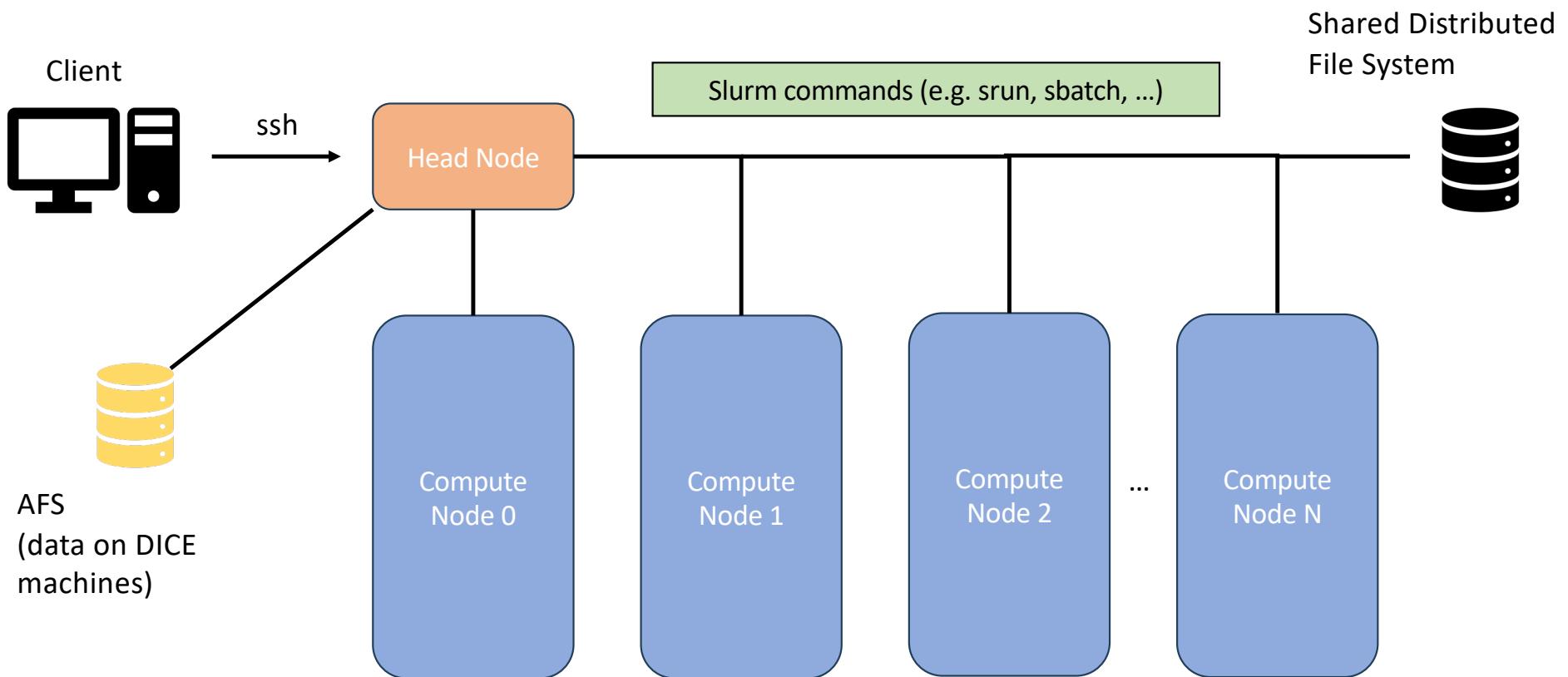


Teaching GPU Cluster

Teaching Cluster Overview



Teaching GPU Cluster

Basic Info:

The teaching cluster has over 120 GPUs in 20 servers (landonia[01-25]) with:

- 100 NVIDIA GTX 1060 6GB GPUs,
- 8 NVIDIA RTX A6000 48GB GPUs,
- a small number of NVIDIA TITAN-X 12GB GPUs.

•

Usage:

Access through **ssh** and use **SLURM** commands.

1. Access the cluster's head node via:

- Connect to the Informatics VPN or use the DICE machine and then run: **ssh <YOUR_UUN>@mlp.inf.ed.ac.uk**
- You can remotely connect to your DICE machine using the command **below** if you are connected to the Informatics VPN or the eduroam Wi-Fi. Once you have accessed the DICE machine, run the command mentioned **above**.

ssh <YOUR_UUN>@ssh.inf.ed.ac.uk

- You can find information about how to connect to Informatics VPN at:
<https://computing.help.inf.ed.ac.uk/openvpn>

Teaching GPU Cluster

Usage (continue):

After successfully connecting to the head node, your terminal will display the following:

```
Welcome to Ubuntu 20.04.6 LTS (GNU/Linux 5.15.0-130-generic x86_64)

This is a cluster head node please do not run compute intensive processes here
this node is intended to provide an interface to the cluster only
Please nice any long running processes
-----
Looking for tips? https://authcomputinghelp.inf.ed.ac.uk/cluster-tips
Jan 2025: Note that some GRES have changed to swap dashes to underscores.

Last login: Mon Jan 13 15:45:11 2025 from openvpn-125-012.inf.ed.ac.uk
[uhtred]s2020153: █
```

2. Run your code with SLURM command:

There are two ways you can run your code:

- Interactive job - allow you to interact directly with the allocated compute resources, good for debugging and testing code.
- Batch job - run the task automatically based on a pre-written script (.sh file) without interaction with the compute node and constant user attention, good for long-running tasks.

Teaching GPU Cluster

Usage (continue):

E.g. Let's say you want to run the code utilizing 1 GPU.

Interactive jobs:

```
srun -p Teaching -w saxa --gres gpu:1 --pty bash
```

After executing this command, run “nvidia-smi”, and you will see output similar to the following:

```
Thu Jan 22 04:24:57 2026
+-----+-----+-----+-----+-----+-----+-----+-----+
| NVIDIA-SMI 570.195.03 | Driver Version: 570.195.03 | CUDA Version: 12.8 |
| GPU  Name  Persistence-M  Bus-Id  Disp.A  Volatile Uncorr. ECC |
| Fan  Temp  Perf  Pwr:Usage/Cap | Memory-Usage | GPU-Util  Compute M. |
| 0  NVIDIA H200  00000000:19:00.0 Off  60MiB / 143771MiB | N/A  Default |
| 1  NVIDIA H200  00000000:3B:00.0 Off  60MiB / 143771MiB | N/A  Default |
| 2  NVIDIA H200  00000000:4C:00.0 Off  60MiB / 143771MiB | N/A  Default |
| 3  NVIDIA H200  00000000:5D:00.0 Off  60MiB / 143771MiB | N/A  Default |
| 4  NVIDIA H200  00000000:9B:00.0 Off  60MiB / 143771MiB | N/A  Default |
| 5  NVIDIA H200  00000000:BB:00.0 Off  60MiB / 143771MiB | N/A  Default |
| 6  NVIDIA H200  00000000:CB:00.0 Off  60MiB / 143771MiB | N/A  Default |
| 7  NVIDIA H200  00000000:DB:00.0 Off  60MiB / 143771MiB | N/A  Default |
+-----+-----+-----+-----+-----+-----+-----+-----+
| MIG devices: |
| GPU  GI  CI  MIG  Memory-Usage  Vol  Shared |
| ID  ID  ID  Dev  BAR1-Usage  SM  Unc  CE  ENC  DEC  OFA  JPG |
| 0  7  0  0  9MiB / 16384MiB  16  0  1  0  1  0  1 |
| 0  0  0  0  0MiB / 32767MiB  0  0  0  0  0  0  0 |
+-----+-----+-----+-----+-----+-----+-----+-----+
| Processes: |
| GPU  GI  CI  PID  Type  Process name  GPU Memory |
| ID  ID          Usage |
+-----+-----+-----+-----+-----+-----+-----+
| No running processes found |
+-----+-----+-----+-----+-----+-----+-----+
```

Teaching GPU Cluster

Usage (continue):

E.g. Let's say you want to run the code utilizing 1 GPU.

Batch jobs: `sbatch --gres=gpu:1 test.sh`

After executing this command, you will see “Submitted batch job [ID]”.

The output file will be saved as `slurm-[ID].out` by default in the directory where you call sbatch. You can set the output directory and file names by using the `--output` option with the sbatch command.

test.sh:

You must include this

```
#!/bin/bash
echo I love Machine Learning Systems.
pwd
```

```
[uhtred]s2020153: sbatch --gres gpu:1 test.sh
Submitted batch job 1945143
[uhtred]s2020153: cat slurm-1945143.out
I love Machine Learning Systems.
/home/s2020153
```

Teaching GPU Cluster

Usage (continue):

Since servers are equipped with different types of GPUs, you can specify the type of GPU you wish to use:

E.g. You want to have 1 NVIDIA Titan X GPU:

```
srun --gres=gpu:titan_x:1 --pty bash
sbatch --gres=gpu:titan_x:1 test.sh
```

```
[uhtred]s2020153: srun --gres=gpu:titan_x:1 --pty bash
srun: job 1944884 queued and waiting for resources
srun: job 1944884 has been allocated resources
[landonia08]s2020153: nvidia-smi
Mon Jan 13 15:46:28 2025
+-----+
| NVIDIA-SMI 550.127.08      Driver Version: 550.127.08    CUDA Version: 12.4 |
|                               +-----+                         +-----+
| GPU  Name                  Persistence-M | Bus-Id      Disp.A  | Volatile Uncorr. ECC | | |
| Fan  Temp     Perf          Pwr:Usage/Cap |          Memory-Usage | GPU-Util  Compute M. |
|                               |             |               |           |          MIG M. |
+-----+-----+-----+-----+-----+-----+-----+-----+
| 0  NVIDIA GeForce GTX TITAN X     Off | 00000000:07:00.0 Off |           N/A |
| 22%  27C     P8          16W / 250W |          0MiB / 12288MiB |     0%    Default |
|                               |                           |           N/A |
+-----+-----+-----+-----+-----+-----+-----+-----+
```

You are allowed to request a maximum of 8 GTX 1060/4 Titan X/1 Titan X Pascal GPU/2 A6000 GPUs at a time.

This year we have 8 NVIDIA H200 for you! (so I guess nobody want to use 1060 any more)

You can get access with `-p Teaching -w saxa` after srun/sbatch command.

NOTE: Please allocate resources according to your specific requirements. Avoid over-allocating to ensure resources are available for others!

Teaching GPU Cluster

Usage (continue):

Other useful SLURM commands:

Check all available GPU types: `scontrol show node | grep gpu`

```
Gres=gpu:titan_x_pascal:1(S:0),gpu:titan_x:2(S:0),gpu:gtx_1060:2(S:1)
CfgTRES=cpu=12,mem=96000M,billing=12,gres/gpu=5
Gres=gpu:gtx_1060:8(S:0-1)
CfgTRES=cpu=12,mem=96000M,billing=12,gres/gpu=8
```

Check current SLURM job status: `squeue`

```
[[uhtred]s2020153: squeue
      JOBID PARTITION      NAME      USER ST      TIME  NODES NODELIST(REASON)
      1563085 General_U collect_ s2263903 PD      0:00      1 (BadConstraints)
      1944830 PGR-Stand ilcc_pre s2148449 PD      0:00      1 (Resources)
```

Cancel a job: `scancel <job_id>`

For more details about the school cluster and running SLURM commands, please refer to:
<https://computing.help.inf.ed.ac.uk/teaching-cluster>.

IMPORTANT: Please read the **Files and Backups** section and **GPU cluster tips** in this link carefully, as it contains crucial information about file storage and can help prevent data loss.

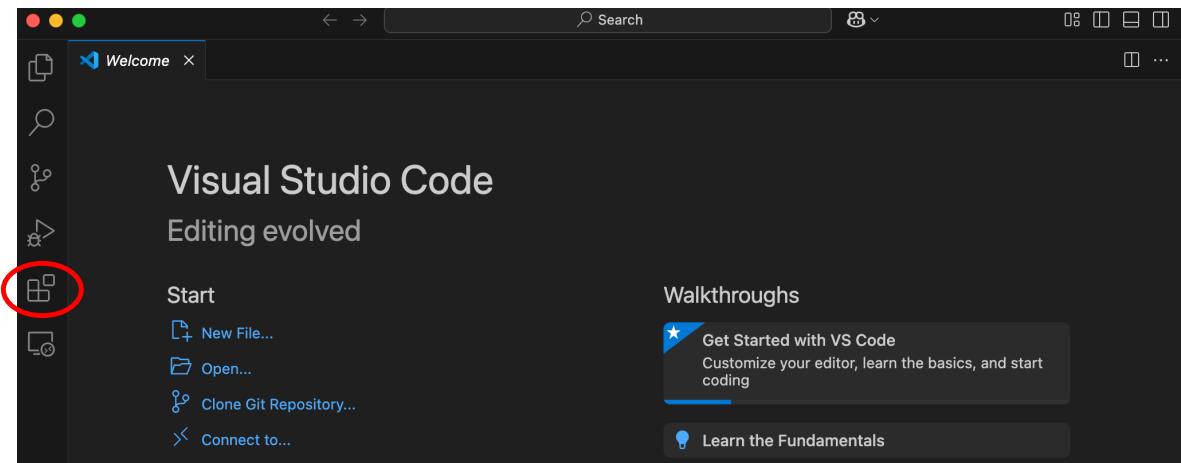
Teaching GPU Cluster

How to write codes in the teaching cluster?

VSCode!

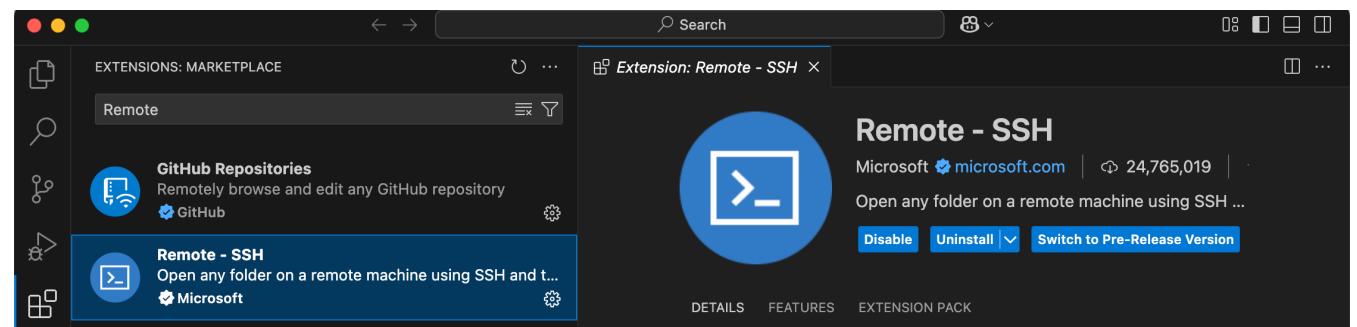
1. Open your VScode

2. Go to Extensions



3. Search for the keyword

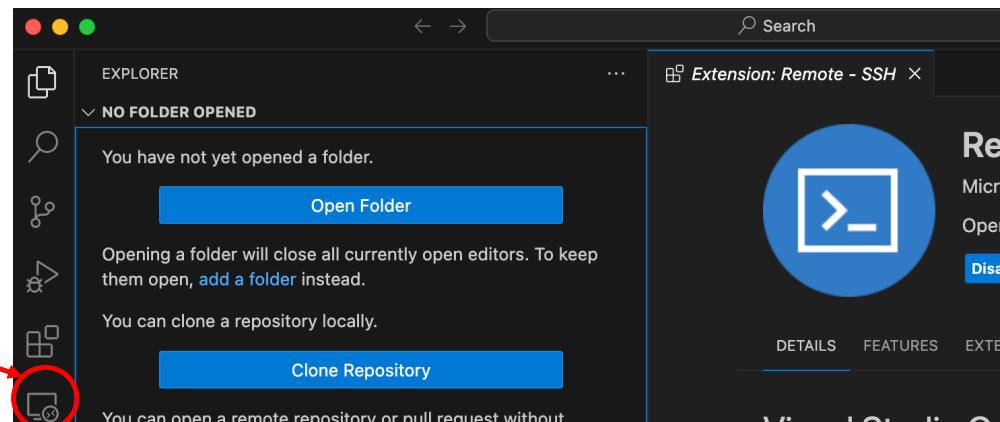
"Remote" and install the "Remote - SSH" extension.



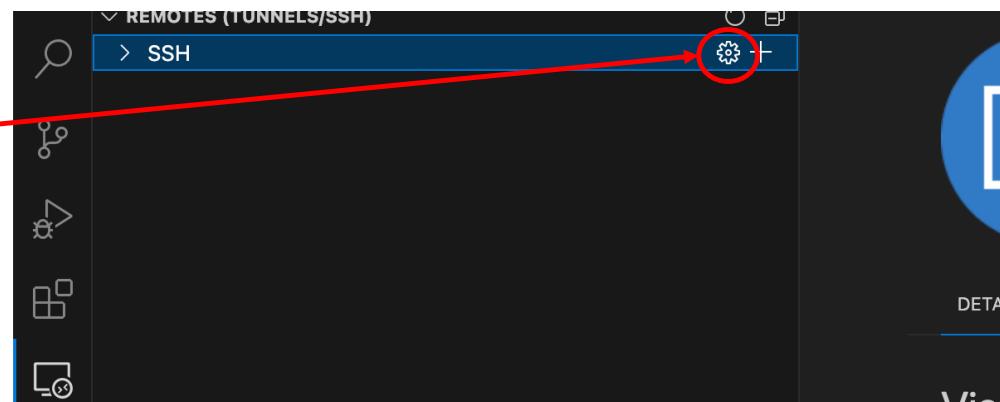
Teaching GPU Cluster

How to write codes in the teaching cluster? (Continue)

4. Once the extension is installed, the REMOTES icon will appear on the left sidebar.



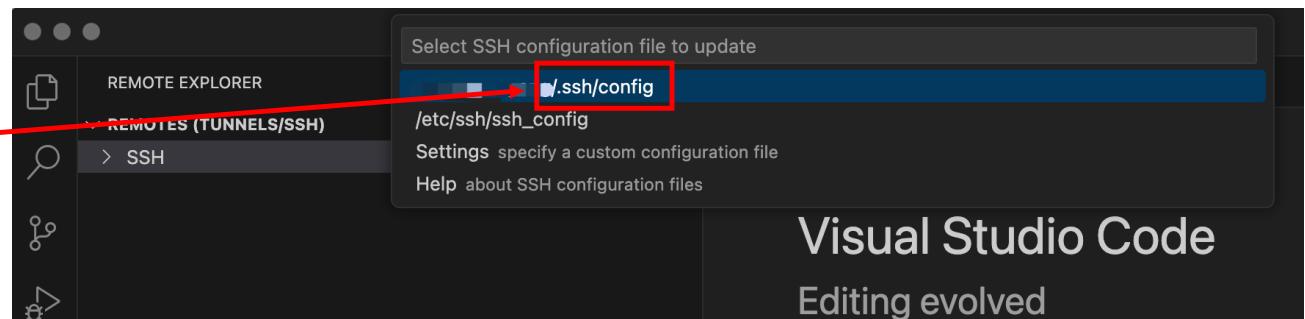
5. Click on the REMOTES icon, then select the Config icon located to the right of "SSH."



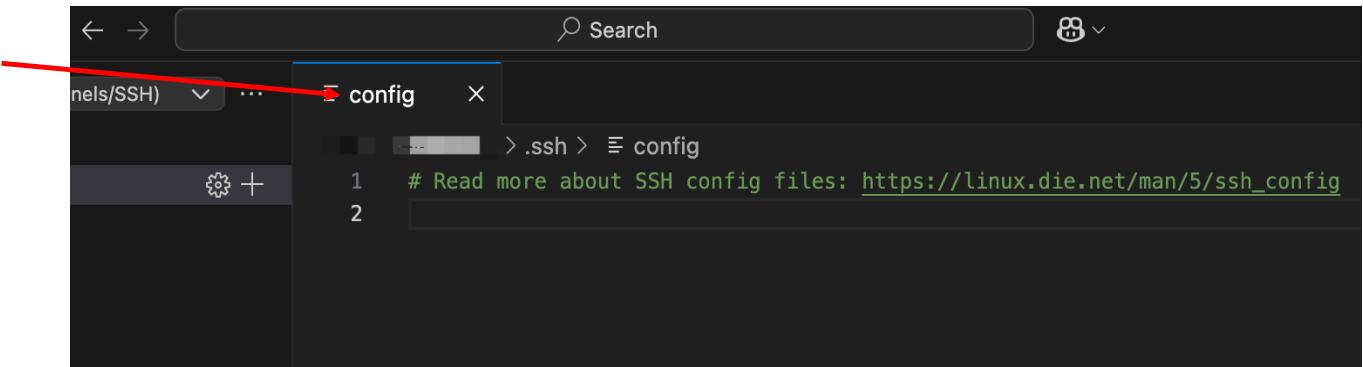
Teaching GPU Cluster

How to write codes in the teaching cluster? (Continue)

6. After clicking the Config icon, choose the configuration file with a path ending in .ssh/config.



7. You will see a page like this.



Teaching GPU Cluster

How to write codes in the teaching cluster? (Continue)

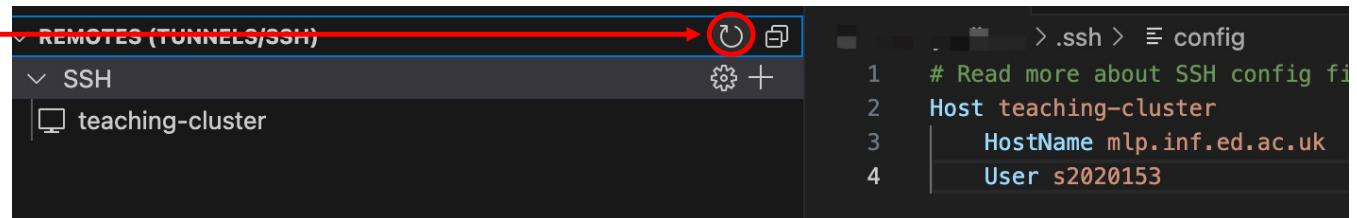
8. In this page, write the SSH configuration text and **save it**:

```
Host teaching-cluster
```

```
HostName mlp.inf.ed.ac.uk
```

```
User <YOUR_UUN>
```

9. Refresh the **REMOTES** page, and a screen similar to this will appear.



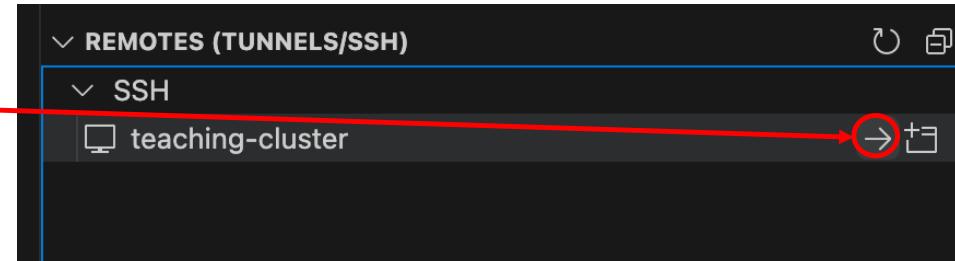
```
REMOTES (TUNNELS/SSH)
SSH
teaching-cluster

> .ssh > config
1 # Read more about SSH config fi
2 Host teaching-cluster
3   HostName mlp.inf.ed.ac.uk
4   User s2020153
```

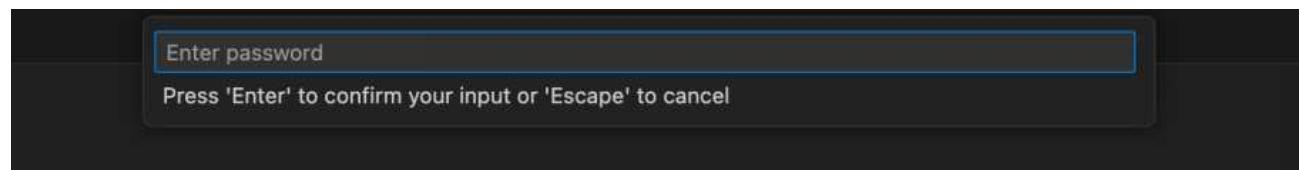
Teaching GPU Cluster

How to write codes in the teaching cluster? (Continue)

10. Connect to the Informatics VPN or Use the DICE machine, then click the right arrow icon to connect to the teaching cluster's head node:



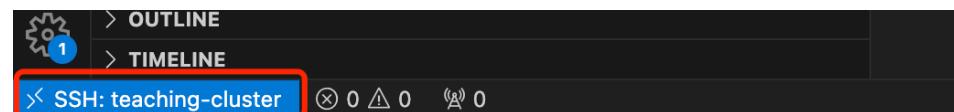
11. Enter the password of your DICE account



Teaching GPU Cluster

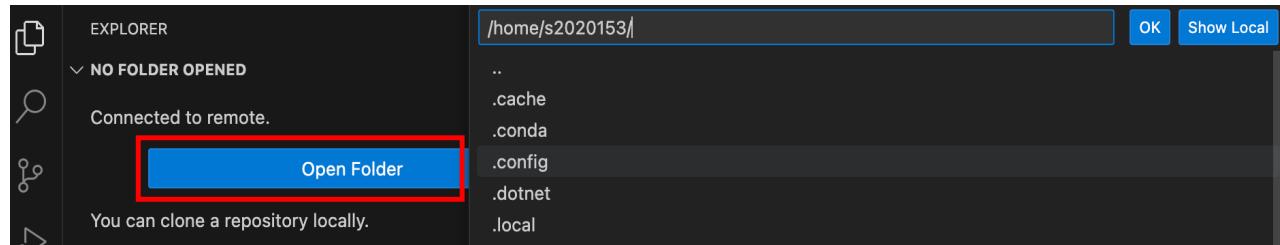
How to write codes in the teaching cluster? (Continue)

12. If this screen appears, congratulations! You have successfully connected to the teaching cluster's head node.



13. Go to your home directory by clicking "Open Folder", and select your home directory (`/home/<YOUR_UUN>`). Click OK.

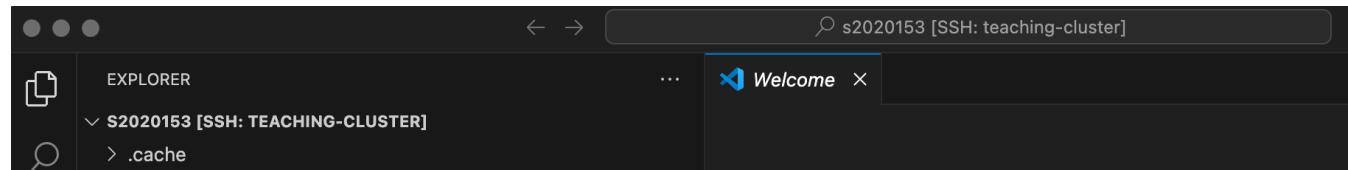
If password required, just type your password of the DICE account again.



Teaching GPU Cluster

How to write codes in the teaching cluster? (Continue)

14. If you can see this screen, you are now ready to write your code in the teaching cluster.



IMPORTANT NOTES:

1. Any changes you make in VSCode, such as writing code, creating or deleting files and folders, will automatically synchronize with the cluster. That's why we recommend it 😊. **Do not forget to save your changes**
2. Only install your environment and write your code on the head node. Actions such as Git operations and similar tasks should also be performed on the head node, as they will not work on the compute nodes.
3. Only run your code on the compute nodes using `srun` or `sbatch`, as the head node has limited computing resources and does not have GPUs. Running `torch.cuda.is_available()` on the head node will return `False`.

If you have further issues about accessing the teaching cluster, please contact the Computing Support Team:
<https://computing.help.inf.ed.ac.uk/>