Introduction to practical sessions, and logistics





Generalities

We have a total of 7 hours of scheduled practical sessions for the MPAS-A part of this course

During the practical sessions, you can work through the exercises at your own pace

Instructors will be in the practical sessions in case you have questions or would like to discuss anything covered in the lectures



Derecho

The practical sessions of this tutorial will be run on Derecho





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323,712 processor cores	3 rd Gen AMD EPYC™ 7763 Milan processors
2,488 CPU-only computation	Dual-socket nodes, 64 cores per socket
nodes	256 GB DDR4 memory per node
82 GPU nodes	Single-socket nodes, 64 cores per socket
	512 GB DDR4 memory per node
	4 NVIDIA 1.41 GHz A100 Tensor Core GPUs per node
	600 GB/s NVIDIA NVLink GPU interconnect
328 total A100 GPUs	40GB HBM2 memory per GPU
	600 GB/s NVIDIA NVLink GPU interconnect



Derecho



Don't forget to enable X11 forwarding when connecting!

ssh -X <username>@derecho.hpc.ucar.edu

On macOS systems you might try -Y rather than -X

ssh -Y <username>@derecho.hpc.ucar.edu



Queues

For part of each day (11:30 am – 5:00 pm) we have 30 nodes reserved for this tutorial

- Submit jobs to the tutorial queue
- This is the default in the job scripts that we provide

Since we have 22 participants and only 30 nodes, please try to only have one job queued via qsub at a time

Feel free to work on exercises outside of scheduled practice sessions

- However, submit jobs to the main queue instead
- All that you need to do is to uncomment one line and comment another in your job script

¹In the tutorial queue



Job scripts

```
#!/bin/bash
                                Job scripts are
                                submitted to the
#PBS -A UMMM0014
                                queuing system
                                with qsub
#PBS -N jobname
#PBS -q main
#PBS -l job priority=premium
#PBS -1 walltime=00:30:00
#PBS -l select=1:ncpus=128:mpiprocs=128
mpiexec ./atmosphere model
```



Checking on queued jobs

The qstat command is used on Derecho to check the status of queued jobs

- qstat with no arguments shows <u>all</u> jobs on the system, which is usually not useful
- qstat -u \$USER shows only your jobs

But...

The information shown by qstat is only refreshed every 30 seconds or so. Therefore, after submitting a job, it may take a bit before you see your job with qstat. If you think your job may have already finished, try qstat -xu \$USER to see running and completed jobs.



Running with qcmd

For some parts of practical exercises, we will run programs on a batch node without a job script via qcmd

Any options that you can provide in a job script can be given to qcmd, e.g.,



Resource allocations

All of our computing will be under the UMMM0014 project

We have 50,000 core-hours

- Divided evenly between 22 participants -> everyone has about 2200 core-hours
- One Derecho node for one wallclock hour = 128 core-hours
- You can continue to work with MPAS-Atmosphere on your own time

The UMMM0014 project is available until the end of September

Use the UMMM0014 project *only for MPAS- Atmosphere* tutorial work, and please don't hesitate to contact one of the tutorial organizers if you have any questions at all!



We have a web page to guide you through some exercises:

https://www2.mmm.ucar.edu/projects/mpas/tutorial/Boulder2025/



Welcome to the MPAS tutorial practice guide

This web page is intended to serve as a guide through the practice exercises of this tutorial. Exercises are split into eight main sections, each of which focuses on a particular aspect of using the MPAS-Atmosphere model.

While going through this practice guide, you may find it helpful to have a copy of the MPAS-Atmosphere Users' Guide available in another window. Click here to open the Users' Guide in a new window.

In case you would like to refer to any of the lecture slides from previous days, you can open the <u>Tutorial Agenda</u> in another window.

You can proceed through the sections of this practical guide at your own pace. It is highly recommended to go through the exercises in order, since later exercises may require the output of earlier ones. Clicking the grey headers will expand each section or subsection.

- 0. Prerequisites and environment setup
- 1. Compiling MPAS, and creating static files and idealized ICs
- 2. Generating intermediate files from GRIB and netCDF datasets



We have a web page to guide you through some exercises:

https://www2.mmm.ucar.edu/projects/mpas/tutorial/Boulder2025/

If you're logged in to Derecho, you can get this URL with:

cat ~duda/url



Blocks of shell commands can be selected all at once

Before going through any of the exercises in subsequer to prepare our environment. Firstly, we will purge all lo point (with no modules loaded); then, we will load seve as well as an MPI implementation:

```
$ module --force purge
$ module load ncarenv/23.09
$ module load craype/2.7.23
$ module load gcc/13.2.0
$ module load ncarcompilers/1.0.0
$ module load cray-mpich/8.1.27
```



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```



Environment setup

As noted in the practice guide, you should **set up your environment every time you log in**:

If you've logged out of Derecho and would like a quick list of commands to set up your environment again after logging back in, you can copy and paste the following commands into your terminal:

```
module --force purge
module load ncarenv/23.09
module load craype/2.7.23
module load gcc/13.2.0
module load ncarcompilers/1.0.0
module load cray-mpich/8.1.27
module load parallel-netcdf/1.12.3
module load netcdf/4.9.2
module load conda
conda activate npl
export PYTHONPATH=/glade/campaign/mmm/wmr/mpas_tutorial/python_scripts
module load ncview
export PATH=/glade/campaign/mmm/wmr/mpas_tutorial/metis/bin:${PATH}
module reload gcc/13.2.0
```



Asking questions

After the conclusion of the tutorial, if you encounter issues or have any questions, the forum might be a good place to discuss those:

https://forum.mmm.ucar.edu/