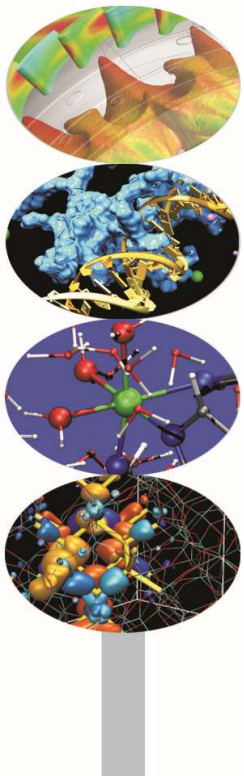


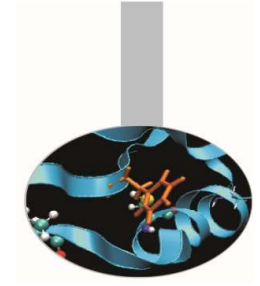
Advanced MPI

- *exercises* -

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Startup notes

Access the server and go on scratch partition:

```
ssh a08tra49@login.galileo.cineca.it
```

```
cd $CINECA_SCRATCH
```

Create a job file <test.job>:

```
#!/bin/bash
```

```
#PBS -A train_cmpD2016
```

```
#PBS -l select=1:ncpus=4:mpiprocs=4:mem=32gb
```

```
#
```

```
cd $PBS_O_WORKDIR
```

```
module load autoloader openmpi
```

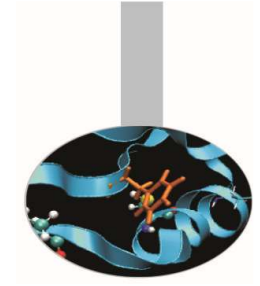
```
rm *.exe
```

```
mpicc c_program.c -o c_program.exe
```

```
mpif90 f_program.f90 -o f_program.exe
```

```
mpirun <program>.exe
```

```
exit
```



Startup notes

Other useful PBS flags:

```
#PBS -N test          <- job name
```

```
#PBS -j oe           <- join stdout and stderr outputs
```

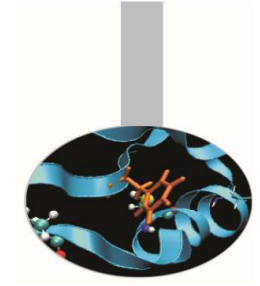
```
#PBS -l walltime=1:00:00      <- max 24 hours
```

Submit the job to the batch queue:

```
qsub <test.job>
```

Check if the job is running:

```
qstat -n -u a08tra49
```



Compiling notes

To compile programs that make use of MPI library:

```
mpif90/mpicc/mpicc -o <executable> <file 1> <file 2> ... <file n>
```

Where: <file n> - program source files

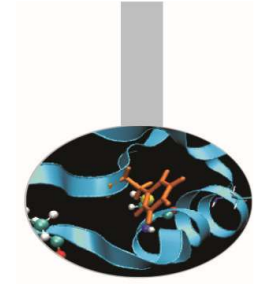
<executable> - executable file

To start parallel execution on one node only:

```
mpirun -np <processor_number> <executable> <exe_params>
```

To start parallel execution on many nodes:

```
mpirun -np <processor_number> -machinefile <node_list_file> \  
    <executable> <exe_params>
```



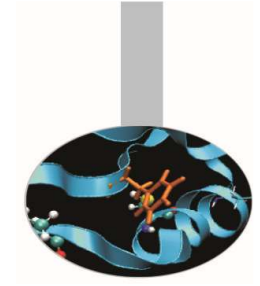
MPI exercises

You can find a list of useful exercises at the address:

<http://www.hpc.cineca.it/content/training-mpi>

Exercise 10: create a cartesian topology and try using `Cart_shift` for neighborhood communications

Exercise 11: MPI data types are useful for managing matrix data whenever not-contiguous data are involved. To solve the exercise define a data type that include two columns(in C)/rows(in Fortran)

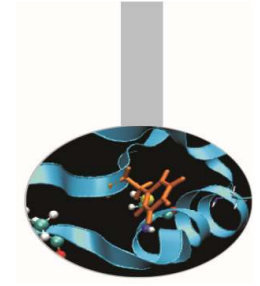


MPI exercises

Exercise 12: *again on matrix data: send and receive buffers do not need to be of the same type. The flag `-std=c99` should be used to compile the C solution program with `gcc`*

Exercise 13: *I/O functions are used to store and retrieve data; the function `File_set_view` is used to get `File_write_all` to store data in the specified order*

Exercise 14: *use of `MPI_Dims_create` is not mandatory; remember to call `File_set_view` before using `File_write_all` and `File_read_all`. Use the flag `-std=c99` to compile with `gcc`*



MPI exercises

Exercise 15: the function Alltoall redistribute an array in an ordered fashion; remember that sendcount and recvcount are the dimensions of the sub-blocks to be sent/received to/from single processes

Exercise 16: the function Allgather performs collection of distributed data to all processes