# Advanced OpenMP Programming

Pangfeng Liu National Taiwan University

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- We will observe the timing when threads access variables at different locations.
- The intuition is that local variables are much easier to access than the global ones, and we need to confirm that.

## Compute Square Roots

- Write a program that computes the square root from 0 to 999999999 and assign the value to a variable v.
- We first observe the timing when we place the variable v at global area.

### Example 1: (assign.c)

```
#include <omp.h>
   #include <stdio.h>
   #include <math.h>
   #define N 999999999
5
   int main()
6
   {
7
     double v;
8
   #pragma omp parallel for
9
     for (int i = 1; i <= N; i++)</pre>
10
        v = sqrt(i);
11
     return 0;
12
   }
```

Private Variables
Thread Creation/Destroying
nowait
Critical Region
Double Ruffer

#### Demonstration

- Run the assign-uni program.
- Run the assign-omp program.

Private Variables
Thread Creation/Destroying
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Critical Region
Double Buffer

### Discussion

• Compare the execution time of assign-uni and assign-omp.



Private Variables
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Double Buffer

### Private Variable

 $\bullet$  We now declare v as private and observe the timing.

### Example 2: (assign-private.c)

```
#include <omp.h>
   #include <stdio.h>
   #include <math.h>
   #define N 999999999
5
   int main()
6
   {
7
     double v;
8
   #pragma omp parallel for private(v)
     for (int i = 1; i <= N; i++)</pre>
9
10
        v = sqrt(i);
11
     return 0;
12
   }
```

Private Variables
Thread Creation/Destroying
nowait
Critical Region
Double Buffer

### Demonstration

• Run the assign-private-omp program.

#### Discussion

- Compare the execution time of assign-private-omp with previous assign-uni and assign-omp.
- What is the reason for this performance difference?

Private Variables
Thread Creation/Destroying
nowait
Critical Region

# Heap Variable

• We now put v into the heap and observe the timing.



### Example 3: (assign-heap.c)

```
#include <omp.h>
   #include <stdio.h>
3
   #include <stdlib.h>
   #include <math.h>
5
   #define N 999999999
6
   int main()
   {
8
     double *v = malloc(sizeof(double));
9
   #pragma omp parallel for
10
     for (int i = 1; i <= N; i++)
11
       *v = sqrt(i);
12
     return 0;
13
   }
```

Private Variables
Thread Creation/Destroying
nowait
Critical Region
Double Buffer

#### Demonstration

• Run the assign-heap-omp program.

#### Discussion

- Compare the execution time of assign-heap-omp with previous programs.
- What is the reason for this performance difference?

# Prime Number Counting

- We want to count the number of prime numbers.
- We start with an array of numbers, and assuming that every number is a prime number.
- We start with the smallest prime number in the array, and mark every multiple of it as composite (non-prime number).
- We repeat this process until no new prime numbers are found.

#### Example 4: (prime.c)

```
#include <stdio.h>
   #include <math.h>
   #include <stdlib.h>
5
   #include <assert.h>
6
   #include "omp.h"
8
   #define N 40000000
9
10
   char notPrime[N];
11
   int nPrime = 0;
12
   char primes[N];
```

```
14
   int main(int argc, char *argv[])
15
   ₹
16
     assert(argc == 2);
17
     int n = atoi(argv[1]);
18
     int bound = round(sqrt(n));
19
20
     for (int i = 2; i <= bound; i++)</pre>
21
        if (!notPrime[i]) {
22
   #pragma omp parallel for
23
          for (int j = 2 * i; j < n; j += i)
24
            notPrime[j] = 1;
25
```

```
int nPrime = 0;
#pragma omp parallel for reduction(+ : nPrime)
  for (int i = 2; i < n; i++)
    if (notPrime[i] == 0)
       nPrime++;

  printf("number of prime is %d\n", nPrime);
  return 0;
}</pre>
```

#### **Variables**

- Array notPrime keeps track of the status of a number. If we know a number i is *not* prime, we set notPrime[i] to 1.
- The prime numbers will be kept in another array primes.
- The range to be tested (n) is given as a command line argument.

# Try All Possibilities

- Try all numbers from 2 to  $\sqrt{n}$ .
- If i is a prime number, mark the all multiple of i as *not* prime.
- It is obvious that the first for loop cannot be parallelized because of dependency, so we parallelize the second for loop.

#### Reduction

- The number of prime number can be obtained by counting the number of zeros in array notPrime.
- We use a reduction on the variable nPrime to simplify the process.

#### Demonstration

- Run the prime-uni program.
- Run the prime-omp program.

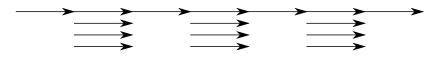
#### Discussion

- Compute the speedup of the previous prime counting program.
- Is there any optimization that can improve the performance?

# **Efficiency**

- The previous program parallelized the inner for loop only.
- The previous program will go through multi-threading every time a prime number is found.
- This incurs overheads of creating and destroying the threads.

# Spawn and Join



- Single thread for the outer loop.
- Multi-thread for the inner loop.

## parallel Once

- We would like to avoid the overheads in creating and destroying threads, so we put a parallel in front of the first for loop.
- The entire two level loop is run by all threads.
- Then we share the workload of the second for loop to improve performance, since most work is done in the second loop.

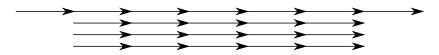
#### **Example 5: (prime-inner.c)**

```
13
   int main(int argc, char *argv[])
14
15
     assert(argc == 2);
16
     int n = atoi(argv[1]);
     int bound = round(sqrt(n));
17
18
19
   #pragma omp parallel
20
     for (int i = 2; i <= bound; i++)</pre>
21
        if (!notPrime[i])
22
   #pragma omp for
23
          for (int j = 2 * i; j < n; j += i)
24
            notPrime[j] = 1;
```

### **Efficient**

- All threads run the first loop, and share the second loop.
- It is OK for all threads to run the first for loop simultaneously, since they will synchronize at every second loop.
- Index variables i and j are private.

## Spawn and Join



- Multi-thread for both outer and inner loops.
- The work in the outer loop is duplicated, but it is more efficient than creating and joining threads.

#### Demonstration

- Run the prime-uni program.
- Run the prime-omp program.
- Run the prime-inner-omp program.

### Discussion

- Compare the execution time of all three prime counting programs.
- Is there any further optimization that can improve the performance?

### Synchronization

- A for pragma will synchronize all threads before leaving the for statement.
- If the for statement is still within the same parallel directive, then a barrier synchronization will do.
- If the for statement is at the end of the parallel directive, threads will be joined by the master thread and destroyed.

#### nowait

- Sometimes we do not wish the the threads to wait for each other.
- Those finish earlier can go on to the next statement to improve performance.
- We can only do this if the following statement does not depend on the previous statement.
- In this case we can use nowait clause.

Performance

Private Variables
Thread Creation/Destroying
nowait
Critical Region
Double Buffer

### nowait

 $1 \mid \mathtt{nowait}$ 



## Two loops

- We place two for directives with a parallel directive.
- The first for has an ascending workload, and the second loop has a descending workload.

#### Example 6: (2for.c)

```
int main(int argc, char *argv[])

{
    assert(argc == 3);
    omp_set_num_threads(atoi(argv[1]));
    int n = atoi(argv[2]);
    printf("# of proc = %d\n", omp_get_num_procs());
    printf("# of loop iterations = %d\n", n);
```

```
15
      double t = omp_get_wtime();
16
   #pragma omp parallel
17
18
   #pragma omp for
19
        for (int i = 0; i < n; i++)</pre>
20
          sleep(i);
21
   #pragma omp for
22
        for (int i = n - 1; i \ge 0; i--)
23
          sleep(i);
24
      }
25
      printf("time = %f\n", omp_get_wtime() - t);
```

#### Demonstration

 Run the 2for-omp program with 4 threads and 8 iterations, and observe the timing.

# Discussion

• Describe the the execution time of 2for-omp and make sure that it is reasonable.

Double Buffer

- In the previous program If the work of the second loop does not depend on the first loop, then we can let the threads go to the second directly.
- Since the first for has an ascending workload, and the second loop has a descending workload, if we let the threads to go to the second loop then the workload imbalance will be reduced.
- We only need to add a nowait clause at the first for directive.

#### Example 7: (2for-nowait.c)

```
15
      double t = omp_get_wtime();
16
   #pragma omp parallel
17
      {
18
   #pragma omp for nowait
19
        for (int i = 0; i < n; i++)</pre>
20
          sleep(i);
21
   #pragma omp for
22
        for (int i = n - 1; i >= 0; i--)
23
          sleep(i);
24
      }
25
      printf("time = %f\n", omp_get_wtime() - t);
```

### Demonstration

• Run the 2for-nowait-omp program with 4 threads and 8 iterations, and observe the timing.

# Discussion

• Describe the the execution time of 2for-omp and make sure that it is reasonable.

#### nowait

- We consider our previous prime number counting program.
- Previously the second marking for will synchronize before going back to the outside for.
- We would like to remove this synchronization, and let each threads to start the "prime" finding as soon as possible.
- This will not cause a race condition since the threads will still synchronize at the beginning of the next inner loop.

#### Example 8: (prime-inner-nowait.c) nowait

```
#pragma omp parallel
for (int i = 2; i <= bound; i++)
    if (!notPrime[i])
#pragma omp for nowait
for (int j = 2 * i; j < n; j += i)
    notPrime[j] = 1;</pre>
```

### Demonstration

• Run prime-inner-omp and prime-inner-nowait-omp and compare their execution time.

# Discussion

• Why can the nowait clause improve performance?

# Compute $\pi$

- We calculate  $\pi$  by integrating  $f(x) = \frac{4}{1+x^2}$ , where x is from 0 to 1.
- Divide the interval into N pieces, and assume the area in each interval is a trapezoid, then sum the area of these N trapezoids into a variable area.
- Note that in each interval you only need to compute  $\frac{4}{1+x^2}$  once.

#### Parallel Version

- We use x to denote the x coordinate, and area for the area in the integral.
- We parallelize the program by adding parallel for pragma and declare x as private.

#### Example 9: (pi.c) Compute $\pi$

```
#include <omp.h>
    #include <stdio.h>
    #define N 100000000
4
    int main()
5
    {
6
      double x;
7
      double area = 0.0;
8
      double t = omp_get_wtime();
9
    #pragma omp parallel for private(x)
10
      for (int i = 0; i < N; i++) {</pre>
11
        x = (i + 0.5) / N:
12
        area += 4.0/(1.0 + x * x):
13
14
      double pi = area / N;
15
      t = omp_get_wtime() - t;
16
      printf("execution time is %f\n", t);
      printf("pi = %f\n", pi);
17
18
      return 0;
19
```

# Demonstration

• Run pi-omp.

### Discussion

- Is the answer from the parallel version correct?
- Find out why the sequential version will not compile and fix it.
   Then compare the the execution time of these two programs.

#### **Atomic**

- We did not declare area private because it has the final global answer.
- As a result the operation on area must be atomic to avoid race condition.
- We simply use critical directive on the loop body to ensure atomic condition.

#### Example 10: (pi-critical.c) Compute $\pi$

```
9
      double x;
10
     double area = 0.0;
11
      double t = omp_get_wtime();
12
   #pragma omp parallel for private(x)
13
      for (int i = 0; i < N; i++)</pre>
14
   #pragma omp critical
15
16
          x = (i + 0.5) / N:
17
          area += 4.0 / (1.0 + x * x):
18
19
      double pi = area / N;
20
      t = omp_get_wtime() - t;
21
      printf("execution time is %f\n", t);
22
      printf("pi = %f\n", pi);
```

# Demonstration

• Run the pi-critical-omp program.

### Discussion

- Is the answer from the parallel version correct?
- Compare the the execution time of with the two previous programs.

### **Atomic**

- It appears that we do not need to make the entire loop body critical because x is already private.
- We now only use critical directive on the area summation.

#### Critical Section

- It is essential to reduce the size of the critical section,
- We want a thread to get through a critical section as soon as possible, so that other threads can get into the critical section as well.

#### Example 11: (pi-critical-small.c) Compute $\pi$

```
9
      double x;
10
     double area = 0.0;
11
      double t = omp_get_wtime();
12
   #pragma omp parallel for private(x)
13
      for (int i = 0; i < N; i++)</pre>
14
15
          x = (i + 0.5) / N:
16
   #pragma omp critical
17
          area += 4.0 / (1.0 + x * x):
18
19
     double pi = area / N;
20
      t = omp_get_wtime() - t;
21
      printf("execution time is %f\n", t);
22
      printf("pi = %f\n", pi);
```

### Demonstration

• Run the pi-critical-small-omp program.

### Discussion

- Is the answer from the parallel version correct?
- Compare the the execution time of with the three previous programs.

# Private Variables

- The performance improvement is very limited by a smaller critical section because it is not very different from the previous implementation.
  - Two statements v.s. one statement.
- If there are many statements in the loop body the benefit will more obvious.

## Critical Section

- The number of critical sections is enormous.
- We would like to remove these time consuming critical sections by letting all threads to work on its integrals by summing into its own area.
- The idea is to use a global array to store the individual area, then the master threads can add sum them up.

Double Buffer

# Array Implementation

- We need an array for threads to store its area.
- During each iteration a thread needs to call omp\_get\_thread\_num() to know where to store its area, which is a significant overheads.

### Example 12: (pi-array.c)

```
8
     double x:
9
     double area[MAXT] = \{0.0\};
10
     double t = omp_get_wtime();
11
   #pragma omp parallel for private(x)
12
     for (int i = 0; i < N; i++) {</pre>
13
       x = (i + 0.5) / N;
14
        area[omp_get_thread_num()] +=
15
          4.0 / (1.0 + x * x);
16
17
     t = omp_get_wtime() - t;
18
     double areaSum = 0.0;
19
     for (int i = 0; i < omp_get_num_procs(); i++)</pre>
20
       areaSum += area[i]:
21
     double pi = areaSum / N;
```

### Demonstration

• Run the pi-array-omp program.

## Discussion

- Is the answer from the parallel version correct?
- Compare the the execution time of with the previous programs.
- Is there any way to reduce the overheads in calling omp\_get\_thread\_num()?

#### Reduction

- We now use a reduction to compute the integral.
- The implementation is much cleaner and (hopefully) with better performance because it has been optimized by the OpenMP library.

#### Example 13: (pi-reduction.c) Compute $\pi$

```
7
     double x:
8
     double area = 0.0;
9
     double t = omp_get_wtime();
10
   #pragma omp parallel for private(x) \
11
     reduction(+ : area)
12
     for (int i = 0; i < N; i++) {</pre>
13
       x = (i + 0.5) / N:
14
        area += 4.0 / (1.0 + x * x):
     }
15
16
     double pi = area / N;
17
     t = omp_get_wtime() - t;
18
     printf("execution time is %f\n", t);
19
     printf("pi = %f\n", pi);
```

### Demonstration

• Run the pi-reduction-omp program.

### Discussion

- Is the answer from the parallel version correct?
- Compare the the execution time of with the previous programs.

# Game of Life

- A two-dimensional board with cells. A cell could either live or dead.
- The status of a cell evolves according to its status and the status of its eight neighbors.
- A dead cell with exactly three live neighbors becomes a live cell.
- A live cell with two or three live neighbors stays live, otherwise it becomes dead.

# Double Buffer

- It is intuitive that we keep the status of cells in a two dimensional array.
- Since the status of cell depends on the status of others, if we modify a cell directly it will affect the computation on other cells. This causes "race" condition during the update.
- Instead we use two arrays A and B to store the status of cells.
   In initially the status is in A.

### **Iterations**

- We repeat the following steps.
  - We set the cell status of B according to A in the first, third, etc. iterations.
  - We set the cell status of A according to B in the second, fourth, etc. iterations.

# Double Buffers

- We use two buffers A and B to avoid data inconsistency in updating the status of cells.
- We use a macro to compute the number of live neighboring cells – a live cell has 1 and a dead cell has 0.
- We pad the broad boundary so that we can use a single macro to compute the number of live neighboring cells

#### Example 14: (life.c)

```
#include <stdio.h>
   #include <stdlib.h>
4
   #include <omp.h>
5
6
   #define MAXN 4096
   #define SIDE (MAXN + 2)
8
   #define nLiveNeighbor(A, i, j) \
10
     A[i + 1][j] + A[i - 1][j] + A[i][j + 1] + 
11
     A[i][j-1] + A[i+1][j+1] + A[i+1][j-1] + 
12
     A[i - 1][j + 1] + A[i - 1][j - 1]
13
14
   char A[SIDE][SIDE];
15
   char B[SIDE][SIDE];
```

Private Variables
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# Print

• We use a simple printing routine to print the status of cells.

```
void print(char A[SIDE][SIDE], int n)
{
  for (int i = 1; i <= n; i++) {
    for (int j = 1; j <= n; j++)
       printf("%2d ", A[i][j]);
    printf("\n");
  }
}</pre>
```

# Input

- if the flag READINPUT is set the main program will read the size of the board, the number of generations, and the cell status from stdin.
- Otherwise it will generate a random input of size 4096 by 4096, and repeat for ten generations.

```
26
   int main()
27
   {
28
      int n, generation, cell;
29
   #ifdef READINPUT
30
      scanf("%d%d", &n, &generation);
31
      for (int i = 1; i <= n; i++)
32
        for (int j = 1; j <= n; j++) {
33
          scanf("%d", &cell);
34
          A[i][j] = cell;
35
36
   #else
37
     n = 4096;
38
      generation = 20;
39
      for (int i = 1; i <= n; i++)
40
        for (int j = 1; j <= n; j++)
41
          A[i][j] = rand() % 2;
42
   #endif
```

### Dead or Alive

- Depending on the generation the program will set the cell status of B with A, or in the other direction.
- A cell will be alive only if it is dead now and has three live neighbors, or it is live and it has two or three live neighbors now.
- We simply use a parallel for directive to distribute the workload.

```
int nln;
  for (int g = 0; g < generation; g++)</pre>
   if (g \% 2 == 0)
#pragma omp parallel for /* from A to B */
      for (int i = 1; i <= n; i++)</pre>
        for (int j = 1; j \le n; j++) {
          nln = nLiveNeighbor(A, i, j);
          B[i][j] = ((A[i][j] == 0 \&\& nln == 3) ||
            (A[i][j] == 1 && (nln == 2 || nln == 3)));
    else
#pragma omp parallel for /* from B to A */
      for (int i = 1; i <= n; i++)</pre>
        for (int j = 1; j \le n; j++) {
          nln = nLiveNeighbor(B, i, j);
          A[i][j] = ((B[i][j] == 0 \&\& nln == 3) ||
            (B[i][j] == 1 && (nln == 2 || nln == 3)));
        }
```

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# Final Status

• if the flag PRINT is set the main program will output the final cell status.

```
63
    #ifdef PRINT
64
      if (generation % 2 == 0)
65
        print(lifeA, n);
66
      else
67
        print(lifeB, n);
68
    #endif
69
      return 0;
70
    }
```

Private Variables
Thread Creation/Destroying
nowait
Critical Region
Double Buffer

### Demonstration

- Run the life-uni program.
- Run the life-omp program.

### Discussion

- Compute the speedup of the parallel program.
- What kind of scheduling policy was used?
- Is there any optimization that can improve the performance?