### How to connect to the system

ssh -X z3XX@login-gfx2.hpc.cam.ac.uk

If you want to use Windows please download **MobaXterm** (http://mobaxterm.mobatek.net)

### Access all material

https://github.com/RSE-Cambridge/2017-cuda-fortran-material

(short url: https://tinyurl.com/CUDAFORTRAN)

git clone https://github.com/RSE-Cambridge/2017-cuda-fortran-material.git

## **Environment HW/SW**

#### HW (Wilkes, Cambridge)

- dual socket 6-core Ivy Bridge (12 cores total)
- 2 NVIDIA GPU K20

#### **SW**

- PGI 16.10 (community edition)
- MKL 11.3
- Open MPI 1.10.3
- CUDA 8.0
- Few custom libraries (for QE-GPU)

#### SLURM commands

Three command to learn ...

sbatch <script>

squeue -u \$USER

scancel <jobid>

# Submission scripts

- submit-monday.sh
- submit-tuesday.sh
- submit-wednesday.sh
- submit-serial.sh
- submit.sh

# Example submission script

```
#!/bin/bash
#SBATCH -J NAME
#SBATCH - A TRAINING-GPU
#SBATCH --nodes=UUUUU
#SBATCH --ntasks=XXXXX
#SBATCH --time=0:05:00
#SBATCH --no-requeue
#SBATCH --partition=tesla
#SBATCH --reservation=cuda_fortran_mon
. /etc/profile.d/modules.sh
module purge
module load default-wilkes
module unload cuda intel/cce intel/fce intel/impi intel/mkl
module load intel/mkl/11.3.3.210
module load pgi/16.10
module load openmpi/pgi/1.10.3
module load cuda/8.0
module load custom/magma/2.0.2 custom/lib-jdr/1.0
cd $SLURM SUBMIT DIR
export EXE="./pw.x"
export PARAMS="-input pw.in"
export OMP NUM THREADS=6
mpirun -np XXXXX -report-bindings -bind-to socket --map-by ppr:1:socket ${EXE} ${PARAMS}
```

## Compile and run QE-GPU code

Two examples: "PW-AUSURF112" & "PW-AUSURF54" (k and gamma)

#### Compile:

```
cd qe-gpu_mini
./configure --enable-parallel --enable-openmp -without-scalapack
cp make.inc-STELLA make.inc
make -j4 pw
```

(executable located in "PW/src/pw.x")

#### Run:

```
export EXE="./pw.x"
export PARAMS="-input ausurf_k.in"

export OMP_NUM_THREADS=6
mpirun -np XXXXX -report-bindings \
  -bind-to socket --map-by ppr:1:socket ${EXE} ${PARAMS}
```