PART – A : Open MP programs

Execution Commands:

For compilation: gcc -fopenmp filename.c

For running: ./a.out

1. Given a n*n matrix arr and a vector vec of length n, their product res=arr*vec . Write a program to implement the multiplication using OpenMP PARALLEL directive.

Code:

```
#include<stdio.h>
#include<stdlib.h>
#include<omp.h>
void main() {
       int m,n;
        printf("Enter the size of square matrix : ");
       scanf_s("%d",&n);
       printf("Enter the size of vector : ");
       scanf_s("%d", &m);
        if (m!=n) {
                printf("Multiplication is not possible.\n");
               exit(0);
        }
        int i=0, j=0;
        int **arr=(int**)malloc(n*sizeof(int*));
        int *vec=(int*)malloc(n*sizeof(int));
        int *res=(int*)malloc(n*sizeof(int));
        omp set num threads(n);
        #pragma omp parallel private(j)
        {
                #pragma omp for
                for (i=0;i<n;i++) {
                        srand(i);
                        arr[i] = (int*) malloc (n*sizeof (int));
                        vec[i]=rand()%100;
                        for (j=0; j < n; j++)
                                arr[i][j]=rand()%100;
        #pragma omp parallel private(j)
                #pragma omp for
                for(i=0;i<n;i++) {
                       res[i]=0;
                        for(j=0;j<n;j++)
                               res[i]+=arr[i][j]*vec[j];
                }
        }
        printf("Matrix * Vector = Resultant Matrix\n");
        for(i=0;i<n;i++) {
                for(j=0;j<n;j++)
                       printf("%3d ",arr[i][j]);
                if(i==n/2)
                       printf(" * %3d = %6d\n", vec[i], res[i]);
```

2. Consider a scenario where a person visits a supermarket for shopping. He purhases various items in different sections such as clothing, gaming, grocery, stationary.

Code:

```
#include<stdio.h>
#include<stdlib.h>
#include<omp.h>
void main() {
    int r,i,ans=0;
    printf("Enter number of sections : ");
    scanf s("%d",&r);
    int **arr=(int**)malloc(r*sizeof(int*));
    int *size=(int*)malloc(r*sizeof(int));
    omp set num threads(r);
        #pragma omp parallel
        #pragma omp for
        for (i=0;i<r;i++) {
            srand(i);
            int j,sum=0;
            size[i]=rand()%20;
                    arr[i] = (int*) malloc(size[i] *sizeof(int));
                    for (j=0;j<size[i];j++) {</pre>
                            arr[i][j]=rand()%100;
                sum+=arr[i][j];
            #pragma omp critical
                ans+=sum;
            }
    for(i=0;i<r;i++) {
        printf("Section - %2d ( %3d Items ) : ",i,size[i]);
        for(int j=0;j<size[i];j++)</pre>
            printf("%3d ",arr[i][j]);
        printf("\n");
    printf("Total Amount : %d",ans);
}
```

3. Consider a person named X on the earth, to find its accurate position on the globe, we require the value of pi.

Code:

a. Using Parellel

```
#include<stdio.h>
```

```
#include<stdlib.h>
#include<time.h>
#include<omp.h>
void main() {
    unsigned long long int num, i;
    printf("Enter the number of steps : ");
    scanf_s("%llu",&num);
    time t st,et;
    st=clock();
    double step=1.0/(double)num, pi=0.0;
    double *local pi=(double*)malloc(num*sizeof(double));
    omp_set_num_threads(num);
    #pragma omp parallel for
    for(i=0;i<num;i++) {
        double x=(i+0.5)*step;
        local pi[i] = (4.0*step) / (1+x*x);
    for(i=0;i<num;i++)</pre>
       pi+=local pi[i];
    et=clock();
    printf("Time Taken : %lf\n", (double) ((double) (et-st)/CLOCKS PER SEC));
    printf("Value of Pi = %lf\n",pi);
}
```

b. Using Atomic

```
#include<stdio.h>
#include<stdlib.h>
#include<time.h>
#include<omp.h>
void main() {
    int num, i;
    printf("Enter the number of steps : ");
    scanf s("%d", &num);
    time t st, et;
    st = clock();
    double step = 1.0 / (double) num, pi = 0.0;
    omp_set_num_threads(num);
    #pragma omp parallel for
    for (i = 0; i < num; i++) {
        double x = (i + 0.5) * step;
        double local pi = (4.0 * step) / (1 + x * x);
        #pragma omp atomic
        pi += local pi;
    }
    et = clock();
    printf("Time Taken : %lf\n", (double)((double)(et - st) /
CLOCKS PER SEC));
   printf("Value of Pi = %lf\n", pi);
```

c. Using Critical

```
#include<stdlib.h>
#include<time.h>
#include<omp.h>
void main() {
    int num, i;
    printf("Enter the number of steps : ");
    scanf s("%d", &num);
    time t st,et;
    st=clock();
    double step=1.0/(double)num,pi=0.0;
    omp set num threads(num);
    #pragma omp parallel for
    for(i=0;i<num;i++) {</pre>
        double x=(i+0.5)*step;
        double local pi=(4.0*step)/(1+x*x);
        #pragma omp critical
            pi+=local pi;
    et=clock();
    printf("Time Taken : %lf\n", (double) ((double) (et-st)/CLOCKS PER SEC));
    printf("Value of Pi = %lf\n",pi);
}
```

d. Using Reduction

```
#include<stdio.h>
#include<stdlib.h>
#include<time.h>
#include<omp.h>
void main() {
    int num, i;
    printf("Enter the number of steps : ");
    scanf s("%d", &num);
    time t st,et;
    st=clock();
    double step=1.0/(double)num,pi=0.0;
    omp set num threads(num);
    #pragma omp parallel for reduction(+:pi)
    for(i=0;i<num;i++) {
        double x=(i+0.5)*step;
        double local pi=(4.0*step)/(1+x*x);
        pi+=local pi;
    }
    et=clock();
    printf("Time Taken : %lf\n", (double) ((double) (et-st)/CLOCKS PER SEC));
    printf("Value of Pi = %lf\n",pi);
}
```

4. Design, Develop and run a multithreaded program to generate and print Fibonacci series, one thread must generate the series upto number and other thread must print them. Ensure proper synchronization.

a. Using single:

#include<stdio.h>

```
#include<stdlib.h>
#include<time.h>
#include<omp.h>
int main() {
       int n, i;
       printf("Number of terms : ");
       scanf s("%d",&n);
       int* a = (int*)malloc(n * sizeof(int));
       a[0] = 0;
       a[1] = 1;
       time_t st, et;
       st = clock();
       omp_set_num_threads(2);
       #pragma omp parallel
               #pragma omp single
                       printf("id of thread involved in the computation of
fibonacci numbers = %d\n", omp_get_thread_num());
                       for (i = 2; i < n; i++)
                              a[i] = a[i - 2] + a[i - 1];
               #pragma omp single
                       printf("id of thread involved in the displaying of
fibonacci numbers = %d\n", omp get thread num());
                       printf("Fibonacci numbers : ");
                       for (i = 0; i < n; i++)
                               printf("%d ", a[i]);
                       printf("\n");
               }
       }
       et = clock();
       printf("Time Taken : %lfms\n", ((double)(et - st)*1000 /
CLOCKS PER SEC));
       return 0;
```

b. <u>Using critical:</u>

```
#include<stdio.h>
#include<stdlib.h>
#include<time.h>
#include<omp.h>
int main() {
       int n, i;
       printf("Number of terms : ");
       scanf_s("%d",&n);
       int* a = (int*)malloc(n * sizeof(int));
       a[0] = 0;
       a[1] = 1;
       time t st, et;
       st = clock();
       omp set num threads(2);
        #pragma omp parallel
                #pragma omp critical
                if(omp get thread num() == 0)
```

```
printf("id of thread involved in the computation of
fibonacci numbers = %d\n", omp_get_thread_num());
                       for (i = 2; i < n; i++)
                               a[i] = a[i - 2] + a[i - 1];
               else if(omp get thread num()==1)
                       printf("id of thread involved in the displaying of
fibonacci numbers = %d\n", omp get thread num());
                       printf("Fibonacci numbers : ");
                       for (i = 0; i < n; i++)
                               printf("%d ", a[i]);
                       printf("\n");
               }
       }
       et = clock();
       printf("Time Taken : %lfms\n", ((double)(et - st)*1000 /
CLOCKS PER SEC));
       return 0;
```

5. University awards gold medal to the student who has scored highest. Write an OpenMP program to find the student with highest cgpa.

```
#include<stdio.h>
#include<stdlib.h>
#include<time.h>
#include<omp.h>
int main() {
       int n, i;
       time_t st, et;
       st = clock();
       printf("Enter the number of students : ");
       scanf_s("%d", &n);
       double* arr = (double*)malloc(n * sizeof(double));
       double arr max = 0;
       #pragma omp parallel for
       for (i = 0; i < n; i++) {
               srand(i);
               arr[i] = (double) (rand() % 10000)/10;
       printf("CGPA of students : ");
       for (i = 0; i < n; i++)
               printf("%.21f ", arr[i]);
       printf("\n");
       #pragma omp parallel for
       for (i = 0; i < n; i++) {
               #pragma omp critical
               if (arr_max < arr[i])</pre>
                       arr_max = arr[i];
        }
       et = clock();
       printf("Student with highest CGPA = %.21f\n", arr max);
       printf("Time Taken : %.21fms\n", ((double)(et - st) * 1000 /
CLOCKS PER SEC));
```

}

6. Multiply two square matrices (1000, 2000 or 3000 dimensions). Compare the performance of a squential and parallel algorithm using openMP.

```
#include<stdio.h>
#include<stdlib.h>
#include<time.h>
#include<omp.h>
void main() {
       int n;
       printf("Enter the dimension of square matrices : ");
       scanf_s("%d", &n);
       int i = 0, j = 0, k = 0;
       int** arr1 = (int**)malloc(n * sizeof(int*));
       int** arr2 = (int**)malloc(n * sizeof(int*));
       int** res = (int**)malloc(n * sizeof(int*));
       omp_set_num_threads(64);
       #pragma omp parallel private(j)
               #pragma omp for
               for (i = 0; i < n; i++) {
                       srand(i);
                       arr1[i] = (int*)malloc(n * sizeof(int));
                       arr2[i] = (int*)malloc(n * sizeof(int));
                       res[i] = (int*)malloc(n * sizeof(int));
                       for (j = 0; j < n; j++) {
                               arr1[i][j] = rand() % 100;
                               arr2[i][j] = rand() % 100;
                       }
                }
       time t st, et;
       st = clock();
        #pragma omp parallel private(j,k)
               #pragma omp for
               for (i = 0; i < n; i++) {
                       for (j = 0; j < n; j++) {
                               res[i][j] = 0;
                               for (k = 0; k < n; k++)
                                       res[i][j] += arr1[i][k] * arr2[k][j];
                       }
                }
        }
       et = clock();
       printf("Time taken by parallel algorithm : f^n, (double) (et -
st) / CLOCKS_PER_SEC);
       st = clock();
        for (i = 0; i < n; i++) {
               for (j = 0; j < n; j++) {
                       res[i][j] = 0;
                       for (k = 0; k < n; k++)
                               res[i][j] += arr1[i][k] * arr2[k][j];
                }
       et = clock();
```

```
printf("Time taken by Sequential algorithm : %lf\n", (double)(et -st) / CLOCKS_PER_SEC); }
```

PART B – MPI Programs

Execution Commands:

For compilation: mpicc filename.c

For running: mpiexec -np <no_of_processes> ./a.out

7. Assume you have n robots which pick mangoes in a farm. Write a program to calculate the total number of mangoes picked by n robots parallelly using MPI.

```
#include<stdio.h>
#include<stdlib.h>
#include<mpi.h>
int main(int argc, char** argv)
    int rank, numproc;
    int sum = 0;
    int total sum = 0;
    MPI Init(&argc, &argv);
    MPI Comm size (MPI COMM WORLD, &numproc);
   MPI Comm rank (MPI COMM WORLD, &rank);
    srand(rank);
    sum = rand() % 100;
    printf("Robot %d picked %d mangoes.\n", rank, sum);
    MPI Reduce (&sum, &total sum, 1, MPI INT, MPI SUM, 0, MPI COMM WORLD);
    if (rank == 0)
        printf("Total Mangoes picked by %d Robots = %d\n", numproc,
total sum);
   MPI Finalize();
```

8. Design a program that implements application of MPI Collective Communications

```
#include <stdio.h>
#include <stdlib.h>
#include <mpi.h>
int main(int argc, char* argv[])
{
       int size, rank;
       MPI Init(&argc, &argv);
       MPI Comm rank (MPI COMM WORLD, &rank);
       MPI Comm size (MPI COMM WORLD, &size);
       float recvbuf, sendbuf[100];
       if (rank == 0) {
               int i;
               printf("Before Scatter : sendbuf of rank 0 : ");
               for (i = 0; i < size; i++) {
                       srand(i);
                       sendbuf[i] = (float)(rand()%1000)/10;
                       printf("%.1f ", sendbuf[i]);
```

9. Implement Cartesian Virtual Topology in MPI.

```
#include <stdio.h>
#include <stdlib.h>
#include <mpi.h>
#define SIZE 16
#define UP 0
#define DOWN 1
#define LEFT 2
#define RIGHT 3
int main(int argc, char* argv[])
       int numtasks, rank, source, dest, outbuf, i, tag = 1, inbuf[4] = {
MPI PROC NULL, MPI PROC NULL, MPI PROC NULL, MPI PROC NULL, }, nbrs[4],
dims[2] = \{ 4, 4 \}, periods[2] = \{ 0, 0 \}, reorder = 0, coords[2];
       MPI Request reqs[8];
       MPI Status stats[8];
       MPI_Comm cartcomm;
       MPI Init(&argc, &argv);
       MPI Comm size(MPI COMM WORLD, &numtasks);
       if (numtasks == SIZE) {
               MPI Cart create (MPI COMM WORLD, 2, dims, periods, reorder,
&cartcomm);
               MPI Comm rank(cartcomm, &rank);
               MPI Cart coords (cartcomm, rank, 2, coords);
               MPI_Cart_shift(cartcomm, 0, 1, &nbrs[UP], &nbrs[DOWN]);
               MPI Cart shift(cartcomm, 1, 1, &nbrs[LEFT], &nbrs[RIGHT]);
               printf("rank= %d coords= %d %d neighbors(u,d,l,r)= %d %d %d
%d\n", rank, coords[0], coords[1], nbrs[UP], nbrs[DOWN], nbrs[LEFT],
nbrs[RIGHT]);
               outbuf = rank;
               for (i = 0; i < 4; i++) {
                       dest = nbrs[i];
                       source = nbrs[i];
                       MPI_Isend(&outbuf, 1, MPI INT, dest, tag,
MPI COMM WORLD, &reqs[i]);
                       MPI Irecv(&inbuf[i], 1, MPI INT, source, tag,
MPI COMM WORLD, &reqs[i + 4]);
               MPI Waitall(8, reqs, stats);
               printf("rank= %d inbuf(u,d,l,r)= %d %d %d %d\n", rank,
inbuf[UP], inbuf[DOWN], inbuf[LEFT], inbuf[RIGHT]);
       }
       else
               printf("Must specify %d tasks. Terminating.\n", SIZE);
       MPI Finalize();
```

10. Design a MPI program to simulate the uses blocking send/receive routines and nonblocking send/receive routines.

```
#include<stdio.h>
#include<stdlib.h>
#include<time.h>
#include<mpi.h>
int main(int argc, char* argv[])
       int numtasks, rank, rc, count, next, prev, sz, inmsg;
       MPI Status Stat;
       time t st, et;
       MPI Init(&argc, &argv);
       MPI_Comm_size(MPI COMM WORLD, &numtasks);
       sz = (numtasks / \overline{2}) * \overline{2};
       MPI Comm rank (MPI COMM WORLD, &rank);
       st = clock();
       if (rank == 0) prev = sz - 1;
       else prev = rank - 1;
       if (rank == sz - 1) next = 0;
       else next = rank + 1;
       if (rank % 2 == 0 \&\& rank < sz) {
               rc = MPI Send(&rank, 1, MPI INT, next, 0, MPI COMM WORLD);
               rc = MPI Recv(&inmsg, 1, MPI INT, prev, 1, MPI COMM WORLD,
&Stat);
       else if (rank % 2 == 1 && rank < sz) {
               rc = MPI Recv(&inmsg, 1, MPI INT, prev, 0, MPI COMM WORLD,
&Stat);
               rc = MPI Send(&rank, 1, MPI INT, next, 1, MPI COMM WORLD);
       MPI Barrier (MPI COMM WORLD);
       et = clock();
       if(rank==0) printf("Time taken by Blocking send/receive : f^*
(double) (et - st) / CLOCKS PER SEC);
       MPI Barrier (MPI COMM WORLD);
       MPI Request reqs[2];
       MPI Status stats[2];
       st = clock();
       if (rank == numtasks - 1) next = 0;
       else next = rank + 1;
       if (rank == 0) prev = numtasks - 1;
       else prev = rank - 1;
       MPI_Irecv(&inmsg, 1, MPI_INT, prev, 0, MPI_COMM_WORLD, &reqs[0]);
       MPI_Isend(&rank, 1, MPI_INT, next, 0, MPI_COMM_WORLD, &reqs[1]);
       MPI Barrier(MPI COMM WORLD);
       et = clock();
       if (rank == 0) printf("Time taken by NonBlocking send/receive :
%lf\n", (double)(et - st) / CLOCKS_PER_SEC);
       MPI Finalize();
}
```
