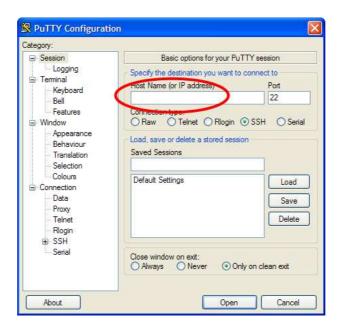
## Lab 1: Introduction to MPI

- 1. Download and install required programs
- 1.1 Download PuTTY from <a href="http://the.earth.li/~sgtatham/putty/latest/x86/putty.exe">http://the.earth.li/~sgtatham/putty/latest/x86/putty.exe</a>, save it on Desktop.
  - 1.2 Download WinSCP from https://winscp.net/eng/download.php, install it.
- 2. Use PuTTY to login to cluster.ict.mahidol



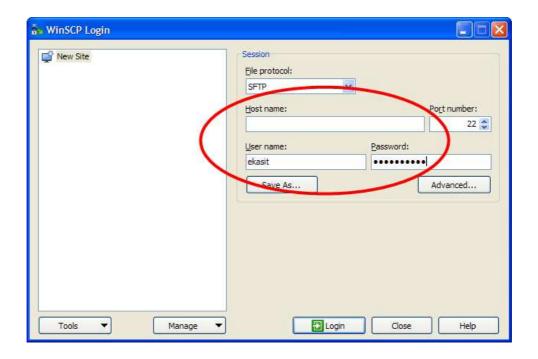
- 2.1 Enter your login and password
- 2.2 For the first time login, press Enter 3 times until you get the command prompt [yourlogin@cluster ~]\$

- 2.3 Type 1s command to list all files in your home directory. For first time login, you should not see any file.
- 3. Create a helloworld.c MPI program.
- 3.1 Use any editor in your machine, e.g notepad, notepad++, editplus, eclipse, or visual C, to create the following program.

```
#include "mpi.h"
#include <stdio.h>

int main(int argc,char *argv[] )
{
   int rank, size;
   char hostname[MPI_MAX_PROCESSOR_NAME];
   int resultlen;
   MPI_Init( &argc, &argv );
   MPI_Comm_rank( MPI_COMM_WORLD, &rank );
   MPI_Comm_size( MPI_COMM_WORLD, &size );
   MPI_Get_processor_name(hostname,&resultlen);
   printf( "Hello! I'm %d of %d running on %s\n",rank,size,hostname);
   MPI_Finalize();
   return 0;
}
```

3.2 Save the program. Use WinSCP to upload it to your home directory on cluster.ict.mahidol.



- 3.3 In PuTTY, type 1s and you must see the helloworld.c in home directory.
- 3.4 Compile the program. Type

mpicc -o helloworld helloworld.c

- 3.5 If there is an error, correct it in the editor and re-upload it again. If no error, you'll see the program helloworld (without .c extension) in your home directory.
- 3.6 Run the program with 2 processes. Type mpirun -np 2 helloworld

You will see the output. Try running program with 3,4,5 and 6 processes (change the –np argument).

4. Create and run the integersum.c program as follows.

```
#define LEFT 1
#define RIGHT 1000
#include <stdio.h>
#include <mpi.h>
int main(int argc, char *argv[])
   int rank, size;
   MPI_Status status;
   int interval;
   int number, start, end, sum, GrandTotal;
   int proc;
   MPI Init( &argc, &argv );
   MPI Comm rank( MPI COMM WORLD, &rank );
  MPI Comm size( MPI COMM WORLD, &size );
   if (rank == 0) {
      GrandTotal = 0;
      for (proc=1; proc<size; proc++) {</pre>
         MPI_Recv(&sum,1,MPI_INT,proc,123,MPI_COMM_WORLD,&status);
         GrandTotal=GrandTotal+sum;
      printf("Grand total = %d \n", GrandTotal);
   }
   else {
      interval=(RIGHT-LEFT+1)/(size-1);
      start=(rank-1)*interval+LEFT;
      end=start+interval-1;
      if (rank == (size-1)) {/* for last block */
         end = RIGHT;
      }
      sum=0; /*Sum locally on each proc*/
      for (number=start; number<=end; number++)</pre>
         sum = sum+number;
      /*send local sum to Master process*/
      MPI Send(&sum, 1, MPI INT, 0, 123, MPI COMM WORLD);
   MPI Finalize();
}
```

5. Modify the integersum.c program such that it prints out the starting number, ending number, and local summation computed by each slave process. The following is one possible output.

```
Rank 2: start 251, end 500, local sum 93875
Rank 1: start 1, end 250, local sum 31375
Rank 4: start 751, end 1000, local sum 218875
Grand total = 500500
Rank 3: start 501, end 750, local sum 156375
```