# **OpenMP loops**

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# Outline

- > Expressing parallelism
  - Understanding parallel threads
- > Memory Data management
  - Data clauses
- > Synchronization
  - Barriers, locks, critical sections
- > Work partitioning
  - Loops, sections, single work, tasks...

### > Execution devices

- Target



# What we saw so far..

### > Threads

- How to create and properly manage a team of threads
- How to join them with barriers

### > Memory

- How to create private and shared variables storages
- How to properly ensure memory consistency among parallel threads

### > Data syncronization

- How to create locks to implement, e.g., mutual exclusion
- How to identify Critical Sections
- How to ensure atomicity on single statements



- But..how can we split an existing workload among parallel threads?
   Say, a loop
- > Typical \$c€nario
  - 1. Analyze sequential code from customer/boss
  - 2. Parallelize it with OpenMP (for a "generic" parallel machine)
  - 3. Tune num\_threads for specific machine
  - 4. Get money/congratulations from customer/boss
  - > Might not be as easy as with PI Montecarlo!

How to do 2. without rewriting/re-engineering the code?



Let's code!

- > Create an array of N elements
  - Put inside each array element its index, multiplied by '2'
  - arr[0] = 0; arr[1] = 2; arr[2] = 4; ...and so on
- > Now, do it in parallel with a team of T threads
  - N = 19, T ≠ 19, N > T
  - Hint: Act on the boundaries of the loop
  - Hint #2: omp\_get\_thread\_num(), omp\_get\_num\_threads()
- > Example:









10

11

14

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# Loop partitioning among threads

- > Case #1: N multiple of T
  - Say, N = 20, T = 4
- > chunk = #iterations for each thread
- > Very simple..

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$$chunk = rac{N}{T}$$
 ;

 $i_{start} = thread_{ID} * chunk;$   $i_{end} = i_{start} + chunk + 1$ 

Let's

code!



- > Case #2: N not multiple of T
  - Say, N = 19, T = 4
- > chunk = #iterations for each thread (but last)
  - Last thread has less! (chunk<sub>last</sub>)



 $chunk = \frac{N}{T} + 1;$   $chunk_{last} = N \% chunk$ 

 $i_{start} = thread_{ID} * chunk;$   $i_{end} = \begin{cases} i_{start} + chunk & if not last thread \\ i_{start} + chunk_{last} & if last thread \end{cases}$ 

Let's

code!



- > Unfortunately, we don't know which thread will be "last" in time
- > But...we don't actually care the order in which iterations are executed
  - If there are not depenencies..
  - And..we do know that

0 <= omp\_get\_thread\_num() < omp\_get\_num\_threads()</pre>

> We choose that last thread as highest number



> Case #1 (N multiple of T)

$$chunk = \frac{N}{T}$$
  $i_{start} = thread_{ID} * chunk;$   $i_{end} = i_{start} + chunk$ 

Case #2 (N not multiple of T)

$$chunk = \frac{N}{T} + 1;$$
  $chunk_{last} = N \% chunk$ 

 $i_{start} = thread_{ID} * chunk;$   $i_{end} = \begin{cases} i_{start} + chunk & if not last thread \\ i_{start} + chunk_{last} & if last thread \end{cases}$ 



# Work sharing constructs

### > A way to distribute work among parallel threads

- In a simple, and "elegant" manner
- Using pragmas
- > OpenMP was born for this
  - OpeMP 2.5 targets regular, loop-based parallelism

### > OpenMP 3.x targets irregular/dynamic parallelism

- We will see it later



# The for construct

```
#pragma omp for [clause [[,] clause]...] new-line
  for-loops
Where clauses can be:
private(list)
firstprivate(list)
lastprivate(list)
linear(list[ : linear-step])
reduction (reduction-identifier : list)
schedule([modifier [, modifier]:]kind[, chunk size])
collapse(n)
ordered[(n)]
nowait.
```

- > The iterations will be executed in parallel by threads in the team
- The iterations are distributed across threads executing the parallel region to which the loop region binds
- > for-loops must have <u>Canonical loop form</u>



# **Canonical loop form**

for (init-expr; test-expr; incr-expr)
 structured-block

- > init-expr; test-expr; incr-expr not void
- > Eases programmers' life
  - More structured
  - Recommended also for "sequential programmers"
- > Preferrable to while and do...while
  - If possible





### > Create an array of *N* elements

- Put inside each array element its index, multiplied by '2'
- arr[0] = 0; arr[1] = 2; arr[2] = 4; ...and so on..

### > Now, do it in parallel with a team of *T* threads

- Using the for construct



# **Data sharing clauses**

#pragma omp for [clause [[,] clause]...] new-line
 for-loops

Where clauses can be:

```
private(list)
firstprivate(list)
lastprivate(list)
linear(list[ : linear-step])
reduction(reduction-identifier : list)
schedule([modifier [, modifier]:]kind[, chunk_size])
collapse(n)
ordered[(n)]
nowait
```

- > first/private, reduction we already know...
  - Private storage, w/ or w/o initialization
- > linear, we won't see



- > A list item that appears in a lastprivate clause is subject
  to the private clause semantics
- Also, the value is updated with the one from the sequentially last iteration of the associated loops



# lastprivate variables and memory

> Create a new storage for the variables, local to threads, and initialize







- > Modify the "PI Montecarlo" exercise
  - Use the for construct
- > Up to now, each threads executes its "own" loop

- i from 0 to 2499
- > Using the for construct, they actually share the loop
  - No need to modify the boundary!!!
  - Check it with printf



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2	K	











2500	
2501	
4999	

N N

7500
7501
9999





### > Create an array of N elements

- Put inside each array element its index, multiplied by '2'
- arr[0] = 0; arr[1] = 2; arr[2] = 4; ...and so on..

### > Declare the array as lastprivate

- So you can print its value after the parreg, in the sequential zone
- Do this at home



# OpenMP 2.5

### > OpenMP provides three work-sharing constructs

- Loops
- Single
- Sections



# The single construct

#pragma omp single [clause [[,] clause]...] new-line
 structured-block

Where clauses can be:

private(list)
firstprivate(list)
copyprivate(list)
nowait

- > The enclosed block is executed by only one threads in the team
- > ...and what about the other threads?



- > Each worksharing construct has an implicit barrier at its end
  - Example: a loop
  - If one thread is delayed, it prevents other threads to do useful work!!
  - Remember: barrier = consistent view of the sh memory





# Nowait clause in the for construct

#pragma omp for [clause [[,] clause]...] new-line
 for-loops

Where clauses can be:

```
private(list)
firstprivate(list)
lastprivate(list)
linear(list[ : linear-step])
reduction(reduction-identifier : list)
schedule([modifier [, modifier]:]kind[, chunk_size])
collapse(n)
ordered[(n)]
nowait
```

- > The nowait clause removes the barrier at the end of a worksharing (WS) construct
  - Applies to all of WS constructs
  - Does not apply to parregs!



> Removed the barrier at the end of WS construct

- Still, there is a barrier at the end of parreg

```
#pragma omp parallel num threads(4)
  #pragma omp for nowait
  for(int i=0; i<N; i++)</pre>
    . . .
  } // no barrier
  // USEFUL WORK!!
} // (implicit) barrier
```



# The sections construct

```
#pragma omp sections [clause[ [,] clause] ... ] new-line
  [#pragma omp section new-line]
    structured-block
  [#pragma omp section new-line]
    structured-block
Where clauses can be:
private(list)
firstprivate(list)
lastprivate(list)
reduction (reduction-identifier : list)
nowait
```

> Each section contains code that is executed by a single thread

- A "switch" for threads
- > Clauses, we already know..
  - lastprivate items are updated by the section executing last (in time)



### > Loops implement <u>data-parallel</u> paradigm

- Same work, on different data
- Aka: <u>data decomposition</u>, <u>SIMD</u>, <u>SPMD</u>
- > Sections implement <u>task-based paradigm</u>
  - Different work, on the same or different data
  - Aka: task decomposition, MPSD, MPMD



## The master construct

#pragma omp master new-line
 structured-block

#### No clauses

### > The structured block is executed only by master thread

- "Similar" to the single construct



### > It is **not** a work-sharing construct

- There is no barrier implied!!



# **Combined parreg+ws**

### > For each WS construct, there is also a compact form

- In this case, clauses to both constructs apply







- > Download the Code/ folder from the course website
- > Compile
- > \$ gcc -fopenmp code.c -o code
- > Run (Unix/Linux)
- \$ ./code
- > Run (Win/Cygwin)
- \$ ./code.exe



# References



- > Course website
  - <u>http://algo.ing.unimo.it/people/andrea/Didattica/HPC/</u>
- > My contacts
  - paolo.burgio@unimore.it
  - <u>http://hipert.mat.unimore.it/people/paolob/</u>
- > Useful links
  - <u>http://www.openmp.org</u>
  - <u>http://www.google.com</u>

