#### Code of the Day

```
void task1() {
  // Lock mutex1 first, then mutex2
  std::lock_guard<std::mutex> lock1(mutex1);
  std::cout << "Task 1 locked mutex1\n";

  // Simulate some work
  std::this_thread::sleep_for(std::chrono::milliseconds(1));

  std::lock_guard<std::mutex> lock2(mutex2);
  std::cout << "Task 1 locked mutex2\n";
}</pre>
```

```
int main() {
  std::thread t1(task1);
  std::thread t2(task2);
  t1.join();
  t2.join();
  return 0;
}
```

```
void task2() {
  // Lock mutex2 first, then mutex1
  std::lock_guard<std::mutex> lock2(mutex2);
  std::cout << "Task 2 locked mutex2\n";

  // Simulate some work
  std::this_thread::sleep_for(std::chrono::milliseconds(1));

  std::lock_guard<std::mutex> lock1(mutex1);
  std::cout << "Task 2 locked mutex1\n";
}</pre>
```

What problem can you see from this code?

Will the problem always occur?

# ECE 455

GPU Algorithm and System Design

[Fall 2025]

OpenMP Basics 09/22/2025

#### Before we get started...

- Quick overview, things discussed last time
  - Thread-level parallelism vs process-level parallelism
    - Process: An isolated virtual memory space that wraps the execution of your program
    - Thread: A basic execution unit living inside a process to run your program, starting with "main" function
  - Parallel programming using C++ thread (std::thread, std::mutex, std::condition\_variable)
- Purpose of today's lecture:
  - Introduce the basic usage of OpenMP
  - Use OpenMP to parallelize a loop of independent work
- Miscellaneous
  - Lab Assignment #2 due on Friday 9/26 at 23:59 PM

# Multi-core parallel computing with OpenMP

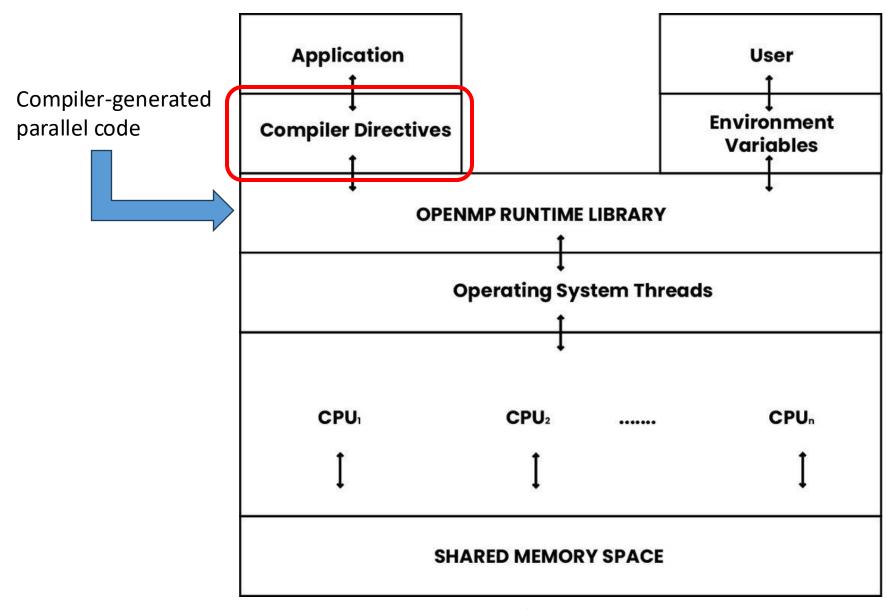
# OpenMP: <a href="https://www.openmp.org/">https://www.openmp.org/</a>

- OpenMP is an Application Programming Interface (API) that supports shared-memory multithreaded programming in C, C++, and Fortran
- Unlike functional-styled programming model like std::thread, OpenMP provides compiler directives that instruct the compiler to insert parallel code
  - aka directive-based programming model counting on compiler to generate parallel code without explicitly managing threads which can sometimes be very massy
- OpenMP has two major advantages

- OpenMP.
- Ease of use: Just tell compiler what and how to parallelize your code
- https://www.openmp.org/

Portability: It's compiler's job to insert platform-dependent parallel code!

# Software Architecture of OpenMP



## History of OpenMP

- Misson statement: standardize directive-based high-level parallelism that is performant, productive and portable across different platforms
- Current spec is 5.2
  - Released in November 2021, <a href="http://www.openmp.org">http://www.openmp.org</a>
  - More than 600 Pages
- Not all compilers are equally up-to-date with the OpenMP standard
  - GNU compiler supports up to OpenMP 4.5, but does not fully support GPU offloading
  - LLVM is the new kind in town, good support, up to features of 5.0 (partial features)
  - Microsoft compiler lagging behind (Version 2.0 as of Jan 2021)

#### Example: Hello World in OpenMP

```
#include <iostream>
#include <omp.h>
int main() {
  #pragma omp parallel num threads(4)
    int me = omp get thread num();
    int NT = omp_get_num_threads();
    std::cout << "Hello World. I'm thread " << me << " out of " << NT << ".\n";</pre>
    for (int i = 0; i < 2; i++)
      std::cout << "Iter:" << i << "\n";</pre>
  std::cout << "All done here..." << std::endl;</pre>
```

All OpenMP directives start with "#pragma" telling the compiler to generate parallel code based on the following instructions

```
Iter:1
OMP parallel region is executed by four threads
```

- Each thread executes the parallel region
- This type of programming approach is referred to single-source multiple-execution model

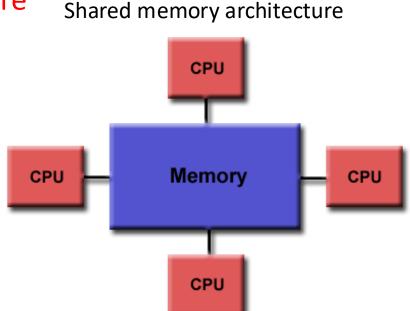
```
$ g++ test.cpp -fopenmp
$ ./a.out
Hello World. I'm thread 0 out of 4.
Iter:0
Iter:1
Hello World. I'm thread 2 out of 4.
Iter:0
Hello World. I'm thread 3 out of 4.
Iter:0
Iter:1
Hello World. I'm thread 1 out of 4.
Iter:0
Iter:1
All done here...
```

#### Example: Hello World in C++ Thread (for Comparison Purpose)

```
#include <iostream>
#include <thread>
                                                              Explicit thread management, creating threads,
int main() {
                                                             assigning a starting function (lambda here), and
  std::thread threads[4];
                                                                          joining created threads
  for (int t = 0; t < 4; t++) {
    threads[t] = std::thread([me=t, NT=4]({
       std::cout << "Hello World. I'm thread " << me << " out of " << NT << ".\n";</pre>
      for (int i = 0; i < 2; i++) {
                                                                $ g++ test.cpp
         std::cout << "Iter:" << i << "\n";</pre>
                                                                $ ./a.out
                                                                Hello World. I'm thread 0 out of 4.
                                                                Iter:0
    }));
                                                                Iter:1
                                                                Hello World. I'm thread 2 out of 4.
                                                                Iter:0
                                                                Iter:1
  for (int t = 0; t < 4; t++) {
                                                                Hello World. I'm thread 3 out of 4.
    threads[t].join();
                                                                Iter:0
                                                                Iter:1
                                                                Hello World. I'm thread 1 out of 4.
  std::cout << "All done here..." << std::endl;</pre>
                                                                Iter:0
                                                                Iter:1
                                                                All done here...
```

#### When to Use OpenMP?

- OpenMP: targets parallelism on a shared-memory architecture
  - Help you program thread-level parallelism (TLP) on multi-core CPUs without explicitly managing threads like using std::thread
    - You know, programming std::thread can be very tedious
  - Suitable for simple, commonly used parallel patterns
- OpenMP + CUDA: targets parallelism on the GPU
  - Modern OpenMP supports heterogeneous execution using GPU
  - Support only limited GPU-parallel primitives

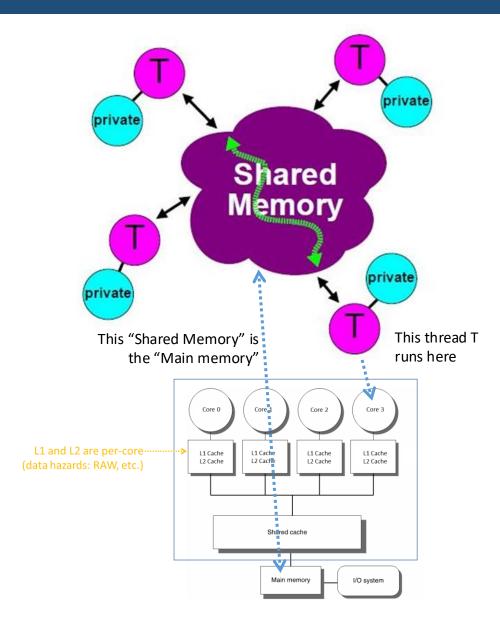


- OpenMP + MPI: targets parallelism on a cluster (distributed computing)
  - MPI handles communication among machines
  - OpenMP handles thread-level parallelism on a machine

#### When to NOT Use OpenMP

- Fine-grained parallelism
  - For tasks with very fine-grained parallelism, the overhead of creating and managing threads might outweigh its benefits
- Complex synchronization
  - If your application requires intricate synchronization patterns, OpenMP's built-in synchronization primitives might not be sufficient (e.g., OpenMP does not have things like std::condition\_variable)
- Dynamic parallelism pattern
  - For algorithms that depends on a lot of runtime variables to generate parallelism, OpenMP's directives may not be sufficient as the compiler has very limited information about what will happen at runtime

#### OpenMP Attributes [important slide]



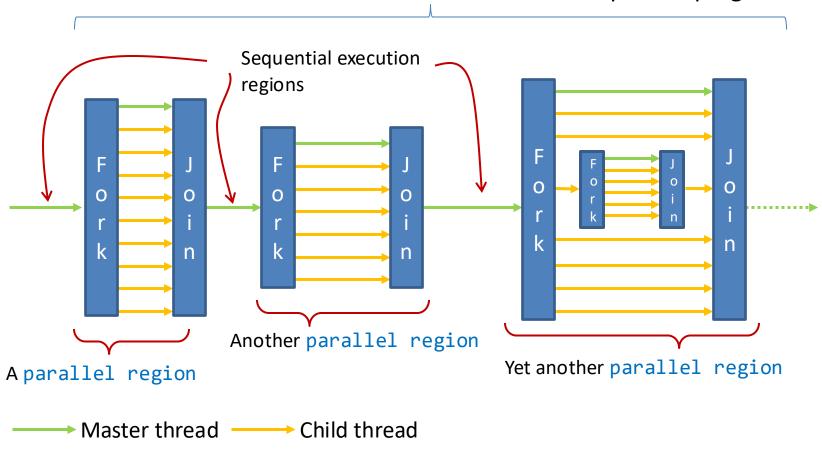
- Threads execute different units of work simultaneously
- Threads have private data inaccessible by other threads
- Threads have access to a pool of memory shared by all threads
- How data is accessed by threads is transparent to programmer
- Synchronization rules are defined by OpenMP

#### OpenMP is Based on the Fork-Join Model

- Master thread spawns a team of child threads based on the number of threads you request
  - Master thread will be part of the team member (Note: This is different from std::thread)
  - When work is done, all forked threads join together automatically (no explicit std::thread::join())
- You can ask for many threads in your OpenMP program, but it doesn't help you run faster if:
  - You go above the number of physical cores on your CPU
    - Ex: when you ask for more than 2x number of physical cores, typically your code slows down as the operating system gets busy servicing your many threads *oversubscription*
  - The parallelizable portion of your program can only support limited parallelism
    - Ex: two independent tasks won't benefit from running them using six threads

#### OpenMP Execution Model

A small time-window into the execution flow of an OpenMP program



### Compiling OpenMP Programs Using the Command Line

• Compiling OpenMP programs is easy as it has been integrated into mainstream compilers

```
• GCC: $ g++ -o integrate_omp integrate_omp.cpp -fopenmp
```

• Clang \$ clang++ -o integrate\_omp integrate\_omp.cpp -fopenmp

• ICC: \$ icpc -o integrate\_omp integrate\_omp.cpp -openmp

• MSVC: \$ cl /openmp integrate\_omp.cpp

### Example: Calculate Entries in a Table Sequentially

• Version 1: Calculate the std::sin values at different input n from 0 to 255

```
#include <omp.h>
constexpr auto PIE = 3.14159265358979323846;

int main()
{
    const int size = 256;
    double sinTable[size]; // sin table to be initialized

    for (int n = 0; n<size; ++n)
        sinTable[n] = std::sin(2 * PIE * n / size);
    // the table is now initialized
}</pre>
```

#### Example: Calculate Entries in a Table in Parallel w/ OpenMP

• Version 2: Calculate these 256 different std::sin values in parallel using OpenMP

```
#include <omp.h>
constexpr auto PIE = 3.14159265358979323846;
int main()
    const int size = 256;
    double sinTable[size]; // sin table to be initialized
    #pragma omp parallel for
    for (int n = 0; n<size; ++n)</pre>
        sinTable[n] = std::sin(2 * PIE * n / size);
    // the table is now initialized
```

### Example: Calculate Entries in a Table using SIMD w/ OpenMP

Version 3: Uses wide registers & vector operations to enable SIMD parallelism

```
#include <omp.h>
constexpr auto PIE = 3.14159265358979323846;
int main()
    const int size = 256;
    double sinTable[size]; // sin table to be initialized
    #pragma omp simd
    for (int n = 0; n<size; ++n)</pre>
        sinTable[n] = std::sin(2 * PIE * n / size);
    // the table is now initialized
```

#### Example: Calculate Entries in a Table using GPU w/ OpenMP

- Version 4: Offloading code to a GPU (not supported by all compilers)
  - The **sin** function must also exist on the target device

```
#include <omp.h>
constexpr auto PIE = 3.14159265358979323846;
int main()
{
    const int size = 256;
    double sinTable[size]; // sin table to be initialized

    #pragma omp target teams distribute parallel for map(from:sinTable[0:256])
    for (int n = 0; n<size; ++n)
        sinTable[n] = std::sin(2 * PIE * n / size);
    // the table is now initialized
}</pre>
```

Note: Not all platforms support this feature yet although it's pretty cool

#### A Few Syntax Details to Start with OpenMP

Picking up the OpenMP header file

```
#include <omp.h> (for C/C++)
```

- Most OpenMP constructs are compiler directives
  - C and C++: the directives start with the keyword pragma:
     #pragma omp construct [clause [clause]...]
  - Construct describes the parallelism type you want to perform
  - Clause further qualifies a directive's behavior

#### Why Compiler Directives?

- One can have the same code, with no modifications, run with or without OpenMP
  - It is compiler's job to figure out how to generate the parallel code based on the provided directives
- When OpenMP is disabled, all pragma statements will be treated as comments
  - No need to maintain separate codebases between sequential and parallel code
- Directives are picked up by the compiler only if instructed to do so
  - Example: Visual Studio you have to have the /openmp flag on

## Commonly used Constructs in OpenMP Compiler Directives

Directive	Description	
<u>atomic</u>	Specifies that a memory location will be updated atomically	
<u>barrier</u>	Synchronizes all threads in a team; all threads pause at the barrier, until all threads execute the barrier	
<u>critical</u>	Specifies that code is only executed on one thread at a time	
<u>flush</u>	Enforces that all threads have the same view of memory for all shared objects	
<u>for</u>	Causes the work done in a for loop inside a parallel region to be divided among threads	
<u>master</u>	Specifies that only the master thread should execute a section of the program	
ordered	Specifies that code under a parallelized for loop should be executed like a sequential loop	
parallel	Defines a parallel region, which is code that will be executed by multiple threads in parallel	
sections	Identifies code sections to be divided among all threads	
<u>single</u>	Indicates that a section of code should be executed on a single thread, not necessarily the master thread	
<u>threadprivate</u>	Specifies that a variable is private to a thread	

#### OpenMP Beyond Directives: User-Level Runtime Routines

- In addition to directives, OpenMP provide user-level API to modify OpenMP parameters:
  - Modify/check the number of threads

```
omp_get_max_threads()  # get the max number of threads
omp_[set|get]_num_threads()  # get the number of threads
omp_get_thread_num()  # get the id of this thread
```

Are we in a parallel region?

```
omp_in_parallel()
```

How many processors in the system?

```
omp_get_num_procs()
```

Explicit locks

```
omp_[set|unset]_lock()
```

Many more...

#### Example: Set/Get Number of Threads in a Parallel Region

```
// omp get num threads.cpp
// compile with: /openmp or -fopenmp
#include <iostream>
#include <omp.h>
int main() {
    std::cout << "Non parallel block, beginning of test: " << omp get num threads() << "\n";</pre>
    omp set num threads(2); // set the maximum number of threads to two
    std::cout << "Non parallel block, after omp set num threads call: " << omp get num threads() << "\n";</pre>
#pragma omp parallel
#pragma omp master
        std::cout << "Inside a parallel block: " << omp get num threads() << "\n";</pre>
    std::cout << "No parallel block here: " << omp get num threads() << std::endl;</pre>
                                                                                         $ g++ test.cpp -fopenmp
    // changed the number of threads to be used inside parallel block
                                                                                         $ ./a.out
    // using a compiler directive
                                                                                         Non parallel block, beginning of test: 1
#pragma omp parallel num threads(3)
                                                                                         Non parallel block, after omp set num threads call: 1
#pragma omp master
                                                                                         Inside a parallel block: 2
                                                                                         No parallel block here: 1
        std::cout << "Second parallel block: " << omp get num threads() << "\n";</pre>
                                                                                         Second parallel block: 3
                                                                                         Outside parallel block: 1
    std::cout << "Outside parallel block: " << omp get num threads() << std::endl;</pre>
```

#### Example: Get the Maximum Number of Threads

```
#include <cstdio>
                                                                                Query the maximum number of threads that
#include <omp.h>
                                                                                can be used to form a new team when
int main() {
                                                                                encountering a parallel region
    //omp set num threads(5);
    std::printf("I can go w/ this many threads:%d\n", omp get max threads());
#pragma omp parallel
#pragma omp master
       std::printf("Here's how many threads I use in this parallel region: %d\n", omp get num threads());
#pragma omp parallel num threads(3)
#pragma omp master
       std::printf("Max. number of threads: %d\n", omp get max threads());
       std::printf("Actual number of threads used in this other parallel region: %d\n", omp get num threads());
    std::printf("Here's the max number of threads at end:%d\n", omp_get_max_threads());
    return 0;
```

```
$ g++ -std=c++17 test.cpp -fopenmp
$ ./a.out
I can go w/ this many threads:8
Here's how many threads I use in this parallel region: 8
Max. number of threads: 8
Actual number of threads used in this other parallel region: 3
Here's the max number of threads at end:8
```

```
$ g++ -std=c++17 test.cpp -fopenmp
$ ./a.out Got this after uncommenting the first line in main() function

I can go w/ this many threads:5

Here's how many threads I use in this parallel region: 5

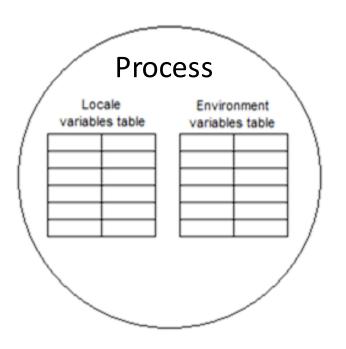
Max. number of threads: 5

Actual number of threads used in this other parallel region: 3

Here's the max number of threads at end:5
```

### Modify OpenMP Parameters using Environment Variables

- Environment Variables are *dynamic named values* that can affect a running processes
  - When operating system spawns a process to run a program, it creates a list of environment variables that store process-specific information in a key-value format



Key	Description of Value
PATH	(:)-separated list of directories in which your system looks for executable files
USER	Logged-in user name
HOME	Path to the logged-in user home directory
UID	Logged-in user ID
SHELL	The current logged-in shell
•••	•••

C++ <u>std::getenv</u> and <u>std::setenv</u> allow you to control environment variables in the current process

#### Two Identical Ways to Modify OpenMP Runtime Parameters

- In OpenMP, you can use both function call and environment variables to modify parameters
  - Example 1: controlling the number of threads the following two methods are equivalent
    - omp\_set\_num\_treads(8)
    - For the Bash shell:

```
export OMP_NUM_THREADS=8
```

- Example 2: controlling nested parallelism the following two methods are equivalent
  - omp set nested(1);
  - export OMP\_NESTED=1
- Both are interchangeable, but OpenMP function call trumps the environment variable setting
  - - → The OpenMP program will use 8 threads

# Commonly Used OpenMP Environment Variables

OMP CANCELLATION:	Set whether cancellation is activated
OMP DISPLAY ENV:	Show OpenMP version and environment variables
OMP DEFAULT DEVICE:	Set the device used in target regions
OMP DYNAMIC:	Dynamic adjustment of threads
OMP MAX ACTIVE LEVELS:	Set the maximum number of nested parallel regions
OMP MAX TASK PRIORITY:	Set the maximum task priority value
OMP NESTED:	Nested parallel regions
OMP NUM THREADS:	Specifies the number of threads to use
OMP PROC BIND:	Whether theads may be moved between CPUs
OMP PLACES:	Specifies on which CPUs the threads should be placed
OMP_STACKSIZE:	Set default thread stack size
OMP_SCHEDULE:	How threads are scheduled
OMP THREAD LIMIT:	Set the maximum number of threads
OMP WAIT POLICY:	How waiting threads are handled

#### OpenMP Programming is Supported by the Three Pillars

- Pillar 1: Compiler directives
  - Instruct compiler what and how to insert parallel code
  - A compiler that doesn't speak OpenMP simply ignores pragmas
    - Code runs just like before; i.e., sequentially
  - Directives have *clauses*, to further qualify a directive's behavior
- Pillar 2: User-level runtime function calls
  - Change OpenMP runtime behaviors programmatically
- Pillar 3: Environment variables
  - Another way of controlling OpenMP behavior without modifying the source code
  - Provides some of the support that the OpenMP functions provide

# OpenMP: Parallel For Construct

#### Parallel-for is the Most Widely used Construct in OpenMP

• Revisit: Calculate 256 different std::sin values in parallel using OpenMP parallel for

```
#include <omp.h>
constexpr auto PIE = 3.14159265358979323846;
int main()
    const int size = 256;
    double sinTable[size]; // sin table to be initialized
    #pragma omp parallel for
    for (int n = 0; n<size; ++n) {</pre>
        sinTable[n] = std::sin(2 * PIE * n / size);
```

- In a nutshell, the scope of an OpenMP directive is the structured block of code that it affects
  - A structured block is the code enclosed by an opening "{" and a closing "}" the only allowed "branches" to go out of the block are exit- or return-styled calls
  - This typically extends from the directive itself to the next immediate matching pair of curly braces "{}"

#### A "structured block"

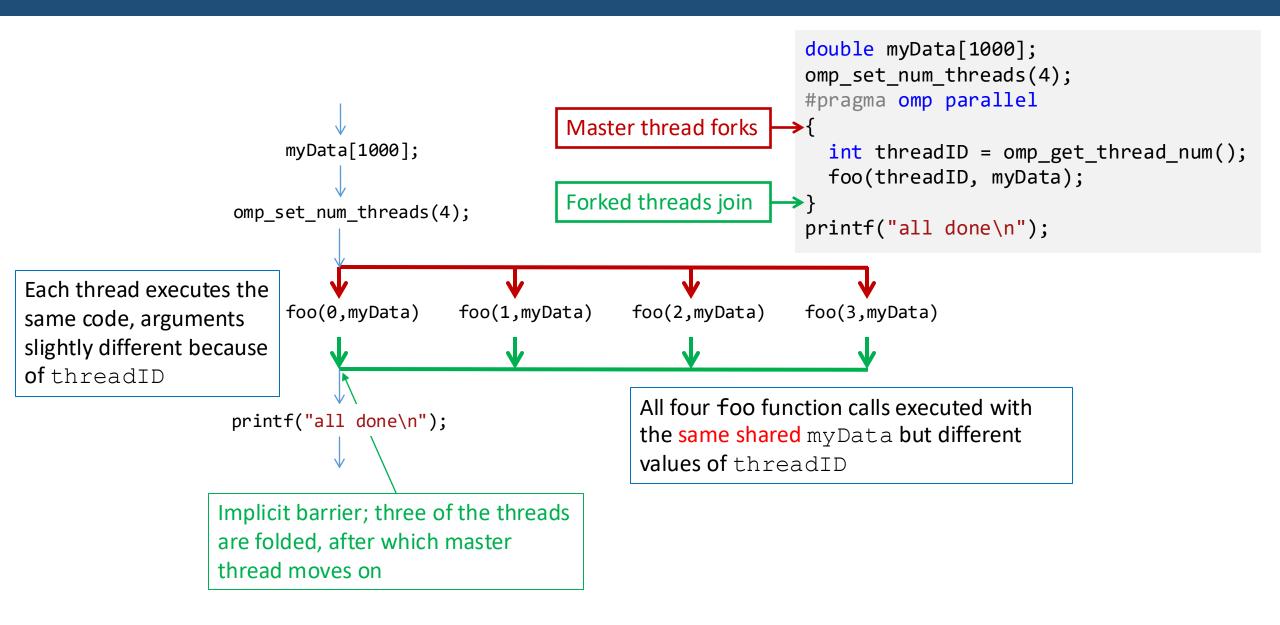
```
#pragma omp parallel
{
   int id = omp_get_thread_num();
   more: res[id] = do_big_job (id);
   if( not_conv(res[id]) )goto more;
}
printf ("I'm outside par. region!\n");
```

<u>IMPORTANT FACT</u>: There is an implicit barrier at the closing "}" where threads executing the block wait on each other to finish – the point at which forked threads will join.

#### Not a "structured block"

```
if (go_now()) goto more;
#pragma omp parallel
{
  int id = omp_get_thread_num();
more: res[id] = do_big_job(id);
  if ( conv (res[id]) ) goto done;
  goto more;
}
done: if (!really_done()) goto more;
```

## Execution Diagram when Entering a Structured Block



#### Example: Printing Messages from Multiple Threads

- The function whatsUpQuestionMark gets called by four parallel threads when the execution hits the parallel region
- Messages get outputted in a garbled way
  - The insertion operator "<<" is a function of std::cout</li>
  - Basically, it looks like the code below:

```
void whatsUpQuestionMark() {
  int me = omp_get_thread_num();
  printf("What's up?, asks thread ");
  printf("%d", me);
  printf("\n", me);
}
```

 Execution order of a function by different threads are non-deterministic

```
#include <omp.h>
#include <iostream>
void whatsUpQuestionMark()
  int me = omp get thread num();
  std::cout << "What's up?, asks thread " << me << "\n";</pre>
int main()
  #pragma omp parallel num threads(4)
      whatsUpQuestionMark();
  std::cout<< "all done...\n";</pre>
  return 0;
```

```
$ g++ test.cpp -fopenmp
$ ./a.out
What's up?, asks thread What's up?, asks thread What's up?, asks thread 03

1
What's up?, asks thread 2
all done...
```

#### What Happens Under the Hood in terms of Generated Code?

```
#pragma omp parallel num_threads(4)
{
    whatsUpQuestionMark();
}
```

- The OMP compiler generates some other code that takes care of the construct above
- When you say num\_threads(4), only three threads are created because the caller thread will be part of the team
- Since creating threads comes with certain cost, the OMP runtime uses a thread pool to reuse threads

```
void wrapper_for_the_structured_block()
{
    whatsUpQuestionMark();
}
```

```
std::thread threads[3];
for (int i = 0; i < 3; ++i) //note that i starts at 1
    threads[i] = std::thread(wrapper_for_the_structured_block);
wrapper_for_the_structured_block();
for (int i = 0; i < 3; ++i) //three threads folded; see Remarks below threads[i].join();</pre>
```

Note: The threads are not always created and joined from the scratch (std::thread and std::thread::join) for a parallel region; they might be recycled back into a thread pool to reduce the cost of repetitively creating threads and joining threads in parallel regions

#### A Short Side Trip: Timing an OpenMP Application

- double omp\_get\_wtime() returns a value in seconds of the time elapsed from some timepoint
  - This timepoint is a consistent point in the past guaranteed not to change during execution of program
- double omp\_get\_wtick() returns the number of seconds between clock ticks
  - This tick provides a higher resolution than wall time by directly measuring the duration in terms of clock ticks

```
#include <omp.h>
int main() {
   int const N = 100000;
   double dummy[N];
   double start = omp get wtime();
   for (int i = 0; i < N; i++) {
       int temp = std::rand();
       dummy[i] = 2.*temp / (std::pow(temp*temp, 1.5) + 0.2);
   double end = omp get wtime();
    std::printf("start = %.16g\n", start);
    std::printf("end = %.16g\n", end);
    std::printf("diff = %.16g\n", end - start);
    double wtick = omp get wtick();
    std::printf("wtick = %.16g\n", wtick);
    std::printf("1/wtick = %.16g\n", 1.0 / wtick);
```

```
start = 1528033.489782439
end = 1528033.510552168
diff = 0.02076972858048975
wtick = 3.011764252346089e-07
1/wtick = 3320313
Press any key to continue . . . _
```

#### Nested Parallelism

- Motivation: You have a parallel region that runs in parallel by, say four threads, and inside that parallel region, there is another code region that might also run in parallel
  - Ex: Multiplying two matrices requires looping through each row and column of the two matrices to calculate dot product

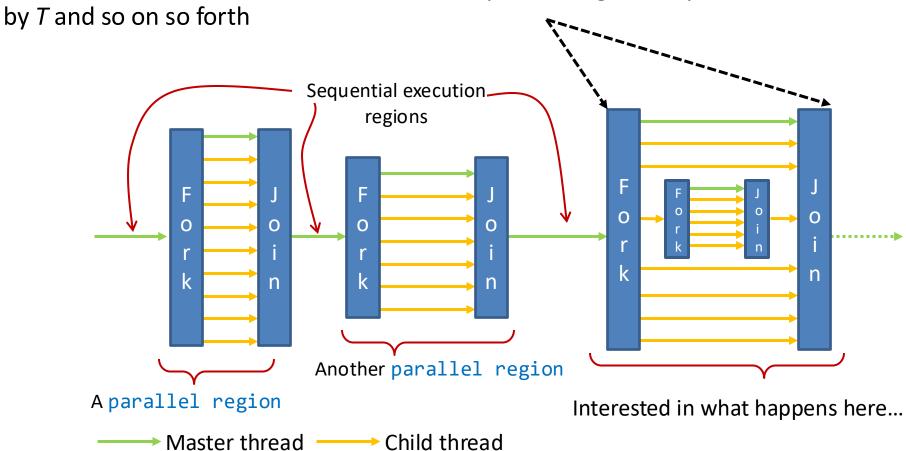
```
void MultiplyMatrices(int nCount, double **matrixA,
                double **matrixB, double **matrixC)
   int i, j, k;
   for (i = 0; i < nCount; i++)
       for (j = 0; j < nCount; j++)
           matrixC[i][j]=0;
            for (k = 0; k < nCount; k++)
               matrixC[i][j] +=
                    matrixA[i][k]*matrixB[k][j];
```

How can we parallelize nested loops using OpenMP?

#### OpenMP Supports Nested Parallelism

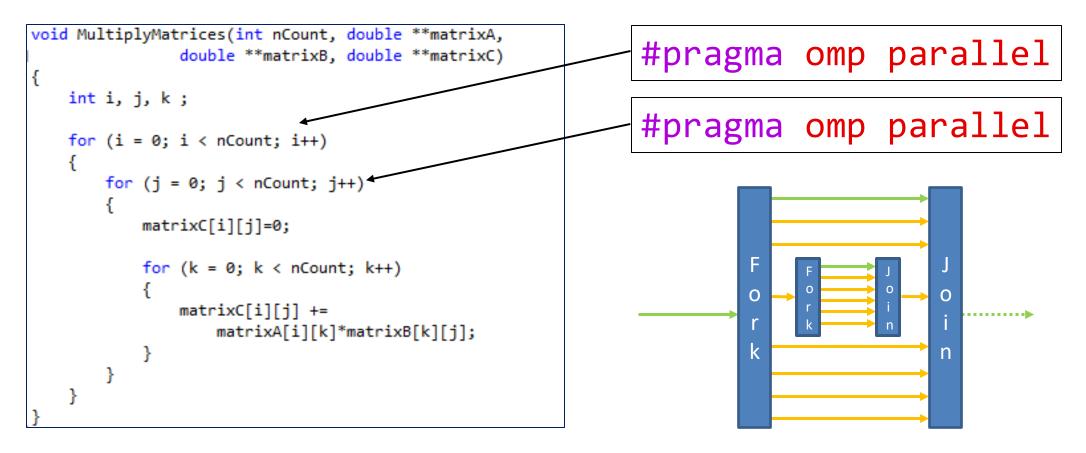
OpenMP parallel construct can be defined in a nested fashion

• When a thread T on a team of threads hits a parallel region, it spawns another teams of threads mastered



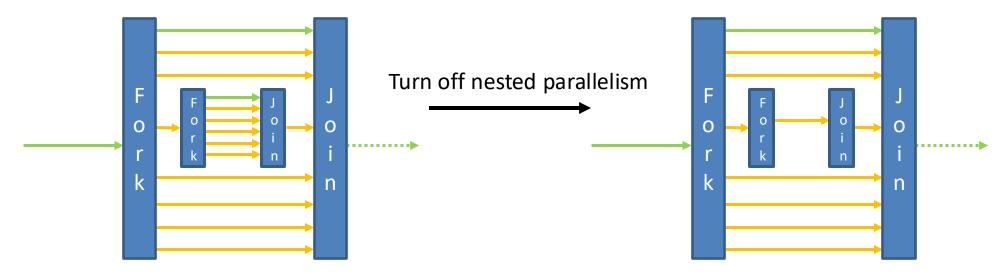
#### Nested Parallelism Revisited

- Nested parallelism is very easy to achieve using OpenMP simply add another parallel-for
  - One to parallelize the first for loop of i
  - Another to parallelize the second for loop of j



#### Nested Parallelism is Not a Silver Bullet ...

- Nested parallelism may not always take effect as you wish
  - Why? Nested parallelism is hard to implement efficiently depends a lot of application workloads
    - Some parts of the code with nested parallelism is ok
    - Some other parts of the code are better off without nested parallelism
- API to control nested parallelism behavior: omp\_set\_nested()
  - Alternatively, you can use the OMP\_NESTED environment variable to globally turn on/off nested parallelism
  - When you turn off nested parallelism, a parallel thread encountering a new parallel region will not be able to spawn new parallel threads that new parallel region will be executed by thread *T* only



#### Control the Behavior of Nested Parallelism

- Relevant API elements
  - omp\_set\_nested(...)
  - omp\_get\_nested(...)
  - omp\_set\_dynamic(...)
  - omp get dynamic(...)

Enables or disables nested parallelism

Control the number of threads in nested parallel regions

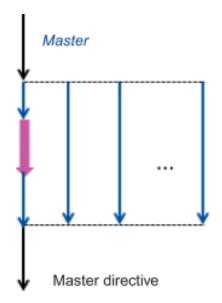
- If argument is nonzero, the runtime library can adjust the number of threads
- If argument is zero, the runtime library cannot dynamically adjust the number of threads

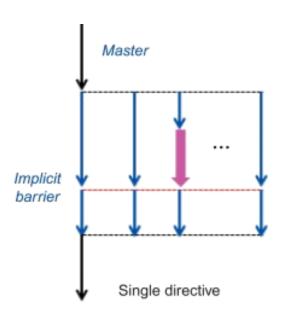
How the runtime decides to adjust the number of threads, it's implementation specific, might change from vendor to vendor.

- There are equivalent environment variables doing the same thing for nested parallelism
  - Example: if you use in bash export OMP\_NUM\_THREADS="4,2", you will get outer level parallelism to run with 4 threads, while the next level of nested parallelism to run with 2 threads. Then, you have 8 threads active: 4 X 2.

## The OpenMP single directive

- single directive defines a section of code to run by only a single thread in a parallel region
  - Similar to master that defines a section of code to run by only the master thread, single allows any thread to run the section of code whichever thread reaches the region first
- Synchronization behavior is a bit different ...
  - single directive has an implicit barrier upon completion of the region, where all threads wait for synchronization
  - master directive doesn't have the barrier. The other threads move and do their business





#### Example: Running a Function Once in a Nested Parallel Region

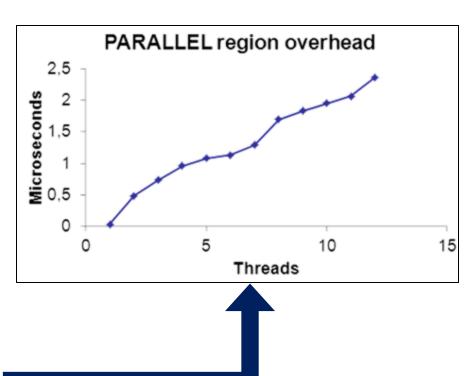
```
#include <omp.h>
// function to print level by just one thread
void report num threads(int whichLevel) {
    #pragma omp single
    printf("Level %d: number of threads in the team - %d\n", whichLevel, omp_get_num_threads());
int main() {
    omp set dynamic(1);
    omp set_nested(1);
    #pragma omp parallel num threads(2)
        report num threads(1);
        #pragma omp parallel num threads(3)
                                                                       Level 1: number of threads in the team - 2
            report num threads(2);
                                                                       Level 2: number of threads in the team - 3
            #pragma omp parallel num_threads(4)
                                                                       Level 2: number of threads in the team -
                                                                       Level 3: number of threads in the team -
                 report num threads(3);
                                                                       Level 3: number of threads in the team -
                                                                       Level 3: number of threads in the team -
                                                                       Level 3: number of threads in the team -
                                                                       Level 3: number of threads in the team -
    return(0);
                                                                       Level 3: number of threads in the team
```

#### Departing Thoughts about Nested Parallelism

- Examples of good reasons to employ nested parallelism
  - Insufficient parallelism at outer level
    - Example: an outer loop of 10 trips on a 64-core machine
  - Load balance scenario:
    - Many-core processors today can handle many threads
      - IBM POWER9 chip: handles 96 threads
    - More threads in flight can help balance the load



- Overhead of nested parallelism often outweighs its advantages
  - Scheduling, synchronization, hierarchical concurrency, etc. in a nested region comes with non-negligible cost!
- Better to redesign your algorithm in a flat one-dimensional loop with just one-level omp parallel for



#### Quiz: Remove Nested Parallelism

- Consider the following parallel loops with two-level nested parallelism:
  - run\_some\_independent\_function is totally independent of i and j

```
#pragma omp parallel for
for (int i = 0; i < N; i++) {
    #pragma omp parallel for
    for (int j = 0; j < M; j++) {
        some_independent_func(i, j);
    }
}</pre>
```

```
#pragma omp parallel for
for (int k = 0; k < N*M; k++) {
   some_independent_func(k/M, k%M);
}</pre>
```



Parallelize the code using just one level of omp parallel for

### Summary: What's Reasonable to Expect from OpenMP

- If you have more than N cores in an SMP setup, it's unlikely that you get a N-times speed-up
- Some of the reasons for lack of scaling:
  - Amdahl Law states that the maximum speed-up is limited to the portion that can be parallelized
  - Cache coherence (multiple caches can have their private copies for the same data)
  - False cache sharing (data too close to each other will cause cache access to be serialized)
- Execution scales somewhat better if you have an embarrassingly-parallel application
  - No dependencies exist at all! Just throw as many threads as possible supported by your machine!
  - Unfortunately, real-world parallel applications all have different types of dependencies