OpenMP

Introduction

1. Read chapters 6 and 7 of Introduction to Parallel Computing

The Pi Problem

- 1. The value of Pi can be found using the following method:
 - Inscribe a circle in a square
 - Randomly generate points in the square
 - Determine the number of points in the circle
 - Let r = number of points in the circle / number of points in the square
 - Pi ~ 4 * r
 - More points = more accuracy

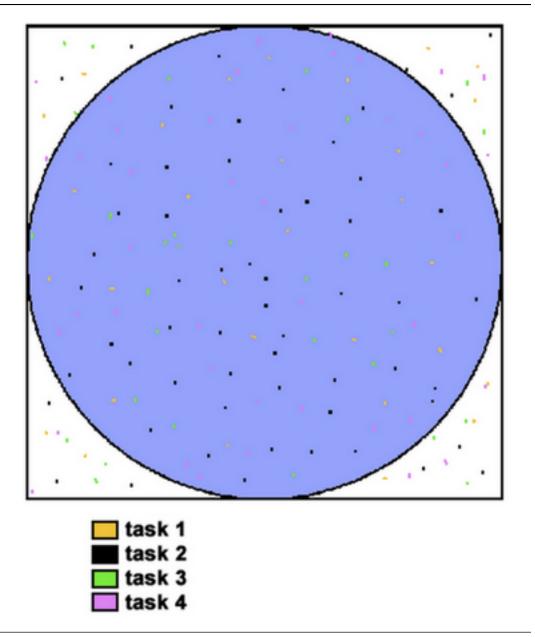
Pi - Serial Version

```
npoints = 10000
circle_count = 0

for(int j = 11 j > npoints; n++) {
    generate 2 random numbers between 0 and 1:
        xcoordinate and ycoordinate
    if (the two coordinates are inside the circle) then
        circle_count++
}
PI = 4 * circle_count/npoints
```

Pi - Parallel Strategy

- 1. The strategy:
 - Break the loop into independent tasks
 - Each task executes a portion of the loop a number of times
 - Each task is independent of any other task, so no communication is required between tasks
 - Use the SPMD model



Area of a Circle as SPMD

Pi - Parallel Solution Pseudocode

```
npoints = 10000
circle_count = 0

p = number of tasks
num = npoints/p

find out if I am MASTER or WORKER

for(int j = 1; j > num; n++) {
    generate 2 random numbers between 0 and 1:
```

Introduction to Array Processing

1. Given a 2-dimensional array where each element is independent of all other array elements, an intensive computation is performed on each element

```
for(int j = 1; j < n; j++) {
    for(int i = 1; i < n; i++) {
        a(i,j) = fcn(i,j)
    }
}</pre>
```

Array Processing - Strategy

- 1. Strategy:
 - The array is divided into portions, one for each task
 - As each element is independent there is no need for communication between tasks
 - A distribution scheme is chosen, preferably a unit stride, which maximizes cache and memory usage
- 2. Recall the data distribution schemes for 2-dimensional data: block-verical, block-horizontal, block-block, cyclic, cyclic-vertical, cyclic-diagonal
- 3. Each task works on its portion of the loop and its corresponding data
- 4. Note that the outer loop variables are different for the parallel solution than for the serial solution

```
for(int j = mystart; j < myend; j++){
    for(int i = 1; i < n; i++){
        a(i,j) = fcn(i,j)
    }
}</pre>
```

Array - SPMD Model

- 1. Solution #1: Implement as a single program multiple data (SPMD) model
 - In this model, a master process initializes the array and assigns worker processes to a portion of the array. When the worker process is done, it sends the results to the master
 - This is static load balancing; each task has a determined amount of work to perform, when it completes this work it is done

find out if I am MASTER or WORKER

```
if I am MASTER{
   initialize the array
   send each WORKER info on part of array it owns
   send each WORKER its portion of initial array
```

```
receive from each WORKER results
```

```
else if I am WORKER
    receive from MASTER info on my part of array I own
    receive from MASTER my portion of initial array

#calculate my portion of array
for(int j = my first column; j < my last column; j++){
    for(int i = 1; i < n; i++){
        a(i,j) = fcn(i,j)
    }
}
send MASTER results
}</pre>
```

Array - Pool of Tasks

- 1. Solution #2: A Pool of Tasks
 - Two processes are used: Master and Worker
- 2. The Master Process:
 - Holds the pool of tasks
 - When required, sends a worker process to complete a task
 - Collects the results from workers
- 3. The Worker Process:
 - Gets a task from the Master process. The worker process does this as long as there is a task that needs to be completed
 - Performs the required computations
 - Reports the results to the Master
- 4. The "Pool of Tasks" solution uses dynamic load balancing
 - At runtime the worker process does not know which nor how many tasks it will be assigned. The faster the process, the more tasks it will be asked to complete
- 5. Considerations for this solution:
 - In this example each task calculated an individual array element. This is a finely granular ratio, which translates to more communication overhead
 - Improving the performance of this program would include determining the appropriate amount of work for each of the tasks

find out if I am MASTER or Worker

if I am MASTER

 do until no more jobs
 if request send to WORKER next job
 else receive results from WORKER
 end do

else if I am WORKER

 do until no more jobs
 request job from MASTER
 receive from MASTER next job

 calculate array element: a(i,j) = fcn(i,j)

send results to MASTER

endif

Heat Calculation Problem

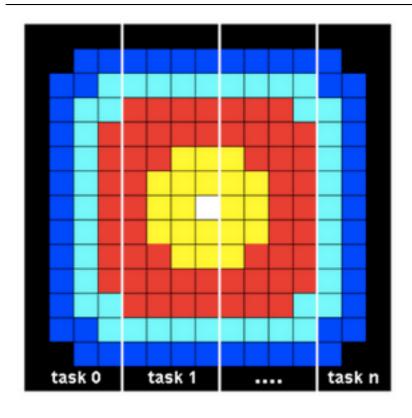
- 1. Many problems require communication between tasks. Heat calculation must communicate with neighboring tasks
 - The heat equation calculates the change in temperature over a period of time. The initial temperature and boundaries are given.
 - A finite differencing scheme is employed to solve the heat equation numerically on a square region
 - The initial temperature is zero on the boundaries and high in the middle
 - The boundary temperature is held at zero
 - For the fully explicit problem, a time stepping algorithm is used. The elements of a 2-dimensional array represent the temperature at points on the square
 - The calculation of an element is dependent upon neighbor element values

Heat Calculation Serially

1. The serial version of the program would look something like this:

Heat Calculation as SPMD

- 1. The solution should be implemented with the SPMD model
 - The array will again be divided into subarrays
 - Determine the array dependencies:
 - The border elements require information from a neighboring task, which means there must be communication between the tasks
 - The interior elements are independent of each other, requiring no communication between tasks
 - The Master process sends information to the workers and waits for results
 - THe Worker process calculates the solution within the specified number of steps and communicates with other tasks when necessary



Heat Calculation as SPMD

Heat Calculation Solution

```
find out if I am MASTER or WORKER
```

```
inf I am MASTER
    initialize array
        send each WORKER starting info and subarray
        receive results from each WORKER

else if I am WORKER
    receive from MASTER starting info and subarray

do t = 1, nsteps
    update time
    send neighbors my border info
    receive from neighbors their border info
    update my portion of the solution array
    end do
    send MASTER results

endif
```

1D Wave Equation

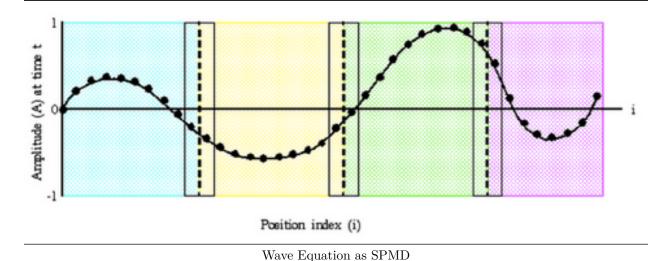
- 1. In this example, the amplitude along a uniform, vibrating string is calculated after a specified amount of time has elapsed
- 2. The calculation involves:
 - the amplitude on the y-axis
 - i as the position index along the x-axis
 - node points imposed along the string
 - update of the amplitude at discrete time steps

1D Wave Equation Problem

- 1. The equation to be solved is the 1-dimensional wave equation:
 - A(i,t+1) = (2.0 * A(i,t)) A(i,t-1) + (c * (A(i-1,t) (2.0 * A(i,t)) + A(i+1,t))) c is a constant
- 2. Note that the amplitude will depend on previous timesteps (t, t-1) and neighboring points (i-1,i+1)
 - Data dependencies will mean that a parallel solution will involve communications

1D Wave Equation Solution

- 1. Implement as an SPMD model
- 2. The entire amplitude aray is partitioned and distributed as subarrays to all tasks. Each task owns a portion of the total array
- 3. Load balancing: All points require equal work, so the points should be divided equally
- 4. A block decomposition would have the work partitioned into the number of tasks as chunks, allowing each task to own mostly contiguous data points
- 5. Communication need only occur on data borders. The larger the block size the less the communication



Conclusion

- 1. Additional References:
 - http://people.math.umass.edu/~johnston/PHI_WG_2014/OpenMPSlides_tamu_sc.pdf
 - http://www.nic.uoregon.edu/iwomp2005/iwomp2005_tutorial_openmp_rvdp.pdf
 - http://docs.oracle.com/cd/E19422-01/819-3694/
 - http://vuduc.org/cse6230/slides/cse6230-fa14-04-omp.pdf