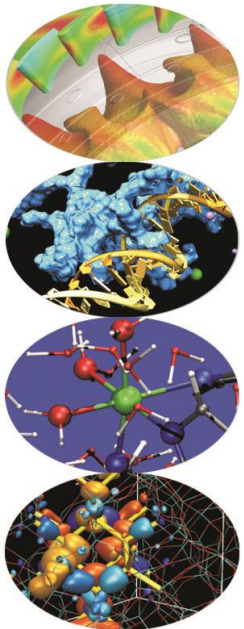
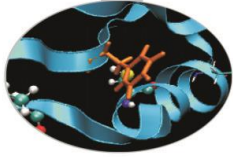


MPI introduction

- *exercises* -





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

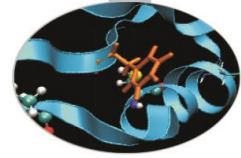
<executable> - executable file

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>
```



Hello world! (Fortran)

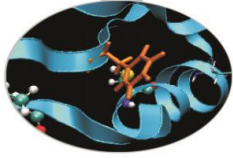
As an ice breaking activity try to compile and run the *Hello* program, either in C or in Fortran.

The most important lines in Fortran code are emphasized:

```
PROGRAM HelloWorld
  INCLUDE 'mpif.h'
  INTEGER my_rank, p
  INTEGER source, dest, tag
  INTEGER ierr, status(MPI_STATUS_SIZE)
  . . .
  CALL MPI_Init(ierr)
  CALL MPI_Comm_rank(MPI_COMM_WORLD, my_rank, ierr)
  CALL MPI_Comm_size(MPI_COMM_WORLD, p, ierr)

  WRITE(*,FMT="(A,I)") "Hello world from process ", my_rank

  CALL MPI_Finalize(ierr)
END PROGRAM HelloWorld
```

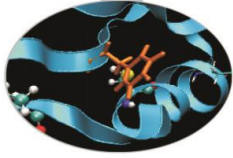


Hello world! (C/C++)

The most important lines in C code are emphasized:

```
#include "mpi.h"
```

```
int main( int argc, char *argv[])  
{  
    int my_rank, numprocs;  
    int dest, tag, source;  
    MPI_Status status;  
  
    MPI_Init(&argc,&argv);  
    MPI_Comm_rank(MPI_COMM_WORLD,&my_rank);  
    MPI_Comm_size(MPI_COMM_WORLD,&numprocs);  
  
    printf("Hello world from process %d\n",my_rank);  
  
    MPI_Finalize();  
    return 0;  
}
```



Hello world! (output)

If the program is executed with one process the output is:

```
Hello world from process 0
```

If the program is executed with four processes the output is:

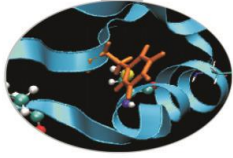
```
Hello world from process 0
```

```
Hello world from process 1
```

```
Hello world from process 2
```

```
Hello world from process 3
```

E1 – exercise – ping-pong

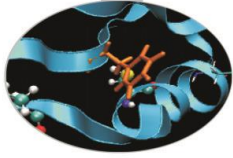


ping-pong is perhaps the simplest example of point to point communication.

In a two process execution of a ping-pong program the process 0 sends a message to process 1 and this sends it back to process 0. This could be easily generalized in a round robin fashion if more than two processes are engaged.

Try modifying the Hello World example in order of realizing round robin communications.

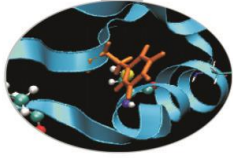
E2 – example – Pi by quadrature



It is known that the mathematical constant π can be approximated by computing the following formula:

$$\pi = 4 \int_0^1 \frac{1}{1+x*x} dx$$

The value of the above integral can be approximated by numerical integration, i.e. by computing the mean value of the function $f(x) = \frac{1}{1+x*x}$ in a number of points and multiplying per the x range. This can be easily done in parallel by dividing the $[0,1]$ range into a number of intervals.



E2 – example – Pi by quadrature

Thus the program may be sketched this way:

- (if `my_rank == 0`) get number of intervals for quadrature
- Broadcast number of intervals to all the processes
- Assign the intervals to the processes (they should not overlap)
- Sum function values in the centre of each interval
- Divide by interval range and multiply by 4

Source code: *Pi_integral*