Homework 3

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1 Algorithm

Given the problem in the assignment we can show that although the initial matrix is symmetrical, after the first iteration we will lose the symmetry so we have to calculate all the values in the matrix.

Given more processors we can split the work to be done. Since we want the divisions to be as cohesive as possible to minimize the communication, we can opt for two ways to do the domain decomposition. Both involve Halo Swapping as it renders the most effective strategy in this case since every element depends on the ones above, below and to the right.

First the serial time for this algorithm will be SerTime = $\Theta(n^2)$ since we will do a constant amount of iterations during which it will update the n^2 elements of the matrix. Now, an ideal parallel algorithm without communication we would get ParTime = $\Theta(n^2/p)$, however any real algorithm will have to communicate so the SpeedUp won't be p.

If we do the division per processors in one dimension, the Parallel time will be as follows: ParTime $=\Theta(n^2/p)+\Theta(n/p)+\Theta(n)=\Theta(n^2/p)+\Theta(n)$. However, if we do the partition in both dimensions (we are allowed to do it since p is given to be a perfect square) the parallel time will be ParTime = $\Theta(n^2/p) + \Theta(n/\sqrt{p})$ which is clearly better than the naive approach shown before.

$\mathbf{2}$ Analysis

The chosen domain decomposition renders the processor communication in a 2D grid. To better the communication involved there should be communication between the first row of processors and the last row, to make use of the modular properties of the algorithm. Therefore, the sought topology is cylindrical.

Given the partition shown we will have for each processor p with coordinates in the processor grid (x, y). The size of the matrix will be $h \times w$ where h and w are:

$$h = \begin{cases} \lceil \frac{n}{\sqrt{p}} \rceil & \text{if } x < \sqrt{p} - 1\\ n \mod \lceil \frac{n}{\sqrt{p}} \rceil & \text{if } x = \sqrt{p} - 1 \end{cases}$$
 (2.1)

$$h = \begin{cases} \lceil \frac{n}{\sqrt{p}} \rceil & \text{if } x < \sqrt{p} - 1\\ n \mod \lceil \frac{n}{\sqrt{p}} \rceil & \text{if } x = \sqrt{p} - 1 \end{cases}$$

$$w = \begin{cases} \lceil \frac{n}{\sqrt{p}} \rceil & \text{if } y < \sqrt{p} - 1\\ n \mod \lceil \frac{n}{\sqrt{p}} \rceil & \text{if } y = \sqrt{p} - 1 \end{cases}$$

$$(2.1)$$

Apart from this $h \times w$ matrix, each processor will have to store three halos corresponding to the following values

• Upper Halo of size $1 \times w$, corresponding to the lower row of the processor above (note that $(\sqrt{p}-1,y)$) is above (0, y) in a cylindrical topology). This values are needed when $i_{local} = 0$ and i'' looks for the value above.

- Lower Halo of size $1 \times w$, corresponding to the upper row of the processor below. This values will be needed hen $i_{\text{local}} = \lceil \frac{n}{\sqrt{p}} \rceil 1$ and i' looks for the value below.
- Right Halo of size $h \times 1$, corresponding to the leftmost column of the processor to the right. This values will be needed when $j_{\text{local}} = \lceil \frac{n}{\sqrt{p}} \rceil 1$ and j' will ask for the value to the right. The processors with coordinates $(x, \sqrt{p} 1)$ will not use this halos since the values of the matrix with coordinates (i, n 1) do not update.

The communication will be composed of three shifts.

- 1. Upward Modular Shift Each processor (x, y) sends its upper row to the processor above $(x 1 \mod \sqrt{p}, y)$ and receives from the one below $(x + 1 \mod \sqrt{p}, y)$ storing the values in the Lower Halo. Since this is a cylindrical topology this operation involves all the processors sending and receiving at the same time.
- 2. **Downward Modular Shift** Similarly, this is the reverse operation. Each processor (x,y) sends its lower row to the processor below $(x+1 \mod \sqrt{p}, y)$ and receives from the one above $(x-1 \mod \sqrt{p}, y)$ storing the values in the Upper Halo.
- 3. Leftward Shift This involves every processor (x, y) sending its leftmost column to the one on their left (x, y 1) side and receiving from the one on their right (x, y + 1) and storing th values in their Right Halo. Note than this an open loop and the processors (x, 0) won't send any information and the processors $(x, \sqrt{p} 1)$ won't receive any information either.

From this we can gather that the communication time is $O(\max(h, w))$ since we are swapping three halos per processor every time the processors communicate. Since $\max(h, w) = \lceil \frac{n}{\sqrt{p}} \rceil$ the time involved in the communication will be $O(\frac{n}{\sqrt{p}})$.

Since the update involved modifying the whole matrix it will take time $O(h \times w)$ per processor. Since $h, w <= \lceil \frac{n}{\sqrt{p}} \rceil$ we can express it as $O(h \times w) = O(n^2/p)$.

Finally the whole parallel time will be

$$ParTime = \Theta(n^2/p) + \Theta(n/\sqrt{p})$$
(2.3)

The SpeedUp will be

SpeedUp =
$$\frac{\text{SerTime}}{\text{ParTime}} = \frac{\Theta(n^2)}{\Theta(n^2/p) + \Theta(n/\sqrt{p})}$$
 (2.4)

SpeedUp
$$\to p$$
 as $n \to \infty$ (2.5)

And the efficiency:

$$\text{Efficiency} = \frac{\text{SerTime}}{p \cdot \text{ParTime}} = \frac{\Theta(n^2)}{\Theta(n^2) + \Theta(n\sqrt{p})} \tag{2.6}$$

Efficiency
$$\to 1$$
 for $p = O(n^2)$ (2.7)

3 Implementation

To implement the algorithm we have used the partition described above and used a cylindrical virtual topology where the row processors are wrapped around but the column processors are not. Furthermore it has been used MPI_SENDRECV for the communication, using a derived datatype for the column communication.

The particular aspects of the whole implementation can be seen in Code A.1

4 Results

The results from the timings are shown in the Table 4.1. The Verification values for the different matrix sizes are shown in the Table 4.2.

n	\mathbf{p}	average	min	max
1000	1	5.71552	5.34630	5.83059
2000	1	22.82041	21.18143	23.31813
1000	4	1.46049	1.41766	1.49389
2000	4	6.04183	5.67908	7.35377
1000	16	0.46770	0.38920	0.54539
2000	16	1.77513	1.50135	2.04136
1000	36	0.30457	0.22946	0.37673
2000	36	1.01406	0.68300	1.30123

Table 4.1: Results in seconds

\mathbf{n}	sum	min
1000	4283810.871966	-892.718330
2000	17631642 148465	-1800 724437

Table 4.2: Verification Values

A Code

```
#include <mpi.h>
1
    #include <ctime>
2
    #include <stdio.h>
3
    #include <stdlib.h>
4
    #include <math.h>
6
    #define ITERATIONS 10
7
    #define LEFT_COLUMN A[1][0]
9
    #define RIGHT_HALO A[1][width]
10
    #define UPPER_ROW A[1][0]
11
    #define UPPER_HALO A[0][0]
12
    #define LOWER_ROW A[height][0]
13
    #define LOWER_HALO A[height+1][0]
14
15
    double** new_matrix(int height, int width);
16
17
    void delete_matrix(double** matrix);
18
    void initialize_matrix(double** A, int size, int height, int width, int* coords);
19
    void transform_matrix(double** A, double** B, int height, int width, int* coords);
20
    void print_matrix(double** A, int height, int width);
21
22
    char* gettime(char *buffer);
23
24
    int main(int argc, char *argv[])
25
26
        ///////// VARIABLES /////////////
27
        char *s;
28
29
        double **A, **B;
30
        int n, m, p, sqrt_p, height, width;
31
        int rank, coords[2];
32
33
        // int output = 0;
34
35
        MPI_Comm matrix_comm;
36
        static int dims [2];
37
        int periods [2] = { 1, 0 };
        int reorder = 1;
39
        int p_up, p_down, p_right, p_left;
40
41
        double local_sum=0.0, local_min;
42
        double starttime, endtime;
43
44
        double global_sum=0.0, global_min;
45
46
        char buffer[80];
47
48
        ////////////// INIT //////////////
49
50
        MPI_Init (&argc, &argv);
51
        MPI_Comm_size(MPI_COMM_WORLD,&p);
52
        sqrt_p = (int)sqrt(p);
53
        dims[0] = sqrt_p;
54
        dims[1] = sqrt_p;
55
```

```
56
         MPI_Cart_create (MPI_COMM_WORLD, 2, dims, periods, reorder, &matrix_comm);
57
58
         MPI_Comm_rank(matrix_comm, &rank);
59
         MPI_Cart_coords (matrix_comm, rank, 2, coords);
60
61
         MPI_Datatype column;
62
         // The dimensions are reversed because MPI_Cart creates the matrix by columns and
63
         we want it by rows
         MPI_Cart_shift(matrix_comm, 1, -1, &p_right, &p_left);
64
         MPI_Cart_shift(matrix_comm, 0, -1, &p_down, &p_up);
65
66
         //////// PROGRAM ////////////
67
68
         n = strtol(argv[1], &s, 10);
69
70
         m = ceil(float(n) / sqrt_p);
71
         height = coords[0] != sqrt_p-1 ? m : n - m*(sqrt_p-1);
72
         width = coords[1] != sqrt_p-1 ? m : n - m*(sqrt_p-1);
73
74
         MPI_Type_vector (height, 1, width+1, MPI_DOUBLE, &column);
75
         MPI_Type_commit (&column);
76
77
         A = \text{new\_matrix}(\text{height+2,width+1});
78
         B = new_matrix(height+2, width+1);
79
80
         initialize_matrix(A, m, height, width, coords);
81
82
         MPI_Barrier (matrix_comm);
83
         if(rank == 0){
             starttime = MPI_Wtime();
85
86
87
         for (int i = 0; i < ITERATIONS; ++i)</pre>
88
89
             //// HALO SWAPPING ////
90
91
             MPI_Sendrecv(&( UPPER_ROW ), width, MPI_DOUBLE, p_up, 0, &( LOWER_HALO ),
92
         width, MPI_DOUBLE, p_down, 0, matrix_comm, MPI_STATUS_IGNORE);
93
             MPI_Sendrecv(&( LOWER_ROW ), width, MPI_DOUBLE, p_down, 0, &( UPPER_HALO ),
94
         width, MPI_DOUBLE, p_up, 0, matrix_comm, MPI_STATUS_IGNORE);
             // Left
95
             MPI_Sendrecv(&( LEFT_COLUMN ), 1, column, p_left, 0, &( RIGHT_HALO ), 1,
96
         column, p_right, 0, matrix_comm, MPI_STATUS_IGNORE);
97
             if(coords[1] != sqrt_p-1) {
98
                  transform_matrix(A,B,height,width,coords);
99
             }else{
100
                  transform_matrix(A, B, height, width-1, coords);
101
102
103
104
105
         ///////// VERIFICATION /////////////
106
107
         local\_min = A[1+0][0];
108
         for (int i = 0; i < height; i++)</pre>
109
110
```

```
for (int j = 0; j < width; j++)</pre>
111
112
                  local_sum += fabs( A[1+i][j]);
113
                  local_min = A[1+i][j] < local_min ? A[1+i][j] : local_min;</pre>
114
              }
115
116
117
         MPI_Reduce(&local_sum, &global_sum, 1, MPI_DOUBLE, MPI_SUM, 0, matrix_comm);
118
         MPI_Reduce(&local_min, &global_min, 1, MPI_DOUBLE, MPI_MIN, 0, matrix_comm);
119
120
         ///////// FINALIZE /////////////
121
         if(rank == 0) {
122
              endtime = MPI_Wtime();
123
             FILE *f = fopen("Results.txt", "a");
124
             if(f != NULL)
125
126
                  fprintf(f, "%s\n", gettime(buffer));
127
                  fprintf(f, "%dx%d matrix\n", n, n);
128
                  fprintf(f,"%d processors\n",p);
                  fprintf(f,"Sum : %f\n", global_sum );
130
                  fprintf(f,"Min : %f\n", global_min );
131
                  fprintf(f, "Time : %f seconds\n", endtime-starttime);
132
                  fprintf(f,"\n=======\n");
133
              }
134
135
136
         delete_matrix(A);
137
         delete_matrix(B);
138
139
         MPI_Finalize();
141
         return 0;
142
143
     }
144
145
     double** new_matrix(int height, int width)
146
147
         double** matrix;
148
         matrix = new double*[height];
149
         matrix[0] = new double[height * width];
150
         for (int i = 1; i < height; i++)</pre>
151
             matrix[i] = matrix[i-1] + width;
152
         return matrix;
153
     }
154
155
     void delete_matrix(double** matrix)
156
157
         delete[] matrix[0];
158
         delete[] matrix;
159
160
161
162
     void initialize_matrix(double** A, int size, int height, int width, int* coords)
163
164
         int Origin_x = coords[0]*size;
165
         int Origin_y = coords[1]*size;
166
         for (int i_G = Origin_x, i = 0; i_G < Origin_x + height; i_G++, i++)
167
168
             for (int j_G = Origin_y, j = 0; j_G < Origin_y + width; j_G++, j++)
169
```

```
170
                   if(coords[0] == coords[1] && i == j){
171
                       A[1+i][j] = i_G * sin(sqrt(i_G));
172
                   }else{
173
                       A[1+i][j] = pow(i_G+j_G,1.1);
174
175
              }
176
177
178
179
          return;
180
     }
181
     void transform_matrix(double** A, double** B, int height, int width, int* coords)
182
183
          double x;
184
          for (int i = 0; i < height; i++)</pre>
185
186
              for (int j = 0; j < width; j++)</pre>
187
188
                   if( ! (coords[0] == coords[1] && i == j) )
189
                        x = 0.0;
191
                        for (int k = 1; k \le 10; k++) {
192
                            x += pow(fabs(0.5 + A[1+i+1][j]), 1.0/double(k));
193
                            x -= pow( fabs( A[1+i-1][j] ) , 1.0/double(k+1)) * pow( fabs( A[1+i-1][j]
194
         i][j+1]), 1.0/double(k+2));
                        }
195
                        x = x < 10.0 ? x : 10.0;
196
                        x = x > -10.0 ? x : -10.0;
197
                       B[1+i][j] = x;
198
                   }else{
199
                       B[1+i][j] = A[1+i][j];
200
201
              }
202
203
          for (int i = 0; i < height; i++)</pre>
204
205
              for (int j = 0; j < width; j++)</pre>
206
207
                   A[1+i][j] = B[1+i][j];
210
211
          return;
^{212}
213
     void print_matrix(double** A, int height, int width)
214
215
          for (int i = 0; i < height+2; i++)</pre>
216
^{217}
              for (int j = 0; j < width+1; j++)</pre>
218
220
                   printf("%f ", A[i][j]);
221
              printf("\n");
222
223
     }
224
225
     char* gettime(char* buffer)
226
227
```

```
time_t rawtime;
228
         struct tm * timeinfo;
229
230
         time (&rawtime);
231
         timeinfo = localtime(&rawtime);
232
233
         strftime(buffer, 80, "%Y-%m-%d %I:%M:%S", timeinfo);
234
         std::string str(buffer);
235
         return buffer;
237
```

Listing A.1: C++ code for the matrix algorithm