

# High Performance Computing

COM403P

## **Week-6**

Matrix Multiplication

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## Objective

Matrix Multiplication for given  $n \times n$  double precision floating point numbers.

## Serial Code

```
#include <stdio.h>
#include <stdlib.h>
#include <omp.h>
#include <time.h>

#define SIZE 100

int main()
{
    double a[SIZE][SIZE], b[SIZE][SIZE], c[SIZE][SIZE], rand_a, rand_b;
    double start, end, exec;

    start = omp_get_wtime();

    for (int i = 0; i < SIZE; i++)
    {
        for(int j = 0; j < SIZE; j++)
        {
            rand_a = rand();
            rand_b = rand();

            a[i][j] = i*rand_a;
            b[i][j] = i*rand_b;

            // for(int k = 0; k < 100000; k++)
            c[i][j] = a[i][j] * b[i][j];
        }
    }

    end = omp_get_wtime();

    exec = end - start;

    printf("Serial Exec time - %f\n", exec);

    return 0;
}
```

## Parallel Code

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

#define MASTER 0
#define FROM_MASTER 1
#define FROM_WORKER 2

#define SIZE 100

int main()
{
    MPI_Init(NULL, NULL);

    long double a[SIZE][SIZE], b[SIZE][SIZE], c[SIZE][SIZE], rand_a, rand_b;
    double start, end, exec;
    int i,j,k;

    int avgrow, extra;
    int offset, mtype;
    int dest, rows;
    int source;
    MPI_Status status;

    start = MPI_Wtime();

    int taskid;
    MPI_Comm_rank(MPI_COMM_WORLD,&taskid);

    int numtasks;
    MPI_Comm_size(MPI_COMM_WORLD,&numtasks);

    int workers = numtasks -1;

    if(taskid == MASTER)
    {
        for (i = 0; i < SIZE; i++)
        {
```

```
        for(j = 0; j < SIZE; j++)
        {
            a[i][j] = (i+j)*1.22;
            b[i][j] = (i+j)*1.22;
        }
    }

//-----
// Master sending Data to Worker tasks

avgrow = SIZE/workers;
extra = SIZE%workers;

offset = 0;
mtype = FROM_MASTER;

for(dest = 1; dest <= workers; dest++)
{
    rows = (dest <= extra)?avgrow+1:avgrow;

    // printf("Sending %d rows to task-%d offset=%d\n",rows,dest,offset);

    MPI_Send(&offset, 1, MPI_INT, dest, mtype, MPI_COMM_WORLD);
    MPI_Send(&rows, 1, MPI_INT, dest, mtype, MPI_COMM_WORLD);
    MPI_Send(&a[offset][0], rows*SIZE, MPI_LONG_DOUBLE, dest, mtype,
MPI_COMM_WORLD);
    MPI_Send(&b[offset][0], rows*SIZE, MPI_LONG_DOUBLE, dest, mtype,
MPI_COMM_WORLD);

    offset += rows;
}

//-----
//Receiving results from Workers

mtype = FROM_WORKER;
for(i = 1; i <= workers; i++)
{
    source = i;
```

```
        MPI_Recv(&offset, 1, MPI_INT, source, mtype, MPI_COMM_WORLD, &status);
        MPI_Recv(&rows, 1, MPI_INT, source, mtype, MPI_COMM_WORLD, &status);
        MPI_Recv(&c[offset][0], rows*SIZE, MPI_LONG_DOUBLE, source, mtype,
MPI_COMM_WORLD, &status);

        // printf("Received results from task=%d\n", source);
    }
    // printf("Result Matrix\n");

    // for(i = 0; i < SIZE;i++)
    // {
    //     // printf("\n");
    //     for(int j = 0; j < SIZE;j++)
    //     {
    //         // printf("%6.2Lf  ",c[i][j]);
    //     }
    // }
    end = MPI_Wtime();

    exec = end - start;
    printf("MPI Exec time - %f\n", exec);
}

//-----
//-----

//-----
//-----

//Worker Task
if(taskid > MASTER)
{
    mtype = FROM_MASTER;

    MPI_Recv(&offset, 1, MPI_INT, MASTER, mtype, MPI_COMM_WORLD,&status);
    MPI_Recv(&rows, 1, MPI_INT, MASTER, mtype, MPI_COMM_WORLD,&status);
    MPI_Recv(&a, rows*SIZE, MPI_LONG_DOUBLE, MASTER, mtype, MPI_COMM_WORLD,&status);
    MPI_Recv(&b, rows*SIZE, MPI_LONG_DOUBLE, MASTER, mtype, MPI_COMM_WORLD,&status);

    for(k = 0; k < SIZE; k++)
    {
```

```
        for(i = 0; i < rows; i++)
        {
            c[i][k] = a[i][k] * b[i][k];
        }
    }

    mtype = FROM_WORKER;

    MPI_Send(&offset,1, MPI_INT, MASTER,mtype, MPI_COMM_WORLD);
    MPI_Send(&rows,1, MPI_INT, MASTER,mtype, MPI_COMM_WORLD);
    MPI_Send(&c, rows*SIZE, MPI_LONG_DOUBLE, dest, mtype, MPI_COMM_WORLD);

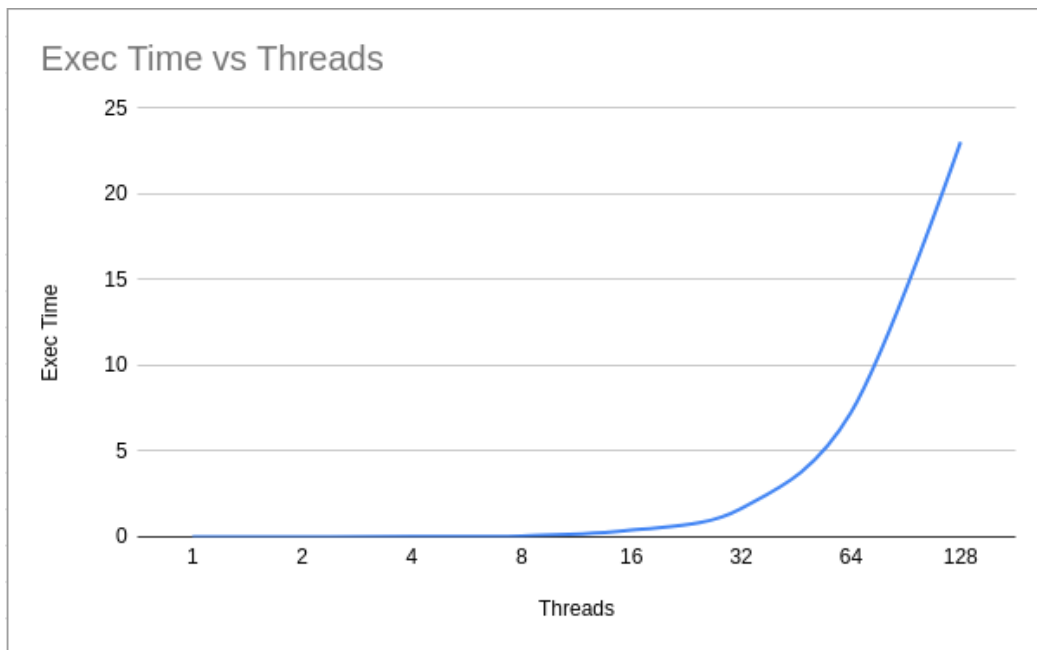
}
MPI_Finalize();

return 0;
}
```

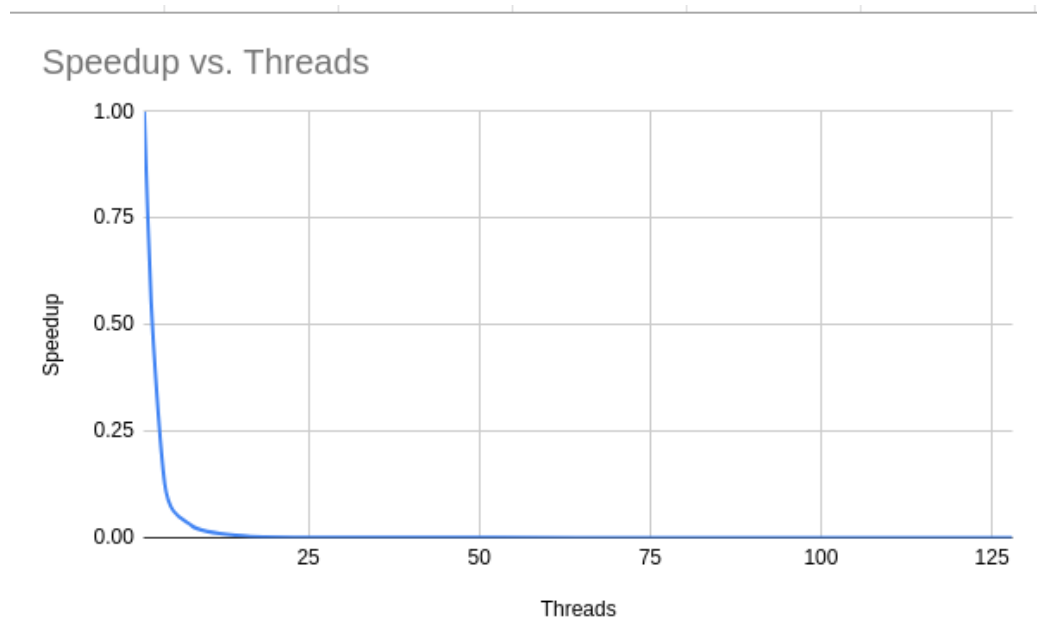
## Observations

no of processors	Exec Time	Speedup	Parallelization Factor
1	0.000359	1	0
2	0.002854	0.5557112824	-0.666246322
4	0.0126	0.1258730159	-4.629676335
8	0.057482	0.02759124596	-11.74779319
16	0.392194	0.004043917041	82.0949979
32	1.652586	0.000959707997	1734.972678
64	7.239875	0.0002190645557	19776.7459
128	23.03238	0.00006885957943	140372.641

## Threads vs Time



## Speedups vs Threads



## **Inferences**

Since MPI is a distributed memory architecture, the communication overhead between nodes causes the parallel code to run slower compared to serial code ( running in 1 node or only in master )