DEPARTMENT OF COMPUTING

IMPERIAL COLLEGE OF SCIENCE, TECHNOLOGY AND MEDICINE

Introduction Lecture and GPU sources

Summary of the GPU resources available to Imperial Students taking Deep Learning module

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1 CSG Guides

CSG source

2 Basic SSH into the labmachine

2.1 Setting up the public key to not type in password

Copy the public key on your machine over to the authorized hosts on the other machine to avoid typing in your password every time you ssh. [source]

```
ssh-copy-id -i ~/.ssh/id_ed25519.pub <usrname>@<host>
```

2.2 Setting up the tunnel from shell1 – another machine

Create an alias on your home machine for the shell you'll use [source1] [source2]

```
Match host shell*
Hostname %h.doc.ic.ac.uk
User az620
```

To begin tunnelling through with ssh [source] and find your favourite [workstation] and set up tunnel with [source3]

```
Match host texel*
Hostname %h.doc.ic.ac.uk
User az620
ProxyJump shell1
```

2.3 Checking if resource being used

```
nvidia-smi && top htop
```

3 SLURM [reference]

The department has a pool of GPUs for deep learning tasks. To reduce complexity of scheduling, slurm is a system that schedules tasks into these resources.

3.1 SSH into the gpu cluster

In the ~/.ssh/config file, add the line below and ssh into it to schedule jobs. This should only be used for scheduling jobs into the slurm scheduler.

```
# should be either 2 or 3

Match host gpucluster*

Hostname %h.doc.ic.ac.uk

User az620

ProxyJump shell1
```

Then submit a pre-existing script using the sbatch command. The output will be stored, by default, in the root of your ~/ directory, with the filename slurm20-{xyz}.out.

See Section 6 for guide on python virtual environment.

3.2 Using CUDA

Many jobs will make use of the Nvidia CUDA toolkit, multiple versions of which are at /vol/cuda. If there ever appears a need to use this toolkit then in the submission bash script add:

```
./vol/cuda/12.0.0/setup.sh # if you're using version 12.0.0
```

Care that you choose a version with which your tensorflow or pytorch are compatible with.

3.3 Template Example bash submission script

This example assumes you have followed the previous steps and installed a python environment (using virtualeny, extra lines may be needed using minconda, check the example script further below) as directed. Please adjust paths if you have an existing python environment, or if you already load your environment in /.bashrc (note: sbatch does not load /.bashrc, source it as per example script). Do not uncomment #SBATCH lines, keep them as below, make sure the #SBATCH directives are directly after #!/bin/bash

Remember to make the sumbission script¹. executable with chmod +x <script_name>.sh

```
#!/bin/bash

#SBATCH — gres=gpu:1

#SBATCH — mail—type=ALL # required to send email notifcations

#SBATCH — mail—user=<your_username> # required to send email notifcations

export PATH=/vol/bitbucket/${USER}/myvenv/bin/:$PATH

# the above path could also point to a miniconda install

# if using miniconda, uncomment the below line

# source ~/.bashrc

source activate

source /vol/cuda/12.0.0/setup.sh

/ usr/bin/nvidia—smi

uptime
```

language=bash

3.4 Scheduling job

- ssh gpucluster2
- cd /vol/bitbucket/\$USER
- sbatch <path to executable>.sh

¹example can be found at /vol/bitbucket/shared/slurmseg.sh

- less slurm-XYZ.out saves the output.
- squeue -1 for queue of running jobs
- scancel <job ID> to delete slurm job

3.5 Creating a SLURM job to act as a VM

```
ssh texel10 # DON'T ssh into a gpucluster machine, it won't work

cd /vol/bitbucket/az620

python -m venv /vol/bitbucket/az620/dlenv

source /vol/bitbucket/az620/dlenv/bin/activate

pip3 install ipykernel jupyterlab jupyterhub

# Add CUDA to the path from /vol/cuda/{version}
```

3.5.1 In Terminal 1

```
ssh gpucluster2
salloc --gres=gpu:1
squeue | grep az620
```

3.5.2 In Terminal 2

If you have defined a config entry for shell* then replace accordingly with az620@shell3.doc.ic.ac.uk \mapsto shell3.

Don't forget leading / in command below

Unmodified:

```
ssh −t −L 10001:localhost:10001 az620@shell3.doc.ic.ac.uk "/vol/linux/bin/slurm_sshtojob.sh −g −w ~/ −p 10001 −e / 
→ vol/bitbucket/{USER}/dlenv"
```

For me:

```
ssh −t −L 10001:localhost:10001 shell3 "/vol/linux/bin/slurm_sshtojob.sh −g −w ~/ −p 10001 −e /vol/bitbucket/

⇒ az620/dlenv"
```

http://localhost:10001/lab?token={from terminal}

```
remember to: scancel <your job id> back in Terminal 1
```

3.5.3 In Jupyter

```
ssh gpucluster1
salloc --gres=gpu:1
squeue
# in VS Code:
ssh -A -J {USER}@shell3.doc.ic.ac.uk {USER}@kingfisher.doc.ic.ac.uk
```

4 SageMaker

Access SageMaker and follow instructions in slides

5 PaperSpace

Remember to close the notebook when finished

Access Paperspace

After access granted, you can access this and use paperspace as your location of the kernel to execute super fast.

6 Python virtual environment

6.1 Where to store data

Data is stored in the /vol/bitbucket/\${USER}. It can be created with mkdir -p ... if it doesnt exist. You should create personal folders here. Otherwise, if you store data in the home directory you risk going over the quota limit [source]². /vol/bitbucket should only be used for temporary storage of material that cna be regenerated (downloads or compilations). To see where you are storing disk space use /vol/linux/bin/usage (disk space) and /vol/linux/bin/nfiles (number of files and directories) [source].

6.1.1 Why do you need a python virtual environment?

Python isn't good at dependency management. pip places all external packages into site-packages/ in the base Python installation. By creating a virtual environment you avoid system polluation (since these packages mix with system-relevant packages causing unexpected side effects)

6.1.2 Creating a Python Virtual Environment in /vol/bitbucket

Use a lab PC to prepare the Python environment; don't use the gpucluster machine for this. You may encounter 'out of space' errors.

Its advised that you create Python virtual environments on /vol/bitbucket. Steps taken from [source]

- 1. create /vol/bitbucket/\${USER}
- 2. create virtual environment

```
export PENV=/vol/bitbucket/${USER}/myenv
python3 -m virtualenv $PENV
ls -al $PENV
```

3. activate your VE:

²you can check disk quota with quota −Q

```
source $PENV/bin/activate
which pip
[should say: /vol/bitbucket/${USER}/myenv/bin/pip]
which python
[should say: /vol/bitbucket/${USER}/myenv/bin/python]
which python3
[should say: /vol/bitbucket/${USER}/myenv/bin/python3]
```

- 4. install packages and verify with which <module name> that it is saved under /vol/bitbucket/\$USER/myenv,
- 5. Or prepare a package requirements file requirements.txt which speicifes a list of packages (optinally with specific version constriants) and install with pip install -r requirements.txt
- 6. once you're done working with the virtual environment you can deactivate it with deactivate
- 7. each time you login to a specific machine redo the activate stage with source /vol/bitbucket/\$USER/myenv (this can be appended to the end of the ~./bash_profile in home dir.)

6.1.3 Why activate and deactivate?

Activating it gives us a new path for the pyhton executable because, in an active environment, the \$PATH environment variable is slightly modified.