Home Assignment 6

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Question 6.1

It is **possible** to reorganize the calculations so that in each iteration we did all of the calculations for each particle before proceeding to the next particle. But we need some modification.

We can **NOT** use the following pseudocode to implement the project:

```
for each timestep
   for each particle {
      Compute total force on particle;
      Find position and velocity of particle;
      Print position and velocity of particle;
}
```

Besides the for loop orders, we **DO** need some modification of the solver. The modification is that:

• Since the update of pos[][] will overwrite the old pos[][], the later computation of other force on the new pos[][] will result in wrong force. We should add an new variable old_pos[][] to record the previous position values and all other particles should use all the old_[][][] to compute forces.

The **NEW** pseudocode is as follow:

```
for each timestep
   for each particle {
       Compute total force on particle (using old position);
       Find new position and velocity of particle;
       Print new position and velocity of particle;
       Old position = New position
}
```

Question 6.6

Using the pesudocode:

```
# pragma omp parallel
    for each timestep {
        if (timestep output) {
            # pragma omp single
                Print positions and velocities of particles;
        }
        # pragma omp for
        for each particle q
                Compute total force on q;
        # pragma omp for
        for each particle q
                Compute position and velocity of q;
    }
```

- It will **DO** cause problems for the first for loop with nowait clause. Without the barrier at the end of the first loop, the computation inside the second loop could overwrite the value of posi[][]. And other threads may using the overwrote posi[][] for the computation inside the first loop and result in wrong force value.
- It will **DO** cause problems for the second *for* loop with *nowait* clause. The print operation need to wait for the finishing and updating of all the threads for the second loop.

Question 6.10

```
int n, thread count, i, chunksize;
double x[n], y[n], a;
...
# pragma omp parallel num threads(thread count) \
default(none) private(i) \
shared(x, y, a, n, thread count, chunksize)
{
# pragma omp for schedule(static, n/thread count)
    for (i = 0; i < n; i++) {
        x[i] = f(i); // f is a function
        y[i] = g(i); // g is a function
}
# pragma omp for schedule(static, chunksize)
    for (i = 0; i < n; i++)
        y[i] += a * x[i];
    }// omp parallel</pre>
```

We have n = 64, $thread_count = 2$, cache line size is 8 doubles, and each core has an L2 cache that can store 131,072 doubles. Since the scale L2 cache is much

larger than the scale of the array. There is no rewrite of L2 cache happens in this problem.

After the first loop the value in each L2 cache is like this:

```
Thread 0
x[] ||00-07||08-15||16-23||24-31||
y[] ||00-07||08-15||16-23||24-31||

Thread 1
x[] ||32-39||40-47||48-55||56-63||
y[] ||32-39||40-47||48-55||56-63||
```

$chunksize = n/thread_count$

The values needed for each thread for the second loop is:

```
Thread 0
x[] ||00-07||08-15||16-23||24-31||
y[] ||00-07||08-15||16-23||24-31||
Thread 1
x[] ||32-39||40-47||48-55||56-63||
y[] ||32-39||40-47||48-55||56-63||
```

After the second loop the value in each L2 cache is as following, the same as before:

```
Thread 0
x[] ||00-07||08-15||16-23||24-31||
y[] ||00-07||08-15||16-23||24-31||

Thread 1
x[] ||32-39||40-47||48-55||56-63||
y[] ||32-39||40-47||48-55||56-63||
```

So 0 L2 cache miss happens.

chunksize = 8

The values needed for each thread for the second loop is:

```
Thread 0
x[] ||00-07||16-23||32-39||48-55||
y[] ||00-07||16-23||32-39||48-55||
Thread 1
x[] ||08-15||24-31||40-47||56-63||
y[] ||08-15||24-31||40-47||56-63||
```

After the second loop the value in each L2 cache is as before following:

```
Thread 0
x[] ||00-07||08-15||16-23||24-31|| + ||32-39||48-55||
y[] ||00-07||08-15||16-23||24-31|| + ||32-39||48-55||

Thread 1
x[] ||32-39||40-47||48-55||56-63|| + ||08-15||24-31||
y[] ||32-39||40-47||48-55||56-63|| + ||08-15||24-31||
```

So **8 L2 cache misses** happen , which happen on reading x[32], x[48], y[48] for thread 0 and x[08], x[24], y[08], y[24] for thread 1.

Question 6.13

Using Figure 6.6 as a guide, sketch the communications that would be needed in an "obvious" MPI implementation of the reduced n-body solver if there were three processes, six particles, and the solver used a cyclic distribution of the particles.

Answer:

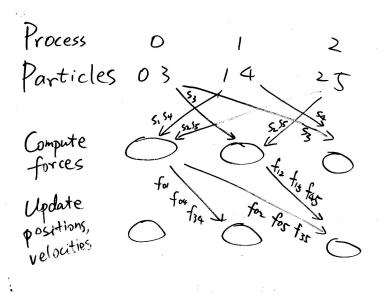


Figure 1: Communication in a possible MPI implementation of the reduced n-body solver with cyclic distribution.

Question 6.15

Assume we have p processors and the array length is N. The global index is i_{global} and the local index is i_{local} . Assuming that the number of processes evenly divides the number of elements in the global array.

6.15.a

The formula for a global index from a local index if the array has a block distribution.

$$i_{global} = local_rank \times \frac{N}{p} + i_{local}$$
 (1)

6.15.b

Then formula for a local index from a global index if the array has a block distribution

$$i_{local} = i_{global} - local_rank \times \frac{N}{p}$$
 (2)

6.15.c

Then formula for a global index from a local index if the array has a cyclic distribution

$$i_{global} = p \times i_{local} + local_rank \tag{3}$$

6.15.d

The formula for a local index from a global index if the array has cyclic distribution

$$i_{local} = \frac{i_{global} - local_rank}{p} \tag{4}$$

Question 6.17

6.17.a

The stack when maximum number of records happens (multiple possibilities):

3-2-NO_CITY-3-NO_CITY-3

Maximum number of records is 6.

6.17.b

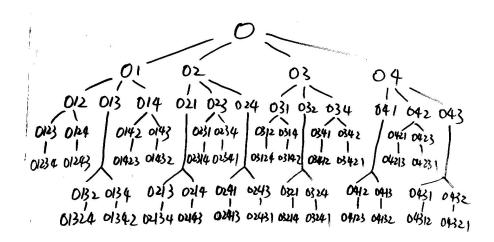


Figure 2: Search tree for five-city TSP.

6.17.c

The stack when maximum number of records happens (multiple possibilities):

 $4 - 3 - 2 - NO_CITY - 4 - 3 - NO_CITY - 4 - NO_CITY - 4$

Maximum number of records is 10.

6.17.d

For a n-city TSP:

Maximum number of records = $1 + 2 + \dots + (n-1) = \frac{n \times (n-1)}{2}$. (5)