Parallel Processing

Lecture 9 Guy Tel-Zur

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Today's agenda

- Final Presentations status
- Continue with OpenMP
- CilkPlus
- Parallel Matlab
- Sorting Algorithms (slides10)
- Load Balancing (slides7)
- Home assignment #3

The Course Roadmap

Algorithms

Embarrassingly Parallel

Partitioning and Divide & Conqur

Synchronous Computations

Load Balancing

Today Sorting

time

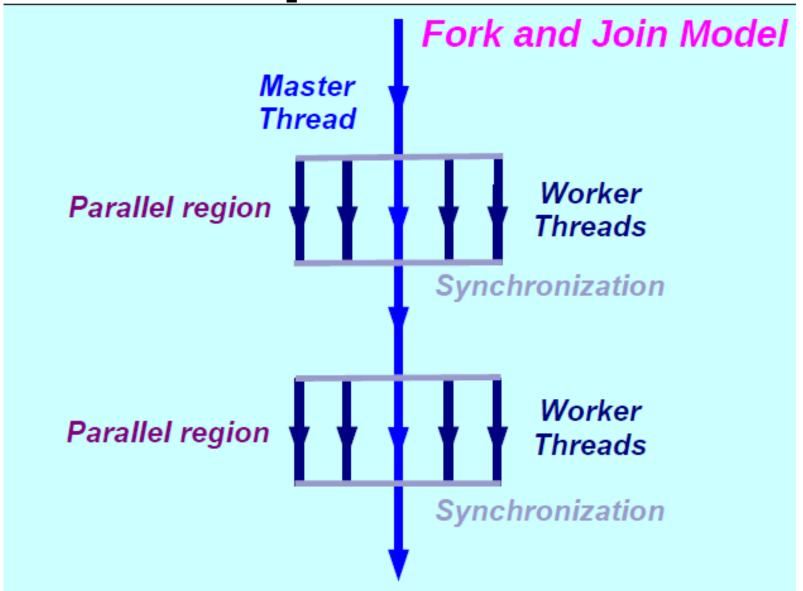
Numerical Algorithms

MPI tip - shortcuts using define

```
#define MASTER 0
#define Bcast(send_data, count, type)
MPI_Bcast(send_data, count, type, MASTER,
MPI_COMM_WORLD) //root --> MASTER
#define Finalize() MPI_Finalize()
#define Init(x,y) MPI_Init(x,y)
#define Rank(x) MPI_Comm_rank(MPI_COMM_WORLD, x)
#define Size(x) MPI_Comm_size(MPI_COMM_WORLD, x)
```



The OpenMP Model



Exercise 2: A simple SPMD pi program

```
#include <omp.h>
static long num_steps = 100000;
                                    double step;
#define NUM_THREADS 2
void main ()
          int i, nthreads; double pi, sum[NUM_THREADS];
          step = 1.0/(double) num_steps;
          omp set num threads(NUM THREADS);
  #pragma omp parallel
         int i, id,nthrds;
        double x:
        id = omp_get_thread_num();
        nthrds = omp_get_num_threads();
        if (id == 0) nthreads = nthrds;
          for (i=id, sum[id]=0.0;i< num_steps; i=i+nthrds) {
                  x = (i+0.5)*step;
                  sum[id] += 4.0/(1.0+x*x);
```

Promote scalar to an array dimensioned by number of threads to avoid race condition.

Only one thread should copy the number of threads to the global value to make sure multiple threads writing to the same address don't conflict.

This is a common trick in SPMD programs to create a cyclic distribution of loop iterations

for(i=0, pi=0.0;i<nthreads;i++)pi += sum[i] * step;



Exercise 3: SPMD Pi without false sharing

```
#include <omp.h>
static long num_steps = 100000;
                                      double step;
#define NUM_THREADS 2
void main ()
                            step = 1.0/(double) num steps;
          double pi;
          omp_set_num_threads(NUM_THREADS);
#pragma omp parallel
                                                        Create a scalar local to
                                                        each thread to
         int i, id,nthrds; double x, sum;
                                                        accumulate partial
         id = omp_get_thread_num();
                                                        sums.
         nthrds = omp_get_num_threads();
         if (id == 0) nthreads = nthrds;
          id = omp get thread num();
         nthrds = omp_get_num_threads();
          for (i=id, sum=0.0;i< num steps; i=i+nthreads){
                                                                      No array, so
                   x = (i+0.5)*step;
                                                                      no false
                   sum += 4.0/(1.0+x^*x);
                                                                      sharing.
                                          Sum goes "out of scope" beyond the parallel
        #pragma omp critical
                                          region ... so you must sum it in here. Must
               pi += sum * step;
                                          protect summation into pi in a critical region so
                                          updates don't conflict
```

Exercise 4: solution

```
#include <omp.h>
     static long num steps = 100000;
                                             double step;
     #define NUM THREADS 2
                                                     For good OpenMP
     void main ()
                                                     implementations,
                                                     reduction is more
              int i; double x, pi, sum = 0.0;
                                                    scalable than critical.
              step = 1.0/(double) num steps;
              omp set num threads(NUM THREADS);
     #pragma omp parallel for private(x) reduction(+:sum)
              for_{i=0;i< num steps; i++}
                      x = (i+0.5)*step;
i private by
                      sum = sum + 4.0/(1.0+x*x);
default
                                             Note: we created a parallel
              pi = step * sum;
                                             program without changing
                                             any code and by adding 4
                                                   simple lines!
```

Parallel Programmers love Monte Carlo Embarrassingly parallel: the

algorithms

```
#include "omp.h"
                                                     embarrassing.
static long num_trials = 10000;
                                            Add two lines and you have a
int main ()
                                                    parallel program.
  long i; long Ncirc = 0; double pi, x, y;
  double r = 1.0; // radius of circle. Side of squrare is 2*r
  seed(0,-r, r); // The circle and square are centered at the origin
 #pragma omp parallel for private (x, y) reduction (+:Ncirc)
  for(i=0;i<num_trials; i++)
   x = random(); y = random();
   if (x*x + y*y) \le r*r Ncirc++;
  pi = 4.0 * ((double)Ncirc/(double)num_trials);
  printf("\n %d trials, pi is %f \n",num_trials, pi);
```

parallelism is so easy its

Compiler notes: Visual Studio

- Start "new project"
- Select win 32 console project
 - Set name and path
 - On the next panel, Click "next" instead of finish so you can select an empty project on the following panel.
 - Drag and drop your source file into the source folder on the visual studio solution explorer
 - Activate OpenMP
 - Go to project properties/configuration properties/C.C++/language ... and activate OpenMP
- Set number of threads inside the program
- Build the project
- Run "without debug" from the debug menu.