

# Embedded computing for scientific and industrial imaging applications

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Lecture 14 - OpenMP, critical sections, parallel for loops

# OpenMP

“Open Specifications for MultiProcessing”

Standard for shared memory parallel programming.  
For shared memory computers, such as multi-core.

Can be used with Fortran (77/90/95/2003), C and C++.

Complete specifications at <http://www.openmp.org>

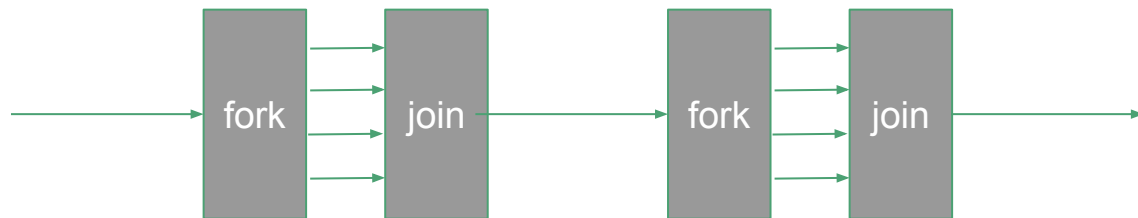
# References

- <http://www.openmp.org>
- <http://www.openmp.org/wp/resources/>
- B. Chapman, G. Jost, R. van der Pas, Using OpenMP: Portable Shared Memory Parallel Programming, MIT Press, 2007.
- R. Chandra, L. Dagum, et. al., Parallel Programming in OpenMP, Academic Press, 2001.
- <https://github.com/OpenMP>

# OpenMP — Basic Idea

Explicit programmer control of parallelization using fork-join model of parallel execution

- all OpenMP programs begin as single process, the master thread, which executes until a parallel region construct encountered
- FORK: master thread creates team of parallel threads
- JOIN: When threads complete statements in parallel region construct they synchronize and terminate, leaving only the master thread.



# OpenMP — Basic Idea

- **Rule of thumb:** One thread per processor (or core),
- User inserts **compiler directives** telling compiler how statements are to be executed
  - which parts are parallel
  - how to assign code in parallel regions to threads
  - what data is private (local) to threads
- Compiler generates explicit threaded code
- Dependencies in parallel parts require synchronization between threads
- User's job to **remove dependencies** in parallel parts or use **synchronization**.  
(Tools exist to look for **race conditions**.)

# OpenMP - compilers

<http://openmp.org/wp/openmp-compilers/>

- MS - [OpenMP in Visual C++](#)

/openmp (Enable OpenMP 2.0 Support)

- Project-> Properties -> C/C++ -> Language
- Change OpenMP Support to Yes(/openmp)

- GNU - gcc

Free and open source

From GCC 6.1, OpenMP 4.5 is fully supported in C and C++.

Compile with -fopenmp to enable OpenMP.

# OpenMP compiler directives

Uses **compiler directives** that start with **#pragma** (!\$ in fortran.)

These look like comments but are recognized when compiled with the flag **/openmp** (VS), **-fopenmp**(gcc).

## OpenMP statements:

OpenMP compiler directives, e.g.

```
#pragma omp parallel do
```

Calls to OpenMP library routines:

```
#include <omp.h>          //need this header  
omp_set_num_threads(2)
```

# OpenMP directives

```
#pragma omp directive [clause ...]  
    if (scalar_expression)  
    private (list)  
    shared (list)  
    default (shared | none)  
    firstprivate (list)  
    reduction (operator: list)  
    copyin (list)  
    num_threads (integer-expression)
```



# A few OpenMP directives

```
#pragma omp parallel [clause]
{      // block of code
}

#pragma omp parallel do [clause]
{      // do loop
}

#pragma omp barrier
// wait until all threads arrive
```

Several others we'll see later...

# OpenMP

API also provides for (but implementation may not support):

- Nested parallelism (parallel constructs inside other parallel constructs)
- Dynamically altering number of threads in different parallel regions

The standard says nothing about parallel I/O.

OpenMP provides "relaxed-consistency" view of memory.

Threads can cache their data and are not required to maintain exact consistency with real memory all the time.

```
#pragma omp flush
```

can be used as a **memory fence** at a point where all threads must have consistent view of memory

# OpenMP test code

```
#include <omp.h>
#include <stdio.h>

int main()
{
    int thread_num;

    omp_set_num_threads(2);
    printf("Testing openmp ...\n");
    #pragma omp parallel
    {
        #pragma omp critical
        {
            thread_num = omp_get_thread_num();
            printf("This thread = %d\n", thread_num);
        }
    }
    return 0;
}
```

# OpenMP test code output

Compiled with OpenMP: from \$CSE6000/codes/OpenMP/Examples/test/test.c

```
$cl.exe /openmp test.c
```

```
$test.exe
```

```
Testing openmp ...
```

```
This thread = 0
```

```
This thread = 1
```

(or threads might print in the other order!)

# OpenMP test code

```
//Specify number of threads to use:  
omp_set_num_threads(2)
```

Can specify more threads than processors, but they won't execute in parallel.

The number of threads is determined by (in order):

- Evaluation of if clause of a directive  
(if evaluates to zero or false  $\Rightarrow$  serial execution)
- Setting the num\_threads clause
- the omp\_set\_num\_threads() library function
- the OMP\_NUM\_THREADS environment variable
- Implementation default

# OpenMP test code

```
#pragma omp parallel
{
    #pragma omp critical
    {
        thread_num = omp_get_thread_num();
        printf("This thread = %d\n", thread_num);
    } //end of omp critical
} //end of omp parallel
```

The `#pragma omp parallel` block **spawns two threads** and each one works independently, doing all instructions in block. Threads are destroyed at `} //end of omp parallel`

However, the statements are also in a `#pragma omp critical` block, which indicates that this section of the code can be executed by only one thread at a time, so in fact they are not done in parallel.

**So why do this?** The function `omp_get_thread_num()` returns a unique number for each thread and we want to print both of these.

# OpenMP test code

Incorrect code without critical section:

```
#pragma omp parallel
{
    thread_num = omp_get_thread_num();
    printf("This thread = %d\n", thread_num);
}
```

Why not do these in parallel?

1. If the prints are done simultaneously they may come out **garbled** (characters of one interspersed in the other).
2. thread\_num is a **shared variable**. If this were not in a critical section, the following would be possible:

Thread 0 executes function, sets thread\_num=0

Thread 1 executes function, sets thread\_num=1

Thread 0 executes print statement: "This thread = 1"

Thread 1 executes print statement: "This thread = 1"

There is a **data race** or **race condition**.

# OpenMP test code

Could change to add a private clause:

```
#pragma omp parallel
{
    #pragma omp critical
    {
        thread_num = omp_get_thread_num();
        printf("This thread = %d\n", thread_num);
    } //end of omp critical
} //end of omp parallel
```

Then each thread has it's own version of the `thread_num` variable.



# OpenMP parallel for loops

```
#pragma omp parallel for
for(i=0; i<N; i++)
{
    //do stuff for each i
}
```

indicates that the for loop can be done in parallel.

## Requires:

- what's done for each value of *i* is independent of others

- Different values of *i* can be done in any order.

The iteration variable *i* is **private** to the thread: each thread has its own version.

By default, all other variables are **shared** between threads unless specified otherwise.

# OpenMP parallel for loops

This code fills a vector  $y$  with function values that take a bit of time to compute:

```
dx = 1.0 / (n+1.0);

#pragma omp parallel for private(x)
for(i = 1; i <= n; i++)
{
    x = i*dx;
    y[i-1] = exp(x)*cos(x)*sin(x)*sqrt(5*x+6.0);
}
```

Elapsed time for  $n = 10^8$ , without OpenMP: about 3.3 sec.

Elapsed time using OpenMP on 2 processors: about 1.9 sec.

# OpenMP parallel for loops

This code is **not correct**:

```
#pragma omp parallel for
for(i = 1; i <= n; i++)
{
    x = i*dx;
    y[i-1] = exp(x)*cos(x)*sin(x)*sqrt(5*x+6.0);
}
```

By default, x is a shared variable.

Might happen that:

- Processor 0 sets x properly for one value of i,
- Processor 1 sets x properly for another value of i,
- Processor 0 uses x but is now incorrect.

# OpenMP parallel for loops

Correct version:

```
#pragma omp parallel for private(x)
for(i = 1; i <= n; i++)
{
    x = i*dx;
    y[i-1] = exp(x)*cos(x)*sin(x)*sqrt(5*x+6.0);
}
```

Now each thread has its own version of x.

Iteration counter i is private by default.

Note that dx, n, y are shared by default. **OK because:**

- dx, n are used but not changed,
- y is changed, but independently for each i

# OpenMP parallel for loops

Incorrect code:

```
dx = 1.0 / (n + 1.0)
#pragma omp parallel for private(x, dx)
for(i = 1; i <= n; i++)
{
    x = i*dx;
    y[i-1] = exp(x)*cos(x)*sin(x)*sqrt(5*x+6.0);
}
```

Specifying dx private won't work here.

This will create a private variable dx for each thread but it will be **uninitialized**.

Will run but give garbage.

# OpenMP parallel for loops

Could fix with:

```
dx = 1.0 / (n + 1.0)
#pragma omp parallel for firstprivate(dx)
for(i = 1; i <= n; i++)
{
    x = i*dx;
    y[i-1] = exp(x)*cos(x)*sin(x)*sqrt(5*x+6.0);
}
```

The `firstprivate` clause creates private variables and initializes to the value from the master thread prior to the loop.

There is also a `lastprivate` clause to indicate that the last value computed by a thread (for  $i = n$ ) should be copied to the master thread's copy for continued execution.

# OpenMP parallel for loops

from \$CSE6000/codes/OpenMP/private1.c

```
n = 7;
y = 2.0;
#pragma omp parallel for firstprivate(y) lastprivate(y)
for(i = 1; i <= n; i++)
{
    y = y + 10.0;
    x[i - 1] = y;
    #pragma omp critical
    {
        printf("i = %d, x[i - 1] = %f\n", i, x[i - 1]);
    }
}
printf("At end y = %f\n", y);
```

# OpenMP parallel for loops

Run with 2 threads: The 7 values of  $i$  will be split up, perhaps

$i = 1, 2, 3, 4$  executed by thread 0,

$i = 5, 6, 7$  executed by thread 1.

Thread 0's private  $y$  will be updated 4 times,  $2 \rightarrow 12 \rightarrow 22 \rightarrow 32 \rightarrow 42$

Thread 1's private  $y$  will be updated 3 times,  $2 \rightarrow 12 \rightarrow 22 \rightarrow 32$

might produce:

$i = 1, x[i - 1] = 12.000000$

$i = 5, x[i - 1] = 12.000000$

$i = 2, x[i - 1] = 22.000000$

$i = 6, x[i - 1] = 22.000000$

$i = 3, x[i - 1] = 32.000000$

$i = 7, x[i - 1] = 32.000000$

$i = 4, x[i - 1] = 42.000000$

At end  $y = 32.000000$

Order might be different but final  $y$  will be from  $i = 7$ .



# OpenMP synchronization

```
#pragma omp parallel for
for(i = 1; i <= n; i++)
{
    /do stuff for each i
}
```

There is an **implicit barrier** at the end of the loop.

The master thread will not continue until all threads have finished with their subset of  $1, 2, \dots, n$ .

**Except if :**

```
#pragma omp parallel for nowait
```

# Conditional clause

Loop overhead may not be worthwhile for short loops.

(Multi-thread version may run slower than sequential)

Can use conditional clause:

```
#pragma omp parallel for if (n > 1000)
for(i = 1; i <= n; i++)
{
    // do stuff
}
```

If  $n \leq 1000$  then no threads are created,  
master thread executes loop sequentially

# Nested loops

```
#pragma omp parallel for private(i)
for(j = 0; j < m; j++)
{
    for(i = 0; i < n; i++)
    {
        a[j * n + i] = 0.0;
    }
}
```

The loop on  $j$  is split up between threads.

The thread handling  $j=0$  does the entire loop on  $i$ ,  
sets  $a[0 * n + 0]$ ,  $a[0 * n + 1]$ , ...,  $a[0 * n + n]$ .

**Note:** The loop iterator  $i$  must be declared **private**!  
 $j$  is private by default,  $i$  is shared by default.

# Nested loops - Which is better? (assume $m \approx n$ )

```
#pragma omp parallel for private(i)
for(j = 0; j < m; j++)
{
    for(i = 0; i < n; i++)
    {
        a[j * n + i] = 0.0;
    }
}
```

```
for(j = 0; j < m; j++)
{
    #pragma omp parallel for
    for(i = 0; i < n; i++)
    {
        a[j * n + i] = 0.0;
    }
}
```

# Nested loops - Which is better? (assume $m \approx n$ )

```
#pragma omp parallel for private(i)
for(j = 0; j < m; j++)
{
    for(i = 0; i < n; i++)
    {
        a[j * n + i] = 0.0;
    }
}
```

```
for(j = 0; j < m; j++)
{
    #pragma omp parallel for
    for(i = 0; i < n; i++)
    {
        a[j * n + i] = 0.0;
    }
}
```

The first has less overhead: Thread created only once.

The second has more overhead: Thread created  $m$  times.

# Nested loops

Incorrect code for replicating first column:

```
#pragma omp parallel for private(j)
for(i = 1; i < n; i++)
{
    for(j = 0; j < m; j++)
    {
        a[j * n + i] = a[j * n + i - 1];
    }
}
```

Corrected: (j's can be done in any order, i's cannot)

```
#pragma omp parallel for private(i)
for(j = 0; j < m; j++)
{
    for (i = 1; i < n; i++)
    {
        a[j * n + i] = a[j * n + i - 1];
    }
}
```

# Reductions

**Incorrect code** for computing  $\|x\|_1 = \sum_i |x_i|$

```
norm = 0.0;
#pragma omp parallel for
for (i = 0; i < n; i++)
{
    norm = norm + fabs(x[i]);
}
```

There is a **race condition**: each thread is updating same shared variable norm.

**Correct code:**

```
#pragma omp parallel for reduction(+ : norm)
for (i = 0; i < n; i++)
{
    norm = norm + fabs(x[i]);
}
```

A reduction reduces an array of numbers to a single value.

# Reductions

A more complicated way to do this:

```
#pragma omp parallel private(mysum) shared(norm)
{
    mysum = 0;
    #pragma omp for
    for(i = 0; i < n; i++)
    {
        mysum = mysum + fabs(x[i]);
    }
    #pragma omp critical
    {
        norm = norm + mysum;
    }
}
```



# Some other reductions

Can do reductions using +, -, \*, &, |, &&, ||

<https://msdn.microsoft.com/en-us/library/2etkydkz.aspx>

General form:

```
#pragma omp parallel for reduction(operator : list)
```

Example with max:

```
int any_zero = 0;
#pragma omp parallel for reduction(||: any_zero)
for(i = 0; i < n; i++)
{
    any_zero = any_zero || (x[i] == 0.0);
}
printf("any zero in x, (1 is true, 0 is false) : %d\n", any_zero);
```