaSLURM\_example.sb
- MatildaSLURM\_example2l.sb
- MatildaSLURM\_examplemod.sb
- output.txt
- readme.txt
- pl roaryscript\_MSU\_Aero2.pl
- slurm-31284.out
- 🔊 slurm-31285.out
- slurm-31286.out
- slurm-31287.out

squeue -u netid to check job status scancel job# to cancel a job

Output is created as expected

Slurm-job#.out: check how the job ran; this is where you will find error messages





You just completed your first HPC analysis