#### Cluster

09 April 2025 16:45

prabal@Prabal:~\$ ssh prghosh@access.grid5000.fr

prghosh@access-south:~\$ ssh Nancy

prghosh@fnancy:~\$ oarsub -I -I host=1,walltime=1:00

prghosh@grvingt-2:~\$ ls

prghosh@grvingt-2:~\$ cd p prabal\_1/ public/ prghosh@grvingt-2:~\$ cd prabal\_1/ prghosh@grvingt-2:~/prabal\_1\$ ls

prghosh@fnancy:~/prabal\_1\$ vim script\_fff.sh

Use ECHAP key and press :wq and ENTER to save the script file in

(.verv) prghosh@fsophia:"/prabal\_ghosh/Inria\_Medical\_Imageing\_Internship\_prabal\$ cd ...
(.verv) prghosh@fsophia:"/prabal\_ghosh/Inria\_Medical\_Imageing\_Internship\_prabal\$ cd ...
(.verv) prghosh@fsophia:"/prabal\_ghosh\$ cd Inria\_Medical\_Imageing\_Internship\_prabal\$ (.verv) prghosh@fsophia:"/prabal\_ghosh\$ cd Inria\_Medical\_Imageing\_Internship\_prabal\$ (.verv) prghosh@fsophia:"/prabal\_ghosh\$ county prghosh@fsophia:"/prabal\_ghosh\$ county prghosh@fsophia:"/prabal\_ghosh\$ county prghosh@fsophia:"/prabal\_ghosh\$ county prghosh@fsophia:"/prabal\_ghosh\$ county prghosh@fsophia:"/prabal\_ghosh\$ county prghosh@fsophia:"/prabal\_ghosh\$ cluster\_lest\_0005.pyp county prghosh@fsophia:"/prabal\_ghosh\$ cluster\_lest\_0005.pyp county prghosh@fsophia:"/prabal\_ghosh\$ cluster\_lest\_0006.pyp county prghosh@fsophia:"/prabal\_ghosh\$ cluster\_lest.sh cluster\_les

oarstat -u # Check status cat OAR\_\${OAR\_JOB\_ID}.out # View output cat OAR\_\${OAR\_JOB\_ID}.err # View errors oarsub -S ./cluster\_1\_test.sh

#### 1. For interactive jobs (if you really need one):

oarsub -I -I host=1/gpu=1,walltime=1:00:00 -p \*gpu='16GB' AND gpu\_compute\_capability\_major>=5" -q production

Then manually run

./cluster\_1\_test.sh

oarstat -u # Check status cat OAR\_\${OAR\_IOB\_ID}.out # View output cat OAR\_\${OAR\_IOB\_ID}.err # View errors

oarsub -I -q production -l gpu=1,walltime=0:05:00

module load conda module load cuda/11.8 conda activate mri\_2025\_4 cd ~/prabal\_ghosh/Inria\_Medical\_Imageing\_Internship\_prabal python cluster\_test\_0005.py

/usr/bin/oarsub -I -q production -I gpu=1,walltime=0:05:00

chmod +x cluster\_1\_test.sh

/path/to/cluster\_1\_test.sh

./cluster\_1\_test.sh

oarsub -S ./cluster\_1\_test.sh

UPLOAD DATA

prabal@Prabal:~\$ scp -r Inria\_Medical\_Imageing\_Internship\_prabal

 $\label{probability} $$ $$ prabal@Prabal:^$ scp test_s0005.ipynb $$ prghosh@access.grid5000.fr.sophia/inria_Medical_Imageing_Internship_prabal_1/$$ $$ $$ prabal_1/$$ $$ prabal_2/$$ $$ p$ 

DOWNLOAD DATA

mri\_2025\_4

this virtual environment is working for cpu in cluster

### **Inside Ubuntu terminal**

prabal@Prabal:~\$ ssh prghosh@access.grid5000.fr prghosh@access-north:~\$ ssh sophia

## Creation of virtual environnement

Conda create -n virtual\_prabal python=3.10.0

Conda activate virtual\_prabal

## Installation of pytorch with gpu cuda 11.8check these 2 methods inside your virtual environment

pip install torch==2.3.1 torchvision==0.18.1 torchaudio==2.3.1 --index-url  $\frac{1}{1000} \frac{1}{1000} \frac{1}{1000}$ 

From <https://pytorch.org/get-started/previous-versions/>

conda install pytorch==2.3.1 torchvision==0.18.1 torchaudio==2.3.1 pytorch-cuda=12.1 -c pytorch -c nvidia

From <https://pytorch.org/get-started/previous-versions/>

pip install TotalSegmentator

Now the setup is done. There might be some errors, which you will have to fix by looking at the error messages during code execution

## This is the python script

### Vim cluster\_test\_0005.py

```
import nibabel as nib
from pathlib import Path
from totalsegmentator.python_api import totalsegmentator
# import numpy as np
# import os
# import os
# import astplotlib.pyplot as plt
input_path_1 =
Path("/home/prghosh/prabal_ghosh/Inria_Medical_Imageing_Internship_prabal/s0005/m
ri.nii.g2")
output_path_1 =
Path("/home/prghosh/prabal_ghosh/Inria_Medical_Imageing_Internship_prabal/s0005/s
egmentations_test_2")
img = nib.load(input_path_1).get_fdata()
print(img.shape)
```

 $\label{eq:print} \text{print}(\texttt{f"*****The .nii files are stored in memory as numpy's: } \{\texttt{type}(\texttt{img})\}. \texttt{******"})$ 

\_\_\_\_\_\_

#### 1. Method1:

## This is the shell script

```
vim prabal_test_shell.sh
```

# Inside this .sh file write the following code

```
#!/bin/bash
#OAR -q production
#OAR -1 host=1/gpu=1
#OAR -1 walltime=00:30:00
#OAR -2 ppu_count > 0
#OAR -0 ppu_count > 0
#OAR -0 OAR %jobid%.out
#OAR -E OAR %jobid%.out
#OAR -E OAR %jobid%.out
#OAR -E OAR %jobid%.err
# display some information about attributed resources
echo "=== Host and GPU Info ==="
hostname
nvidia-smi
nvcc --version
echo "=== Loading environment ==="
module load conda
module load cuda/11.8
conda activate mri_2025_4
echo "=== Checking PyTorch GPU Availability ==="
python3 -c "import torch; print('CUDA available:', torch.cuda.is_available());
print('Device:', torch.cuda.get_device_name(a))"
echo "=== Starting TotalSegmentator Job ==="
cd /home/prghosh/prabal_ghosh/Inria_Medical_Imageing_Internship_prabal
#Run your script
python cluster_test_0005.py
echo "=== Done ==="
conda deactivate
```

# Now to run the shell script use the Following commands

```
chmod +x prabal_test_shell.sh # is to check that its exécutable or not

oarsub -S ./prabal_test_shell.sh # to run the script

oarstat -u # Check status
cat OAR_$[OAR_JOB_ID]\out # View output
cat OAR_$[OAR_JOB_ID]\out # View errors
```

## 2. Method2:

If you want to run it interactively, use the following commands — you don't need to write any shell script file.

oarsub -I -q production -l gpu=1,walltime=0:05:00

module load conda module load cuda/11.8 conda activate mri\_2025\_4 cd "/prabal\_ghosh/Inria\_Medical\_Imageing\_Internship\_prabal python cluster\_test\_0005.py