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File: piCalculate.c
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if (n == 0) exit(0);

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/********************
This exercise presents a simple program to determine the value of pi.
The algorithm suggested here is chosen for its simplicity. The method
evaluates the integral of 4/(1+x^*x) between 0 and 1. The method is
simple: the integral is approximated by a sum of n intervals; the
approximation to the integral in each interval is (1/n)*4/(1+x*x). The
master process (rank 0) asks the user for the number of intervals; the
master should then broadcast this number to all of the other processes.
Each process then adds up every n'th interval (x = rank/n,
rank/n+size/n,...). Finally, the sums computed by each process are added
together using a reduction.
You may want to use these MPI routines in your solution:
MPI Bcast MPI Reduce
#include "mpi.h"
#include <math.h>
#include <stdio.h>
#include <stdlib.h>
#define MAX NAME 80
                   /* length of characters for naming a process */
int main(argc, argv)
int argc;
char *argv[];
{
    int n,
    rank, /* rank variable to identify the process */
    numprocs, /* number of processes */
    len; /* variable for storing name of processes */
    double PI25DT = 3.141592653589793238462643; /* 25-digit-PI*/
    double mypi, /* value from each process */
    pi, /* value of PI in total*/
    step, /* the step */
    sum, /* sum of area under the curve */
     startTime, /* starting time */
      endTime; /* ending time */
    char name[MAX_NAME]; /* char array for storing the name of
each process */
    /*Initialize MPI execution environment */
    MPI_Init(&argc,&argv);
    MPI_Comm_size(MPI_COMM_WORLD,&numprocs);
    MPI_Comm_rank(MPI_COMM_WORLD,&rank);
    MPI_Get_processor_name(name, &len);
    if (rank == 0) {
         printf("Number of processing units: %d\n", numprocs);
         printf("Enter the number of intervals: (0 quits) ");
         scanf("%d",&n);
    MPI_Bcast(&n, 1, MPI_INT, 0, MPI_COMM_WORLD);
```

```
//Broadcast the number of bins to all processes
     /* This broadcasts an integer which is n, from the master to all
processes */
     //Calculating for each core
     startTime = MPI_Wtime();
          = 1.0 / (double) n;
     sum = 0.0;
     for (i = rank + 1; i \le n; i += numprocs) {
         x = step * ((double)i - 0.5);
          sum += 4.0 / (1.0 + x*x);
     mypi = step * sum;
     printf("This is my sum: %.16f from rank: %d name: %s\n", mypi,
rank, name);
     //Now we can reduce all those sums to one value which is pi
     MPI_Reduce(&mypi, &pi, 1, MPI_DOUBLE, MPI_SUM, 0,
             MPI_COMM_WORLD);
     if (rank == 0){
          printf("pi is approximately %.16f, Error is %.16f\n", pi,
fabs(pi - PI25DT));
          endTime = MPI_Wtime();
          printf("Time of calculation PI is: %f\n", endTime-startTime);
     //Terminate MPI execution environment
     MPI Finalize();
     return 0;
}
#OUTPUT:
```

1 Node:

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🛛 🗐 📵 student@node2: /home/ibm
student@node2:/home/ibm$ mpiexec -np 12 -host node1 ./piCalculate
Number of processing units: 12
Enter the number of intervals: (0 quits) 1000
This is my sum: 0.2640498684137977 from rank: 0 name: node1
This is my sum: 0.2638827058653734 from rank: 1 name: node1
This is my sum: 0.2637153742335374 from rank: 2 name: node1
This is my sum: 0.2607168601078259 from rank: 8 name: node1
This is my sum: 0.2605503582258470 from rank: 9 name: node1
This is my sum: 0.2635478743516046 from rank: 3 name: node1
This is my sum: 0.2613812068029122 from rank: 4 name: node1
This is my sum: 0.2612153709208150 from rank: 5 name: node1
This is my sum: 0.2610493672886828 from rank: 6 name: node1
This is my sum: 0.2603836919273075 from rank: 10 name: node1
This is my sum: 0.2602168620455291 from rank: 11 name: node1
This is my sum: 0.2608831967398937 from rank: 7 name: node1
pi is approximately 3.1415927369231262, Error is 0.0000000833333331
Time of calculation PI is: 0.092713
student@node2:/home/ibm$
```

2 Nodes Machine File:

machine.txt:

node1:4 node2:4

```
student@node2: /home/ibm
student@node2:/home/ibm$ mpiexec -np 12 -f machine.txt ./piCalculate
Number of processing units: 12
Enter the number of intervals: (0 quits) 1000
This is my sum: 0.2640498684137977 from rank: 0 name: node1
This is my sum: 0.2613812068029122 from rank: 4 name: node2
This is my sum: 0.2637153742335374 from rank: 2 name: node1
This is my sum: 0.2608831967398937 from rank: 7 name: node2
This is my sum: 0.2610493672886828 from rank: 6 name: node2
This is my sum: 0.2612153709208150 from rank: 5 name: node2
This is my sum: 0.2635478743516046 from rank: 3 name: node1
This is my sum: 0.2607168601078259 from rank: 8 name: node1
This is my sum: 0.2603836919273075 from rank: 10 name: node1
This is my sum: 0.2638827058653734 from rank: 1 name: node1
This is my sum: 0.2602168620455291 from rank: 11 name: node1
This is my sum: 0.2605503582258470 from rank: 9 name: node1
pi is approximately 3.1415927369231262, Error is 0.00000008333333331
Time of calculation PI is: 0.047710
student@node2:/home/ibm$
```

3 Nodes:

```
student@node2: /home/ibm
student@node2:/home/ibm$ mpiexec -np 12 -host node1,node2,node3 ./piCalculate
Number of processing units: 12
Enter the number of intervals: (0 quits) 1000
This is my sum: 0.2637153742335374 from rank: 2 name: node3
This is my sum: 0.2612153709208150 from rank: 5 name: node3
This is my sum: 0.2607168601078259 from rank: 8 name: node3
This is my sum: 0.2602168620455291 from rank: 11 name: node3
This is my sum: 0.2640498684137977 from rank: 0 name: node1
This is my sum: 0.2638827058653734 from rank: 1 name: node2
This is my sum: 0.2635478743516046 from rank: 3 name: node1
This is my sum: 0.2608831967398937 from rank: 7 name: node2
This is my sum: 0.2610493672886828 from rank: 6 name: node1
This is my sum: 0.2603836919273075 from rank: 10 name: node2
This is my sum: 0.2605503582258470 from rank: 9 name: node1
This is my sum: 0.2613812068029122 from rank: 4 name: node2
pi is approximately 3.1415927369231262, Error is 0.0000000833333331
Time of calculation PI is: 0.015973
student@node2:/home/ibmS
```