OpenMP Reduction Case Study: Trapezoid Integration Example

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Don’t do it this way!

const double A = 0.;
const double B = M_PI;
double dx = ( B - A ) / (float) ( numSubdivisions - 1 );
double sum = ( Function(A) + Function(B) ) / 2.;
omp_set_num_threads( numThreads );
#pragma omp parallel for default(none), shared(dx,sum)
for( int i = 1; i < numSubdivisions - 1; i++ )
{
    double x = A + dx * (float) i;
    double f = Function( x );
    sum += f;
}
sum *= dx;

• There is no guarantee when each thread will execute this line
• There is not even a guarantee that each thread will finish this line before some other thread interrupts it.

Assembly code:

Load sum
Add f
Store sum

What if the scheduler decides to switch threads right here?

Find the area under the curve $y = \sin(x)$ for $0 \leq x \leq \pi$

using the Trapezoid Rule

Exact answer: $\int_0^\pi \sin(x)dx = -\cos x \bigg|_0^\pi = 2.0$

The answer should be 2.0 exactly, but in 30 trials, it’s not even close. And, the answers aren’t even consistent. How do we fix this?

0.469635 0.398893
0.517984 0.446419
0.438668 0.431204
0.437553 0.501783
0.398761 0.334996
0.506564 0.484124
0.489211 0.506362
0.584610 0.448226
0.476670 0.434737
0.530668 0.44919
0.500062 0.442342
0.672593 0.548837
0.411158 0.363092
0.408718 0.544778
0.523448 0.356299
The answer should be 2.0 exactly, but in 30 trials, it’s not even close. And, the answers aren’t even consistent. How do we fix this?

There are Three Ways to Make the Summing Work Correctly:

#1: Atomic

```c
#pragma omp parallel for shared(dx)
for( int i = 0; i < numSubdivisions; i++ )
{
    double x = A + dx * (float) i;
    double f = Function( x );
    #pragma omp atomic
    sum += f;
}
```

- More lightweight than critical (#2)
- Uses a hardware instruction CMPXCHG (compare-and-exchange)
- Can only handle these operations:
  - x++, ++x, x--, --x
  - x op= expr , x = x op expr, x = expr op x
  - where op is one of: +, -, *, /, &, |, ^, <<, >>

#2: Critical

```c
#pragma omp parallel for shared(dx)
for( int i = 0; i < numSubdivisions; i++ )
{
    double x = A + dx * (float) i;
    double f = Function( x );
    #pragma omp critical
    sum += f;
}
```

- More heavyweight than atomic (#1)
- Allows only one thread at a time to enter this block of code (similar to a mutex)
- Can have any operations you want in this block of code

#3: Reduction

```c
#pragma omp parallel for shared(dx), reduction(+:sum)
for( int i = 0; i < numSubdivisions; i++ )
{
    double x = A + dx * (float) i;
    double f = Function( x );
    sum += f;
}
```

- OpenMP creates code to make this as fast as possible
- Reduction operators can be: +, - , *, & , | , ^ , && , || , max , min
Speed of Reduction vs. Atomic vs. Critical

(up = faster)

So, do it this way!

const double A = 0.;
const double B = M_PI;
double dx = (B - A) / (float)(numSubdivisions - 1);
omp_set_num_threads(numThreads);
double sum = (Function(A) + Function(B)) / 2.;
#pragma omp parallel for default(none), shared(dx), reduction(+:sum)
for(int i = 1; i < numSubdivisions - 1; i++)
{
    double x = A + dx * (float)i;
    double f = Function(x);
    sum += f;
}
sum *= dx;

O(N) vs. O(log₂N)

2. Reduction automatically creates a binary tree structure, like this, to add the N running sums in log₂N time instead N time.

Serial addition:
Adding 8 numbers requires 7 steps
Adding 1,048,576 (1M) numbers requires 1,048,575 steps

Parallel addition:
Adding 8 numbers requires 3 steps
Adding 1,048,576 (1M) numbers requires 20 steps

Two Reasons Why Reduction is so Much Better in this Case

1. Reduction secretly creates a temporary private variable for each thread’s running sum. Each thread adding into its own running sum doesn’t interfere with any other thread adding into its own running sum, and so threads don’t need to slow down to get out of the way of each other.

2. Reduction automatically creates a binary tree structure, like this, to add the N running sums in log₂N time instead N time.
If You Understand NCAA Basketball Brackets, You Understand Power-of-Two Reduction

Computer Graphics

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Why Not Do Reduction by Creating Your Own sums Array, one for each Thread, Like This?

```c
float *sums = new float [omp_get_num_threads()];
for (int i = 0; i < omp_get_num_threads(); i++)
    sums[i] = 0.;

#pragma omp parallel for private(myPartialSum),shared(sums)
for (int i = 0; i < N; i++)
    {
        myPartialSum = ...  
        sums[omp_get_thread_num()] += myPartialSum;
    }

float sum = 0.;
for (int i = 0; i < omp_get_num_threads(); i++)
    sum += sums[i];

delete [] sums;
```

- This seems perfectly reasonable, it works, and it gets rid of the problem of multiple threads trying to write into the same reduction variable.

- The reason we don’t do this is that this method provokes a problem called False Sharing. We will get to that when we discuss caching.