

## Project 1, due on 02/20.

### Problem 1. Parallel Numerical Integration for Undergraduate Students.

Evaluate  $\int_{2.0}^{5.0} \int_{2.0}^{5.0} \ln(x + 2y)e^{\sqrt{x^2+y^2}} dydx$ . Implement a parallel code using composite Gaussian quadrature rule to approximate this definite integral.

Suppose  $P$  processes are used and the integration domain  $[2.0,5.0] \times [2.0,5.0]$  is partitioned into  $M \times M$  grid blocks. Each of the processes is assigned with a sub-region  $[x_{i,l}, x_{i,u}] \times [y_{i,l}, y_{i,u}]$ , which is partitioned into  $(M/\sqrt{P}) \times (M/\sqrt{P})$  blocks. Here  $i = 0, 1, \dots, \sqrt{P}-1$ .

We apply the 2D Gaussian quadrature rule to each of these grid blocks to compute a numerical quadrature value. The approximation to the given integral is obtained by summing up these numerical quadrature values.

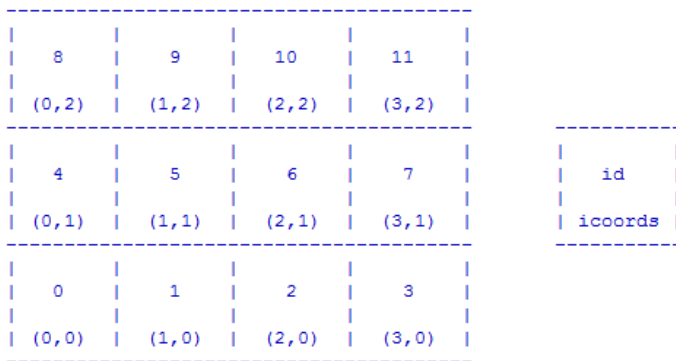
Use the framework implemented in `parallel_trapezoidal.c`. Modify the script “`HPCC_1.sh`” to submit your runs.

1. Use point-to-point communication, specifically, non-blocking send and blocking receive to transfer the quadrature values computed by each of the processes to process 0 and let process 0 compute the sum of these quadrature values.
2. Use  $M = 10000, 20000$  and  $40000$  to do the calculation respectively. For each computation, use 4, 16 and 64 processors respectively. Find the overall the wall clock times spent by the computation, and the communication respectively. Make a table to list the results.

**Hand-In.** Turn in the hardcopy of all your source code, and the report which contains results and a description of your implementation on point-to-point communication. Email the source code.

### Coding Hints.

1. Defines an assignment of processes to subdomains.



**find\_Cartesian\_coordinates()**. This function identifies the coordinates of the subdomain with process rank “id”. The identification of the id and subdomain coordinates are given by

$$\text{id} = \text{icoords}[0] + \text{icoords}[1]*\text{gmax}[0] + \text{icoords}[2]*\text{gmax}[0]*\text{gmax}[1].$$

This defines the natural lexicographical assignment of id numbers to the subdomains, as illustrated for the above 4x3 partition.

```

typedef struct {
    int   gmax[3]; /* # of subdomains in each dir */
    int   nn;      /* total number of nodes      */
    int   dim;
} PP_GRID;

```

```

void find_Cartesian_coordinates(
    int   id,
    PP_GRID *pp_grid,
    int   *icoords)
{
    int   dim = pp_grid->dim;
    int   d, G;

    for (d = 0; d < dim; d++)
    {
        G = pp_grid->gmax[d];
        icoords[d] = id % G;
        id = (id - icoords[d])/G;
    }
} /*end find_Cartesian_coordinates*/

```

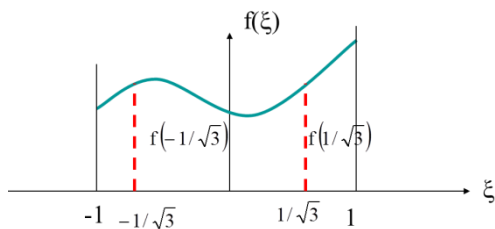
## 2. Compute subdomain boundaries

For process with rank “id” and coordinate  $(i, j)$  on the process grid,

$$x_{i,l} = 2.0 + i * (5.0 - 2.0)/\sqrt{P}$$

## 3. Quadrature rules

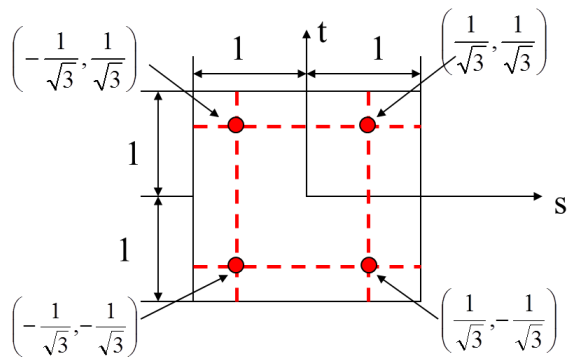
### 1D 2-point Gaussian quadrature rule



$$\int_{-1}^1 f(\xi) d\xi \approx f\left(\frac{1}{\sqrt{3}}\right) + f\left(-\frac{1}{\sqrt{3}}\right)$$

### Gaussian Quadrature Rule in 2D

$$I = \int_{-1}^1 \int_{-1}^1 f(s, t) ds dt$$



$$\begin{aligned}
 I &= \int_{-1}^1 \int_{-1}^1 f(s, t) ds dt \\
 &\approx \int_{-1}^1 \left( \sum_{j=1}^M w_j f(s, t_j) \right) ds \\
 &\approx \sum_{i=1}^M \sum_{j=1}^M w_i w_j f(s_i t_j) \\
 &= \sum_{i=1}^M \sum_{j=1}^M w_{ij} f(s_i t_j)
 \end{aligned}$$

Where  $w_{ij} = w_i w_j$ .

## Problem 2. Parallel Explicit Finite Difference Scheme for Solving 1D

### Heat Equation for Graduate Students.

Consider to solve  $\begin{cases} u_t(x, t) = u_{xx}(x, t), & 0 \leq x \leq 2\pi, t > 0 \\ u(x, 0) = \sin(x) & 0 \leq x \leq 2\pi \end{cases}$  with periodic boundary condition by the explicit finite difference scheme. Compute the solution for  $t = 2.0$ .

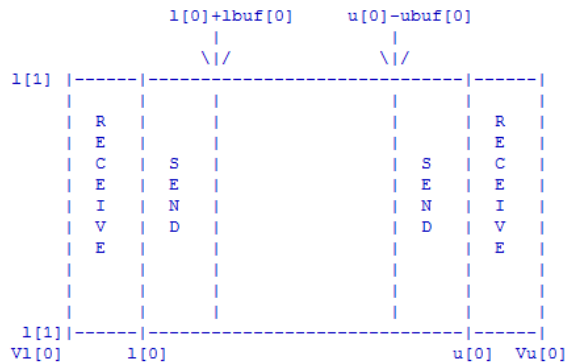
The exact solution is given by  $u(x, t) = e^{-t} \sin(x)$ .

Assume we use  $M + 1$  grid points. The grid space then is  $\Delta x = \frac{2\pi}{M}$ . The grid points are  $x_k = k\Delta x$ ,  $k = 0, \dots, M$ . Let  $\Delta t$  be the time step size. For stability, we should satisfy  $\frac{\Delta t}{\Delta x^2} \leq 0.5$ .

Let  $v_k^n \approx u(k\Delta x, n\Delta t)$  be the approximate solution. The explicit scheme is

$$v_k^{n+1} = v_k^n + \frac{\Delta t}{\Delta x^2} (v_{k+1}^n - 2v_k^n + v_{k-1}^n) \text{ for } k = 0, \dots, M.$$

Implement a parallel version of the scheme to solve the above problem by arbitrary number of grid points  $M + 1$  using  $P$  processors. Assume  $M + 1 \gg P$ . Use non-blocking send and blocking receive for message passing. Use  $M = 1000, 2000, 4000, 8000$  respectively to do the mesh refinement study. Compute  $L_{2,\Delta x}$  error with respect to the mesh refinement. Do each of these calculations with 4, 8, 16 processors respectively. Make a table to list the wall clock time space on computation and communication respectively. Hint: the code to compute work load assignment is `work_division.c`.



Let  $[l[0], u[0]]$  be a subdomain assigned to a process. For convenience of computation, a virtual domain is defined to hold the ghost points for updating solutions defined on grid points within  $[l[0], u[0]]$ . This virtual domain is defined as  $[vl[0], vu[0]]$ . Here  $vl[0] = l[0] - N * \Delta x$ , and  $vu[0] = u[0] + N * \Delta x$ .  $N$  is the number of ghost points.

The communication routine looks like the following:

```

void scatter_states(double *soln)
{
    int    myid, side;
    int    me[3];

    MPI_Comm_rank(MPI_COMM,&myid);
    find_Cartesian_coordinates(myid,pp_grid,me);

    for (side = 0; side < 2; ++side)
    {
        MPI_Barrier(MPI_COMM);
        pp_send_interior_states(me, side,soln);
        pp_receive_interior_states(me ,(side+1)%2,soln);
    }
}

void pp_send_interior_states(
    int    *me,
    int    side,
    double *soln)
{
    int    myid, him[3], dst_id;

    MPI_Comm_rank(MPI_COMM,&myid);
    dst_id = neighbor_id(him,me,0,side,pp_grid);

    /* Next collect soln points to be sent and call MPI_Isend() to send the data
       to the process with rank dst_id */
}

void pp_receive_interior_states(
    int    *me,
    int    side,
    double *soln)
{
    int    myid, him[3], src_id;
    MPI_Comm_rank(MPI_COMM,&myid);
    src_id = neighbor_id(him,me,0,side,pp_grid);

    /* Next call MPI_Recv() to receive the data
       from the process with rank src_id */
}

```

```

int neighbor_id(
    int      *him,
    int      *me,
    int      dir,
    int      side,
    PP_GRID  *pp_grid)
{
    int      *G = pp_grid->gmax;
    int      i, dim = pp_grid->dim;

    for (i = 0; i < dim; i++)
        him[i] = me[i];

    him[dir] = (me[dir] + 2*side - 1);
    if (him[dir] < 0)
        him[dir] = G[dir] - 1;
    if (him[dir] >= G[dir])
        him[dir] = 0;

    return domain_id(him,G,dim);
}      /*end neighbor_id*/

int domain_id(
    int      *icoords,
    int      *G,
    int      dim)
{
    int      tmpid;
    int      i;

    tmpid = icoords[dim-1];
    for (i = dim-2; i >= 0; i--)
        tmpid = icoords[i] + G[i]*tmpid;
    return tmpid;
}      /*end domain_id*/

```

**Hand-In.** Turn in the hardcopy of all your source code, and the report which contains results and algorithmic notes on both computation and communication. Email the source code.