IN3200/IN4200 Exercise Set 11

Exercise 1

Write a simple MPI program to measure the overhead of point-to-point communication involving `MPI_Send` and `MPI_Recv`. (Hint: you can start two MPI processes, between which a message is bounced back and forth a number of times.)

Exercise 2

Write a parallel version of the following `trapezoidal` function using MPI. (Hint: The `for`-loop needs to be modified a little bit to possess parallelism.)

```
double trapezoidal (int n) {
  double result = 0.0;
  double h = 1.0/n;
  double x;
  int i;

  x = 0.0;
  for (i=1; i<n; i++) {
    x += h;
    result += exp(5.0*x)+sin(x)-x*x;
  }

  x = 0.;
  result += 0.5*(exp(5.0*x)+sin(x)-x*x);

  x = 1.0;
```

```
  result += 0.5*(exp(5.0*x)+sin(x)-x*x);

  return (h*result);
}
```

Exercise 3

Write an MPI program that parallelizes the following numerical computation:

```
#include <malloc.h>
#include <math.h>

/* allocating three 1D arrays um, u, up of length M+2 */
/* ... */

double x, dx = 1.0/(M+1);
double t, dt = dx;
double *tmp;
int i;

for (i=0; i<=M+1; i++) {
  x = i*dx;
  um[i] = sin(2.0*M_PI*x);
}

for (i=1; i<=M; i++)
  u[i] = um[i] + 0.5*(um[i-1]-2*um[i]+um[i+1]);

u[0] = u[M+1] = 0.0;

t = dt;
while (t<1.0) {
  t += dt;
  for (i=1; i<=M; i++)
    up[i] = um[i]+u[i-1]+u[i+1];
  up[0] = up[M+1] = 0.0;

  /* shuffle the three arrays */
  tmp = um;
```

```
  um = u;
  u = up;
  up = tmp;
}
```

Hint: The above computation may arise from numerically solving a very simple 1D wave equation. Ignoring the mathematical and numerical details, it is sufficient to notice that three arrays are involved, where computing values in array `up` needs values from arrays `u` and `um`. In particular, values of array `up` can be computed in any random order (thus the existence of parallelism).