# Lab 1 - Instructions and Summary

## Compiling a MPI program

```
Load Intel compilers
```

```
module load intel/psxe-2015
```

Compile for speed -Ofast and add information about line numbers -g

```
mpicc -g -Ofast -o myprogram myprogram.c
mpifc -g -Ofast -o myprogram myprogram.f90
```

Compile for debugging -00 and add information about line numbers -g

```
mpicc -g -00 -o myprogram myprogram.c
mpifc -g -00 -o myprogram myprogram.f90
```

## Running an MPI program on Yeti

Job script for an MPI run on normal partion

```
#!/bin/bash
#SBATCH -p normal
#SBATCH -t 0:02:00
#SBATCH -N 2
#SBATCH --job-name=mpi_hello
#SBATCH --output=out.mpi_hello
#SBATCH --ntasks-per-node=20
module load intel/psxe-2015
srun --mpi=pmi2 ./mpi_hello
```

### Writing a parallel "hello world" program

- All process print: "Hello World from process = %d processor %s"
- Processor 0 prints: "Number of mpi processes = %d"

#### Fortran 90

```
use mpi
MPI_COMM_WORLD
MPI_MAX_PROCESSOR_NAME
MPI_INIT(IERROR) ! integer argument
MPI_COMM_SIZE(COMM, SIZE, IERROR) ! all arguments are integer
MPI_COMM_RANK(COMM, RANK, IERROR) ! all arguments are integer
MPI_FINALIZE(IERROR) ! integer argument
```

### $\mathbf{C}$

```
MPI_Comm MPI_COMM_WORLD
MPI_MAX_PROCESSOR_NAME
int MPI_Init(int *argc, char ***argv)
int MPI_Comm_size(MPI_Comm comm, int *size)
int MPI_Comm_rank(MPI_Comm comm, int *rank)
int MPI_Finalize()
```

# Writing a MPI version of the PI program

- Process 0 reads the number of integration intervalls
- Process 0 broadcasts this number to the rest of the processes
- Each process computes its share of PI
- Process 0 receives all shares and calculates an approximation of pi
- Process 0 prints the result and the error

### Fortran 90