MPI (Message Passing Interface)

Outline

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 - Area_curve_mpi
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- Run the job script
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 - Hello_world_mpi
 - Area curve code
 - BioFilm
 - More
- Run job
- Output

Structure of MPI code

```
#include <mpi.h>
ierr=MPI_Init(&argc, &argv);

    ierr=MPI Comm size(MPI COMM WORLD,&npes);

ierr=MPI_Comm_rank(MPI_COMM_WORLD,&iam);
ierr=MPI Finalize();
```

Structure of MPI Code

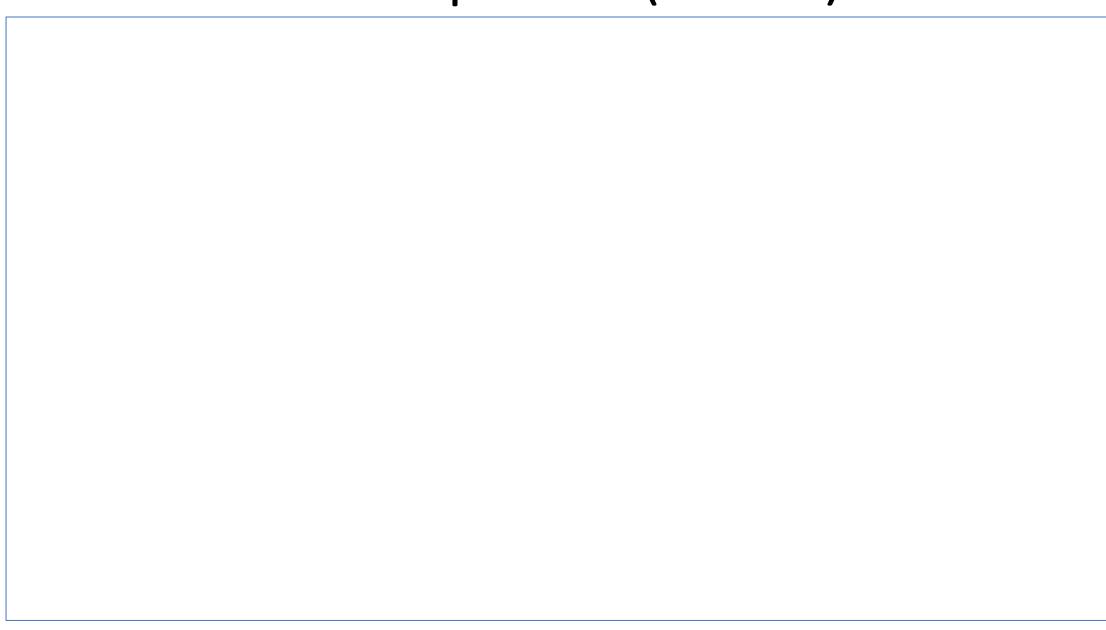
- A parallel "Hello, World" program
- – Initialize MPI
- Have each node print out its node number
- Quit MPI

- #include <mpi.h>
- •
- ierr=MPI_Init(&argc, &argv);
- ierr=MPI_Comm_size(MPI_COM M_WORLD,&npes);
- ierr=MPI_Comm_rank(MPI_CO MM_WORLD,&iam);
- •
- ierr=MPI_Finalize();

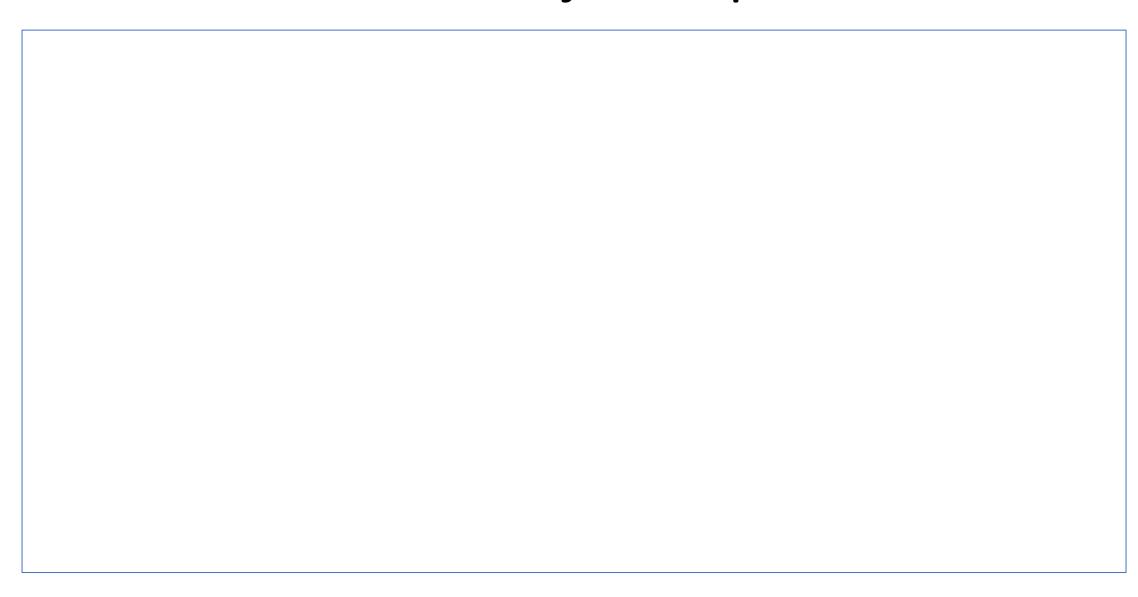
Load the module (software application)

- module load OpenMPI
- sbatch myJobScript
- squeue –u haboudj

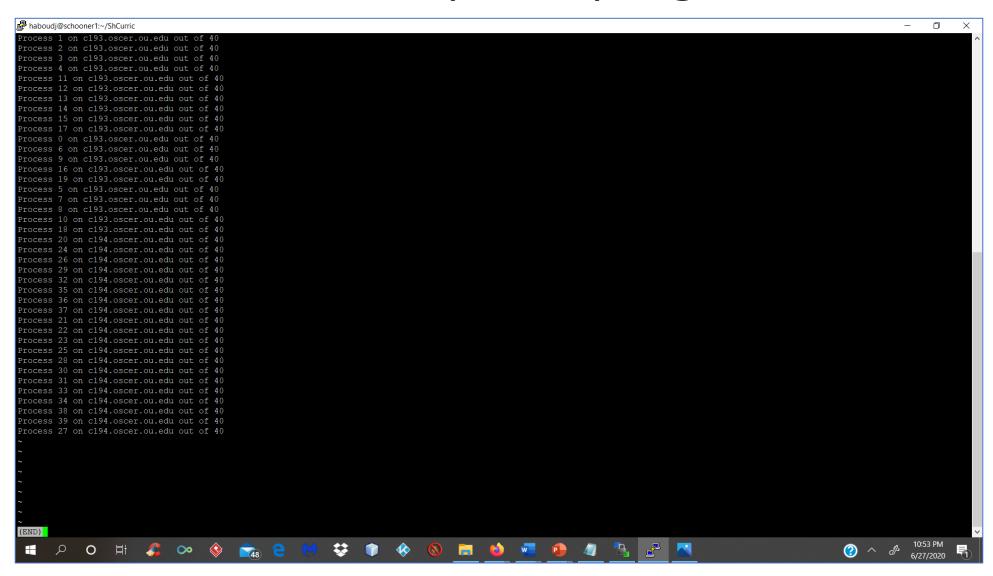


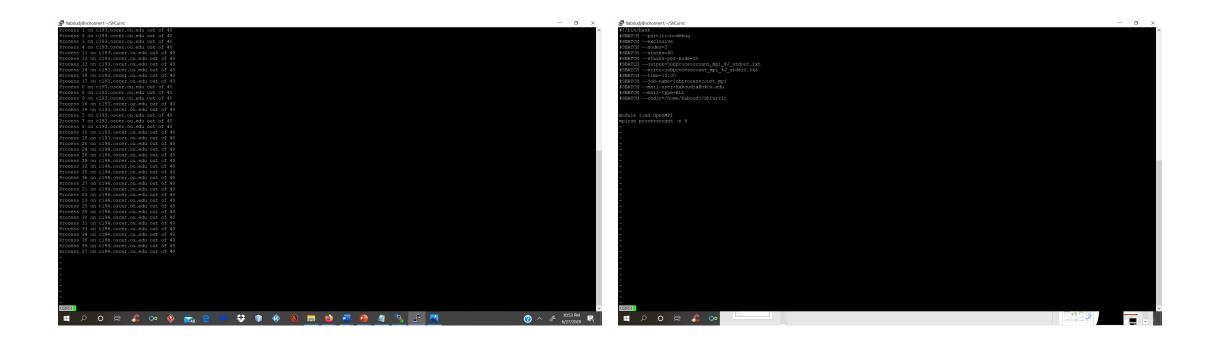


Run the job script



Screenshots of example of program codes run





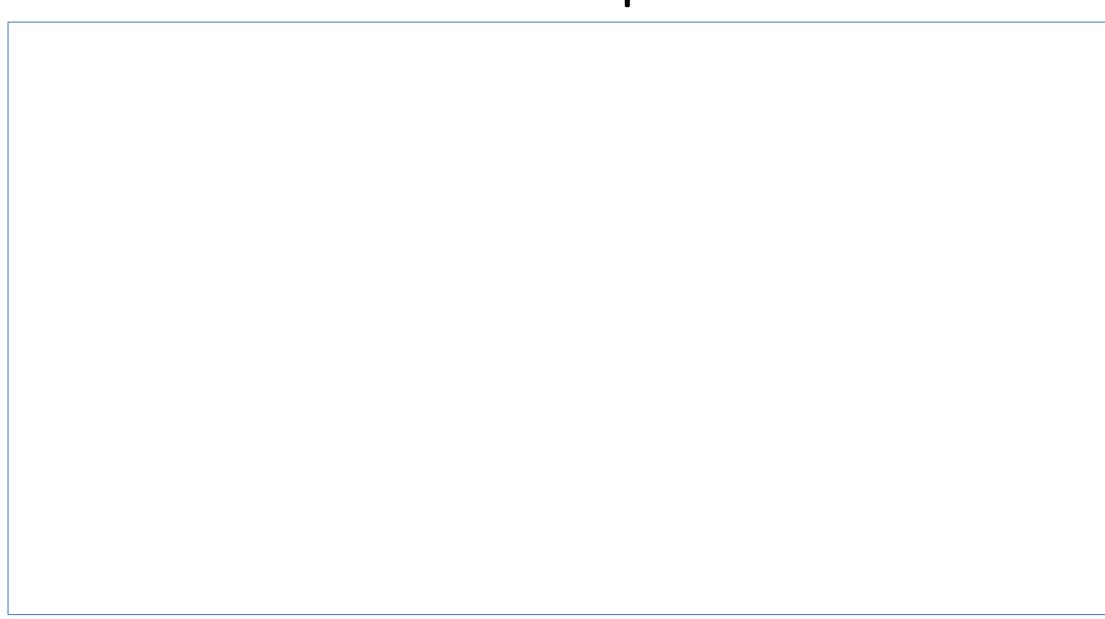
Steps to Use a Super Computer (Example Template)

ssh username@hostname.ncsa.illinois.edu<ENTER> git clone \ https://github.com/aaronweeden/pi2018-submitting.git<ENTER> cd pi2018-submitting<ENTER> cc -o test.exe test.c<ENTER> cat test.pbs<ENTER> #!/bin/bash #PBS -I nodes=2:ppn=32:xe **#PBS -I walltime=00:05:00 #PBS -N test** cd \$PBS_O_WORKDIR aprun -n 4 ./test.exe|sort

Running Job Script

- cat test.pbs<ENTER>
- #!/bin/bash
- #PBS -l nodes=2:ppn=32:xe
- #PBS -I walltime=00:05:00
- #PBS -N test
- cd \$PBS_O_WORKDIR
- aprun -n 4 ./test.exe|sort

Error and Output Files



References

- 1. https://people.sc.fsu.edu/~jburkardt/c src/laplace mpi/laplace mpi.html
- 2. Please see the TACC CS395 COURSE MATERIALS s1.
- 3. https://www.mpi-forum.org/docs/mpi-1.1/mpi1-report.pdf