

Requesting and using GPUs

Both Curnagl and Urblauna have nodes with GPUs.

You can find a detailed description of Curnagl GPUs [here](#) and Urblauna GPUs [here](#)

An introductory tutorial video on using HPC clusters is available [here](#).

Requesting GPUs

In order to access the GPUs they need to be requested via SLURM as one does for other resources such as CPUs and memory.

The flag required is `--gres=gpu:1` for 1 GPU per node, you can use any number between 1 and N (`--gres=gpu:N`). Please check cluster documentation.

An example job script is as follows:

```
#!/bin/bash -l

#SBATCH --cpus-per-task 12
#SBATCH --mem 64G
#SBATCH --time 12:00:00

# GPU partition request only for Curnagl
#SBATCH --partition gpu

#SBATCH --gres gpu:1
#SBATCH --gres-flags enforce-binding

# Set up my modules

module purge
module load my list of modules
module load cuda

# Check that the GPU is visible
```

```
nvidia-smi

# Run my GPU enable python code

python mygpucode.py
```

If the `#SBATCH --gres gpu:1` is omitted then no GPUs will be visible even if they are present on the compute node.

If you request one GPU it will always be seen as device 0.

The `#SBATCH --gres-flags enforce-binding` option ensures that the CPUs allocated will be on the same PCI bus as the GPU(s) which greatly improves the memory bandwidth. This may mean that you have to wait longer for resources to be allocated but it is strongly recommended.

Using CUDA

In order to use the CUDA toolkit there is a module available

```
module load cuda
```

This loads the nvcc compiler and CUDA libraries. There is also a cudnn module for the DNN tools/libraries

Containers and GPUs

Singularity containers can make use of GPUs but in order to make them visible to the container environment an extra flag "--nv" must be passed to Singularity

```
module load singularity

singularity run --nv mycontainer.sif
```

The full documentation is at <https://sylabs.io/guides/3.5/user-guide/gpu.html>

you can find [here](#), examples of using GPUs from containers.

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