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High Performance Computing Lab

Assignment - 8

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MPI Programming

1. Study and implement 2D Convolution using MPI. Use different number of processes and analyze the performance.

Code:

```
#include<assert.h>
#include<mpi.h>
#include<stdio.h>
#include<stdlib.h>
#include<string.h>
#include<sys/time.h>

#defineDEFAULT_ITERATIONS1

intconv_column(int*,int,int,int*,int*,int);
intconv(int*,int,int,int*,int);
int*check(int*,int,int,int*,int);
intconv_column(int*sub_grid,inti,intnrows,int
DIM,int*kernel,intkernel_dim) {
intcounter=0;
```

```
intnum pads= (kernel dim-1) /2;
for (intj=1; j< (num pads+1); j++) {
+j*kernel dim];
*kernel[(((kernel dim-1) * (kernel dim+1)) /2) -
j*kernel diml;
counter=counter+sub grid[i] *kernel[(((kernel dim-1)
* (kernel dim+1)) /2)];
returncounter;
intconv(int*sub grid,inti,intnrows,int
DIM, int*kernel, intkernel dim) {
intcounter=0;
intnum pads= (kernel dim-1) /2;
// convolve middle column
counter=counter+conv column(sub grid, i, nrows, DIM,
kernel, kernel dim);
// convolve left and right columns
for (intj=1; j< (num pads+1); j++) {</pre>
// get last element of current row
intend= (((i/DIM) +1) *DIM) -1;
if (i+j-end \le 0) { // if column is valid
counter=counter+conv column(sub grid, i+j, nrows,
```

```
intfirst= (i/DIM) *DIM;
if (i-j-first>=0) {
DIM, kernel, kernel dim);
int*check(int*sub grid,intnrows,int
intval;
intnum pads= (kernel dim-1) /2;
int*new grid=calloc(DIM*nrows, sizeof(int));
for (inti= (num pads*DIM); i< (DIM*</pre>
kernel dim);
new grid[i- (num pads*DIM)] =val;
intmain(intargc, char**argv) {
// MPI Standard variable
intID, j;
intiters=0;
intnum iterations;
intDIM;
```

```
intGRID WIDTH;
intKERNEL DIM;
intKERNEL SIZE;
num iterations=DEFAULT ITERATIONS;
if (argc >= 3) {
DIM=atoi(argv[1]);
GRID WIDTH=DIM*DIM;
KERNEL DIM=atoi(argv[2]);
KERNEL SIZE=KERNEL DIM*KERNEL DIM;
if (argc==4) {
num iterations=atoi(argv[3]);
    } else {
printf("Invalid command line arguments");
MPI Finalize();
exit(-1);
intmain grid[GRID WIDTH];
memset(main grid, 0, GRID WIDTH*sizeof(int));
for (inti=0; i<GRID WIDTH; i++) {</pre>
main grid[i] =1;
intnum pads= (KERNEL DIM-1) /2;
intkernel[KERNEL SIZE];
memset(kernel, 0, KERNEL SIZE*sizeof(int));
for (inti=0; i<KERNEL SIZE; i++) {</pre>
kernel[i] = 1;
```

```
MPI Statusstatus;
// MPI Setup
MPI Init (NULL, NULL);
MPI Comm size (MPI COMM WORLD, &num procs); // Set the
num procs
MPI Comm rank (MPI COMM WORLD, &ID);
doublestart time=MPI Wtime();
assert(DIM%num procs==0);
intupper[DIM*num pads];
intlower[DIM*num pads];
int*pad row upper;
int*pad row lower;
intstart= (DIM/num procs) *ID;
intend= (DIM/num procs) -1+start;
intnrows=end+1-start;
intnext= (ID+1) %num procs;
intprev=ID!=0?ID-1:num procs-1;
for (iters=0; iters<num iterations; iters++) {</pre>
memcpy(lower, &main grid[DIM* (end-num pads+1)],
sizeof(int) *DIM*num pads);
```

```
pad row lower=malloc(sizeof(int) *DIM*num pads);
memcpy(upper, &main grid[DIM*start], sizeof(int)
*DIM*num pads);
pad row upper=malloc(sizeof(int) *DIM*num pads);
if (num procs>1) {
if (ID\%2 == 1) {
MPI Recv(pad row lower, DIM*num pads, MPI INT, next,
1, MPI COMM WORLD, &status);
MPI Recv(pad row upper, DIM*num pads, MPI INT, prev,
1, MPI COMM WORLD, &status);
            } else {
MPI Send(upper, DIM*num pads, MPI INT, prev, 1,
MPI COMM WORLD);
MPI Send(lower, DIM*num pads, MPI INT, next, 1,
MPI COMM WORLD);
if (ID%2==1) {
MPI Send(upper, DIM*num pads, MPI INT, prev, 0,
MPI COMM WORLD);
MPI Send(lower, DIM*num pads, MPI INT, next, 0,
MPI COMM WORLD);
            } else {
MPI Recv(pad row lower, DIM*num pads, MPI INT, next,
0, MPI COMM WORLD, &status);
MPI Recv(pad row upper, DIM*num pads, MPI INT, prev,
0, MPI COMM WORLD, &status);
        } else {
pad row lower=upper;
pad row upper=lower;
```

```
intsub grid[DIM* (nrows+ (2*num pads))];
if (ID==0) {
memset(pad row upper, 0, DIM*sizeof(int) *num pads);
if (ID== (num procs-1)) {
memset(pad row lower, 0, DIM*sizeof(int) *num pads);
memcpy(sub grid, pad row upper, sizeof(int)
memcpy(&sub grid[DIM*num pads],
&main grid[DIM*start], sizeof(int) *DIM*nrows);
memcpy(&sub grid[DIM* (nrows+num pads)],
pad row lower, sizeof(int) *DIM*num pads);
int*changed subgrid=check(sub grid, nrows, DIM,
if (ID!=0) {
MPI Send(changed subgrid, nrows*DIM, MPI INT, 0, 11,
MPI COMM WORLD);
MPI Recv(&main grid[0], DIM*DIM, MPI INT, 0, 10,
MPI COMM WORLD, &status);
        } else {
for (inti=0; i<nrows*DIM; i++) {</pre>
main grid[i] =changed subgrid[i];
for (intk=1; k<num procs; k++) {</pre>
MPI Recv(&main grid[DIM* (DIM/num procs) *k],
nrows*DIM, MPI INT, k, 11, MPI COMM WORLD, &status);
```

```
for (inti=1; i<num procs; i++) {</pre>
MPI Send (main grid, DIM*DIM, MPI INT, i, 10,
MPI COMM WORLD);
// Output the updated grid state
if (ID==0) {
doubleend=MPI Wtime();
printf("Matrix DIM: %d\n", DIM);
printf("Kernel DIM: %d", KERNEL DIM);
printf("\nConvolution Output: \n");
for (j=0; j<GRID WIDTH; j++) {</pre>
if (j%DIM==0) {
printf("\n");
printf("%d ", main_grid[j]);
printf("\n");
printf("Execution Time: %f\n", end-start time);
if (num procs>=2) {
free(pad row upper);
free(pad row lower);
MPI Finalize();
```

Output:

```
PS D:\Assignments> mpiexec -n 2 .\Q1.exe 10 8
Matrix DIM: 10
Kernel DIM: 8
Convolution Output:
16
    20
        24
            28
                28
                    28
                        28
                            24
                                20
                                    16
20
    25
        30
            35
                35
                    35
                        35
                            30
                                25
                                    20
24
    30
        36
            42
                42
                    42
                        42
                            36
                                30
                                    24
28
    35
        42
                            42
                                35
                                    28
            49
                49
                    49
                        49
28
    35
        42
            49
                49
                    49
                        49
                            42
                                35
                                    28
28
    35
                                35
                                    28
        42
            49
                49
                    49
                        49
                            42
                    49
                            42
28
    35
        42
            49
                49
                        49
                                35
                                    28
24
    30
        36
            42
               42 42
                        42
                            36
                                30
                                    24
20
    25
               35 35
                                    20
       30
            35
                        35
                            30
                                25
16
    20 24
            28
                28
                    28
                            24
                        28
                                20
                                    16
Execution Time: 0.000978
```

For 500*500:

```
PS D:\Assignments> mpiexec -n 1 .\Q1.exe 500 8
Execution Time: 0.033009
PS D:\Assignments> mpiexec -n 2 .\Q1.exe 500 8
Execution Time: 0.017275
PS D:\Assignments> mpiexec -n 4 .\Q1.exe 500 8
Execution Time: 0.014928
```

PS D:\Assignments> mpiexec -n 20 .\Q1.exe 500 8 Execution Time: 0.023533

р	Time
1	0.033

2	0.0173
4	0.0149
20	0.0235

Conclusion:

Firstly with increasing p, execution time decreases then after saturation point it increases again.

2. Implement dot product using MPI. Use different number of processes and analyze the performance.

Code:

#include<math.h>

```
#include<mpi.h>
#include<stdio.h>
#include<stdlib.h>
#include<time.h>
#include<unistd.h>

#defineNELMS100000
#defineMASTER0
#defineMAXPROCS16

intdot_product();
voidinit_lst();
voidprint_lst();
intmain() {
inti, n, vector_x[NELMS], vector_y[NELMS];
```

```
int prod, sidx, eidx, size;
intpid, nprocs, rank;
doublestime, etime;
MPI Status status;
MPI Comm world;
   n = 100000;
if (n > NELMS) {
printf("n=%d> N=%d n", n, NELMS);
exit(1);
MPI Init (NULL, NULL);
    world = MPI COMM WORLD;
MPI Comm size(MPI COMM WORLD, &nprocs);
MPI Comm rank (MPI COMM WORLD, &pid);
int portion = n /nprocs;
sidx=pid* portion;
eidx=sidx+ portion;
init lst(vector x, n);
init lst(vector y, n);
inttmp prod[nprocs];
for (i=0; i< nprocs; i++)
tmp prod[i] =0;
stime=MPI Wtime();
if (pid== MASTER) {
        prod =dot product(sidx, eidx, vector x, vector y,
n);
```

```
for (i=1; i< nprocs; i++)
MPI Recv(&tmp prod[i-1], 1, MPI INT, i, 123,
MPI COMM WORLD, &status);
   } else {
        prod =dot product(sidx, eidx, vector x, vector y,
n);
MPI Send(&prod, 1, MPI INT, MASTER, 123, MPI COMM WORLD);
if (pid== MASTER) {
for (i=0; i< nprocs; i++)
           prod +=tmp prod[i];
etime=MPI Wtime();
if (pid== MASTER) {
// print lst(vector x,n);
// print lst(vector y,n);
printf("pid=%d: final prod=%d\n", pid, prod);
printf("pid=%d: elapsed=%f\n", pid, etime-stime);
MPI Finalize();
intdot product(int s, int e, int x[], int y[], int n) {
inti, prod =0;
for (i= s; i< e; i++)
        prod = prod + x[i] * y[i];
return prod;
```

```
voidinit_lst(int*l,int n) {
  inti;
  for (i=0; i< n; i++)
  *l++=i;
}
voidprint_lst(int l[],int n) {
  inti;

for (i=0; i< n; i++) {
  printf("%d ", l[i]);
    }
  printf("\n");
}</pre>
```

Output:

```
PS D:\Assignments> mpiexec -n 2 .\Q2.exe
pid=0: final prod=216474736
pid=0: elapsed=0.000839
PS D:\Assignments> mpiexec -n 4 .\Q2.exe
pid=0: final prod=216474736
pid=0: elapsed=0.001077
PS D:\Assignments> mpiexec -n 8 .\Q2.exe
pid=0: final prod=216474736
pid=0: elapsed=0.002098
PS D:\Assignments> mpiexec -n 16 .\Q2.exe
pid=0: final prod=216474736
pid=0: final prod=216474736
pid=0: final prod=216474736
pid=0: elapsed=0.003044
PS D:\Assignments>
```

р	time
1	0.00025
2	0.00083
4	0.00072
8	0.00210
16	0.00304

Conclusion: The increase or decrease in time is not regular it is increasing and decreasing.

3. Implement Prefix sum using MPI. Use different number of processes and analyze the performance.

Code:

```
MPI Comm rank (MPI COMM WORLD, &my rank);
MPI Comm size(MPI COMM WORLD, &p);
intprefix arr[p];
if (my rank==0) {
inti;
for (i=0; i<p; ++i) {
prefix arr[i] =i+1;
doublestart=MPI Wtime();
// all call scatter
MPI Scatter(prefix arr, 1, MPI_INT, &value, 1,
MPI INT, 0, MPI COMM WORLD);
receve a value from someone and add to ours
       otherwise, we send to the chosen one
inti;
intlogn=log2(p);
for (i=0; i<=logn; i++) {
```

```
intlower bound=pow(2, i);
intupper bound=p-lower bound;
if (upper bound<lower bound) {</pre>
upper bound=lower bound;
if (my rank<lower bound) {</pre>
intsend= (int) (my rank+pow(2, i));
if (send>=p)
continue;
printf("%d sending to %dn", my rank,
(int) (my rank+pow(2, i)));
MPI Send(&value, 1, MPI INT, (int)(my rank+pow(2,
i)), 0, MPI COMM WORLD);
        } elseif (my rank>=upper bound) {
intrecv= (int) (my rank-pow(2, i));
if (recv>=p)
continue;
intrecv value;
printf("%dreceving..\n", my rank);
MPI Recv(&recv value, 1, MPI INT, (my rank-pow(2,
i)), 0, MPI COMM WORLD, &status);
value+=recv value;
        } else {
intsend= (int) (my rank+pow(2, i));
intrecv= (int) (my rank-pow(2, i));
if (send>=p||recv>=p)
continue;
```

```
printf("%d sending to %d\n", my rank,
(int)(my rank+pow(2, i)));
MPI Send(&value, 1, MPI INT, (int)(my rank+pow(2,
i)), 0, MPI COMM WORLD);
printf("%dreceving..\n", my rank);
intrecv value;
MPI Statusstatus;
MPI Recv(&recv value, 1, MPI INT, (my rank-pow(2,
i)), 0, MPI COMM WORLD, &status);
value+=recv value;
// after algorithm, each processor hols its own
prefix sum
// we gather at rank
intgather[p];
MPI Gather(&value, 1, MPI INT, gather, 1, MPI INT, 0,
MPI COMM WORLD);
if (my rank==0) {
doubleend=MPI Wtime();
printf("Execution Time: %f\n", end-start);
   /* shut down MPI */
MPI Finalize();
return0;
```

Output:

```
PS D:\Assignments> mpiexec -n 4 .\Q3.exe
3 receving...
3 receving...
1 sending to 2
1 receving..
1 sending to 3
2 sending to 3
2 receving...
2 receving...
0 sending to 1
0 sending to 2
Execution Time: 0.001033
PS D:\Assignments> mpiexec -n 8 .\Q3.exe
1 sending to 2
1 receving...
1 sending to 3
1 sending to 5
7 receving...
7 receving...
7 receving...
2 sending to 3
2 receving..
2 sending to 4
2 receving...
2 sending to 6
0 sending to 1
0 sending to 2
0 sending to 4
Execution Time: 0.001867
5 sending to 6
5 receving...
5 sending to 7
5 receving...
5 receving...
3 sending to 4
3 receving...
3 sending to 5
3 receving..
3 sending to 7
```

6 sending to 7

р	time
2	0.0006
4	0.0010
8	0.0018
16	0.0033

Conclusion: With increasing p value time is gradually increasing.