Introduction to Parallel Computing (CMSC416)



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Announcements

Assignment I is due on: March 7, 11:59 pm ET

- Good-faith attempt of each assignment is required
- Questions?
- Quiz I was due today, Feb. 23, at I I AM
 - You should be able to see your score, and correct answers, in ELMS
 - Questions?

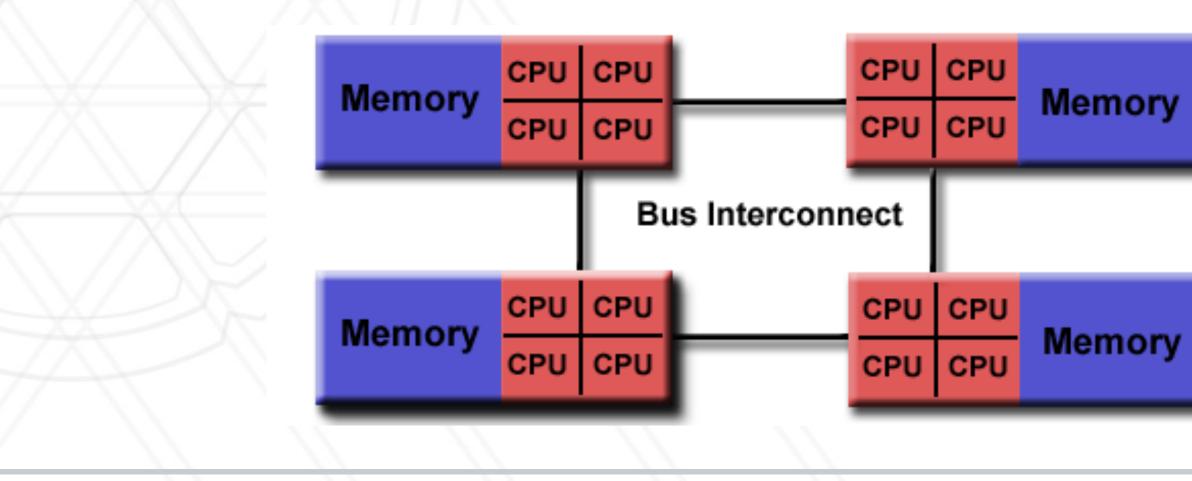




Shared memory programming

- All entities (threads) have access to the entire address space
- Threads "communicate" or exchange data by sharing variables
- User has to manage data conflicts







OpenMP

- OpenMP is an example of a shared memory programming model
- Provides on-node parallelization
- Meant for certain kinds of programs/computational kernels
 - Ones that use arrays and loops



Potentially easy to implement an application in parallel with small code changes



OpenMP

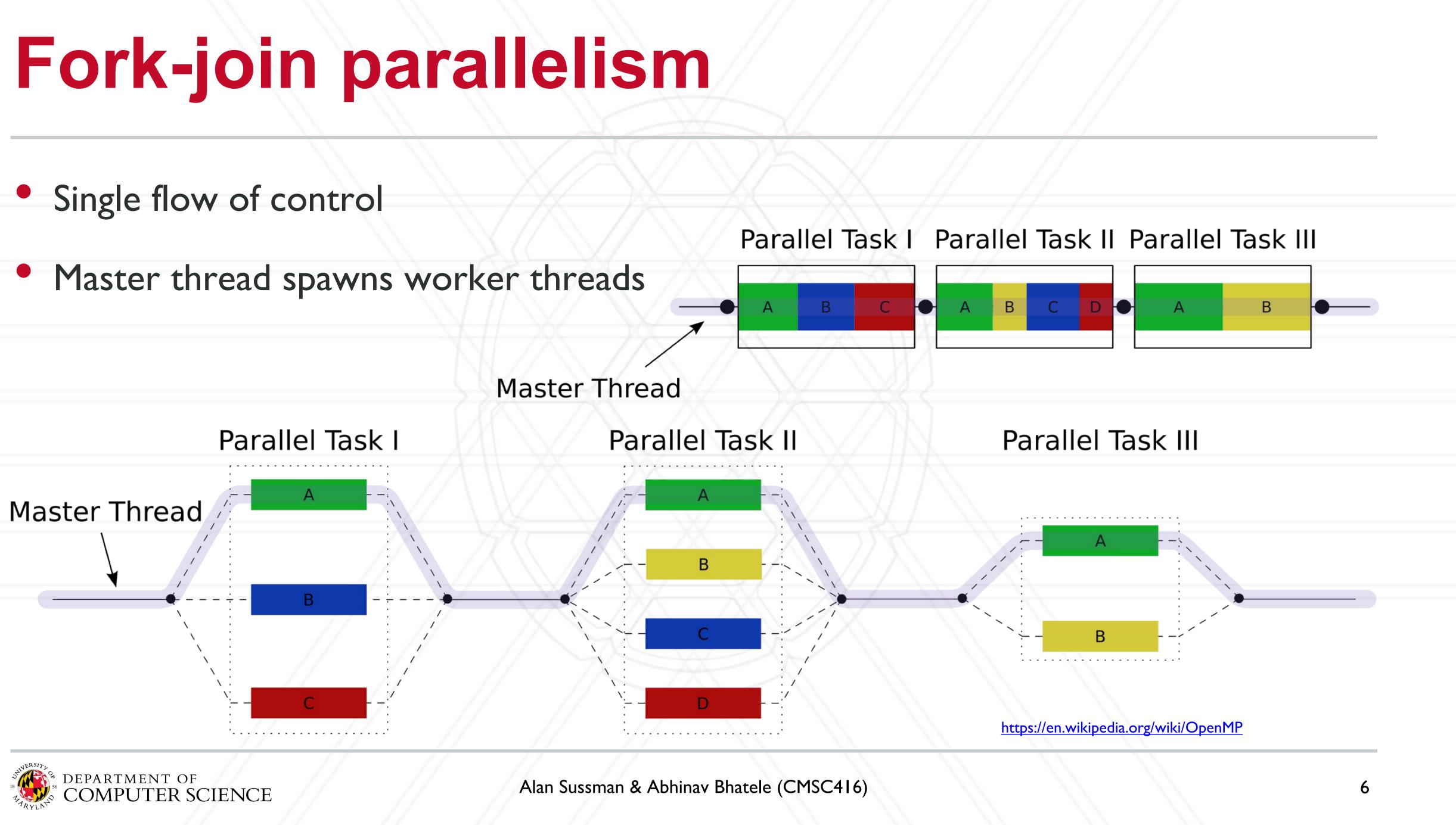
- code
- in the code and how to parallelize them
- Compiler converts code to multi-threaded code
- Fork/join model of parallelism



OpenMP is a language extension and library that enables parallelizing C/C++/Fortran

Programmer uses compiler directives and library routines to indicate parallel regions









Race conditions when threads interact

- Unintended sharing of variables can lead to race conditions
- Race condition: program outcome depends on the scheduling order of threads
 - Defined as one or more threads accessing a memory location with at least one of them performing a write, and without proper synchronization
- How can we prevent data races?
 - Use synchronization
 - Change how data is stored





OpenMP pragmas

- Pragma: a compiler directive in C or C++
- Mechanism to communicate with the compiler
- Compiler may ignore pragmas



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#pragma omp construct [clause [clause] ...]



Hello World in OpenMP

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

```
int main (void)
    #pragma omp parallel
    printf("Hello, world.\n");
    return 0;
```

- Compiling: gcc -fopenmp hello.c -o hello
- Setting number of threads: export OMP_NUM_THREADS=2





Parallel for

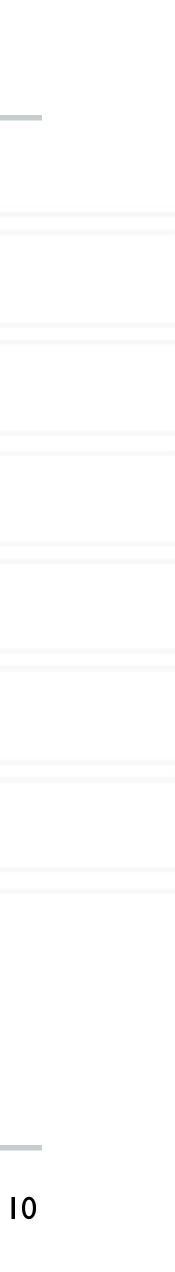
parallel

```
#pragma omp parallel for [clause [clause] ... ]
     • • •
    do work
    • • •
```



Directs the compiler that the immediately following for loop should be executed in

for (i = init; test expression; increment expression) {



Parallel for example

int main(int argc, char **argv) int a[100000]; #pragma omp parallel for

for (int i = 0; i < 100000; i++) { a[i] = 2 * i;

return 0;

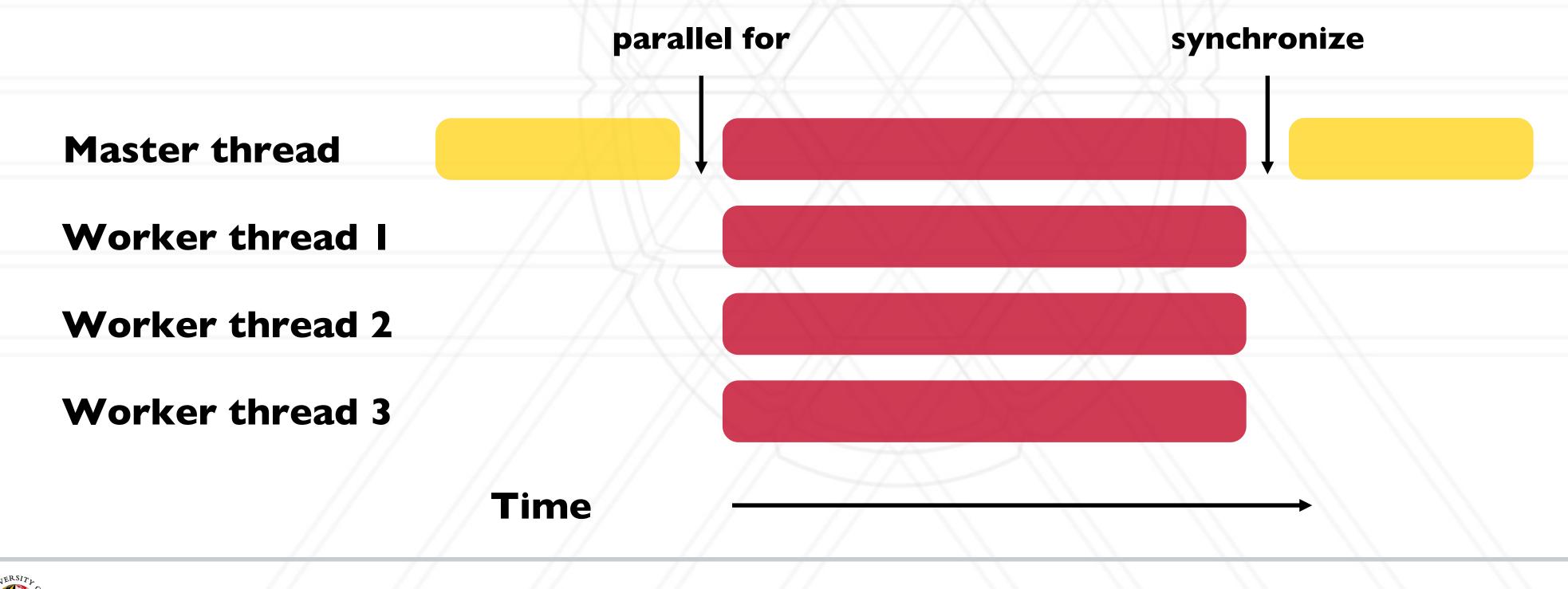




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Parallel for execution

- Master thread creates worker threads
- All threads divide iterations of the loop among themselves









Number of threads

Use environment variable

export OMP NUM THREADS=X

• Use void omp_set_num_threads(int num threads)

Set the number of OpenMP threads to be used in parallel regions

• int omp_get_num_procs(void);

Returns the number of available processors/cores

Can be used to decide the number of threads to create





Data sharing defaults

- Most variables are shared by default
- Global variables are shared
- Exception: loop index variables are private by default
- (thread-private)





• Stack variables in function calls from parallel regions are also private to each thread



saxpy (single precision a*x+y) example

#pragma omp parallel for for (int i = 0; i < n; i++) {</pre> z[i] = a * x[i] + y[i];





Overriding defaults using clauses

- Specify how data is shared between threads executing a parallel region
- private(list)
- shared(list)
- default(shared | none)
- reduction (operator: list)
- firstprivate(list)
- lastprivate(list)

https://www.openmp.org/spec-html/5.0/openmpsu106.html#x139-5540002.19.4





private clause

- Each thread has its own copy of the variables in the list
- Private variables are uninitialized when a thread starts
- region has been executed



• The value of a private variable is unavailable to the master thread after the parallel



default clause

Determines the data sharing attributes for variables for which this would be implicitly determined otherwise





Anything wrong with this example?

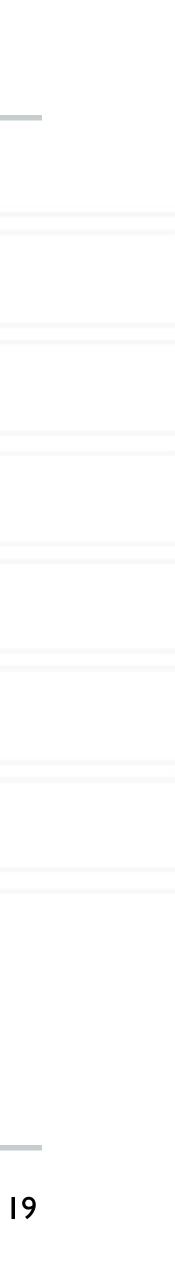
val = 5;

#pragma omp parallel for private(val) for (int i = 0; i < n; i++) { ... = val + 1;





The value of val will not be available to threads inside the loop



Anything wrong with this example?

#pragma omp parallel for private(val) for (int i = 0; i < n; i++) { val = i + 1;

printf("%d\n", val);



The value of val will not be available to the master thread outside the loop



firstprivate clause

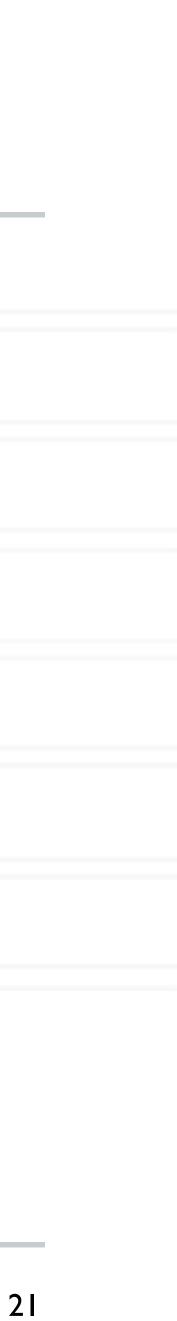
entry to the parallel section

val = 5;

#pragma omp parallel for firstprivate(val) for (int i = 0; i < n; i++) {</pre> ... = val + 1;



Initializes each thread's private copy to the value of the master thread's copy, on



lastprivate clause

- to the master's copy
- Last iteration determined by sequential order

#pragma omp parallel for lastprivate(val) for (int i = 0; i < n; i++) { val = i + 1;

printf("%d\n", val);



Writes the value belonging to the thread that executed the last iteration of the loop





reduction(operator: list) clause

Reduce values across private copies of

Operators: +, -, *, &, |, ^, &&, ||, max, min

#pragma omp parallel for reduction(+: val) for (int i = 0; i < n; i++) { val += i;

printf("%d\n", val);

https://www.openmp.org/spec-html/5.0/openmpsu107.html#x140-5800002.19.5



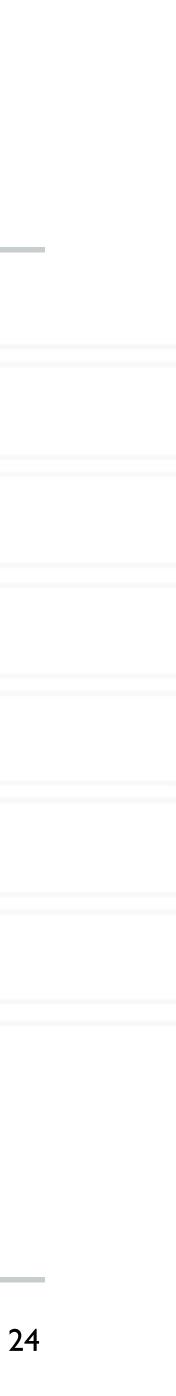
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Loop scheduling

- Assignment of loop iterations to different worker threads
- Default schedule tries to balance iterations among threads
- User-specified schedules are also available





User-specified loop scheduling

Schedule clause

- type: static, dynamic, guided, runtime
- static: iterations divided as evenly as possible (#iterations/#threads)
 - chunk < #iterations/#threads can be used to interleave threads</p>
- dynamic: assign a chunk size block to each thread
 - When a thread is finished, it retrieves the next block from an internal work queue
 - Default chunk size = I



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schedule (type[, chunk])



Other schedules

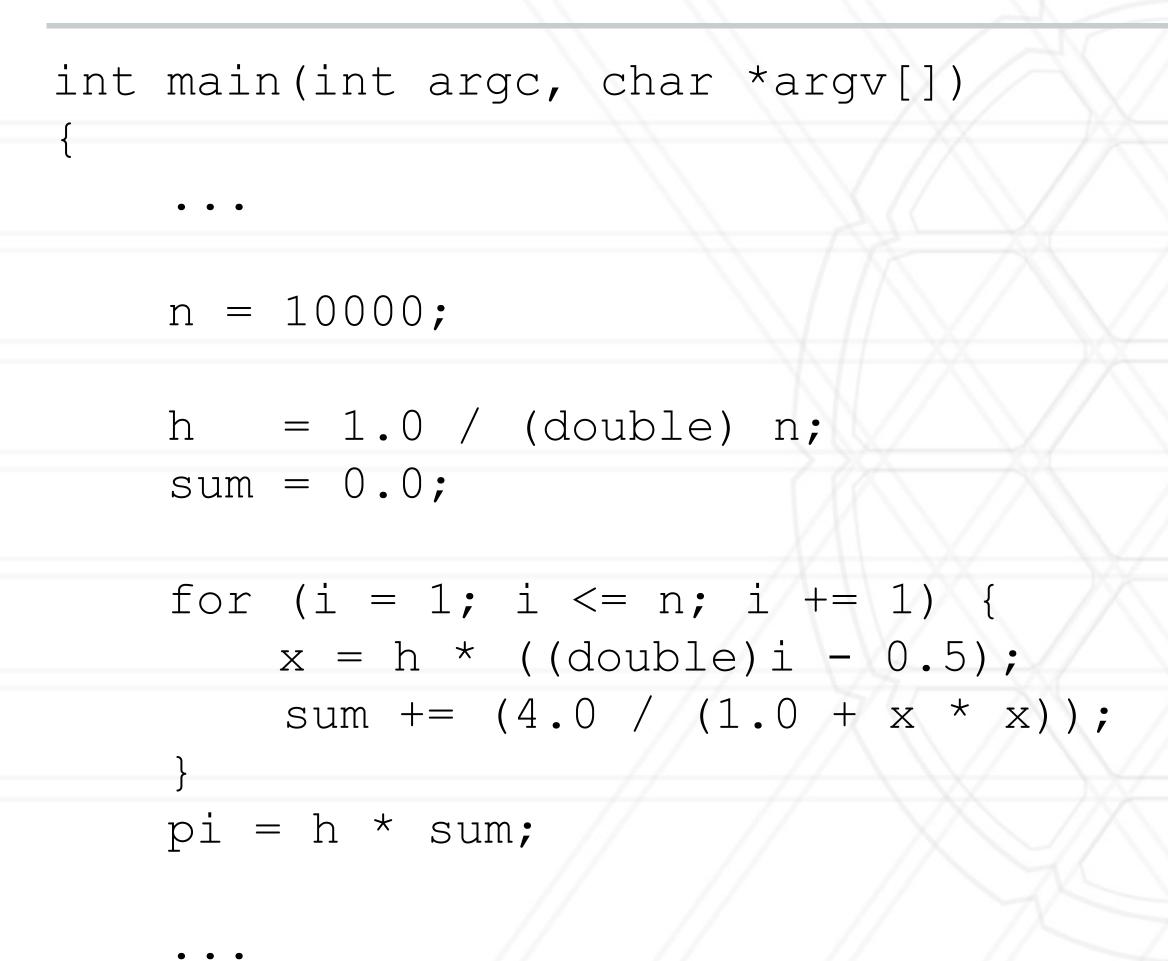
- guided: similar to dynamic but start with a large chunk size and gradually decrease it for handling load imbalance between iterations
- auto: scheduling delegated to the compiler
- runtime: use the OMP_SCHEDULE environment variable

https://software.intel.com/content/www/us/en/develop/articles/openmp-loop-scheduling.html

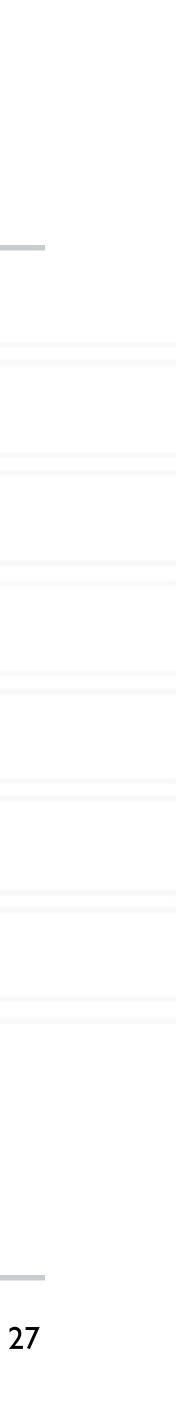




Calculate the value of $\pi = \int_0^1 \frac{4}{1+x^2}$







Calculate the value of $\pi = \int_0^1 \frac{4}{1+x^2}$

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

n = 10000;h = 1.0 / (double) n;sum = 0.0;

for (i = 1; i <= n; i += 1) { x = h * ((double)i - 0.5);sum += (4.0 / (1.0 + x * x));pi = h * sum;



• • •

• • •

#pragma omp parallel for firstprivate(h) private(x) reduction(+: sum)



Parallel region

• All threads execute the structured block

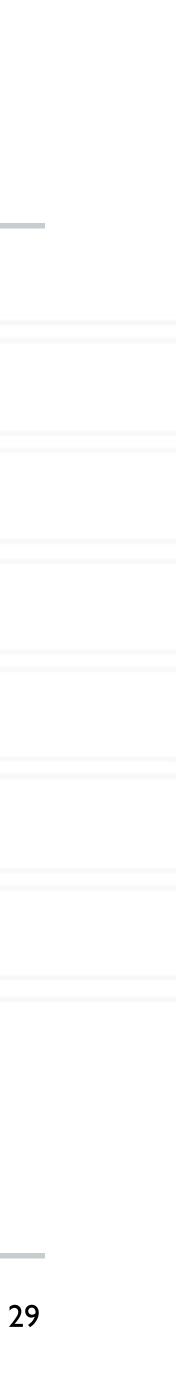
structured block

• Number of threads can be specified just like the parallel for directive



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#pragma omp parallel [clause [clause] ...]



Synchronization

- Concurrent access to shared data may result in inconsistencies
- Use mutual exclusion to avoid that
- critical directive
- atomic directive
- Library lock routines

https://software.intel.com/content/www/us/en/develop/documentation/advisor-user-guide/top/appendix/adding-parallelism-to-your-program/replacing-annotations-with-openmp-code/adding-openmp-code-tosynchronize-the-shared-resources.html



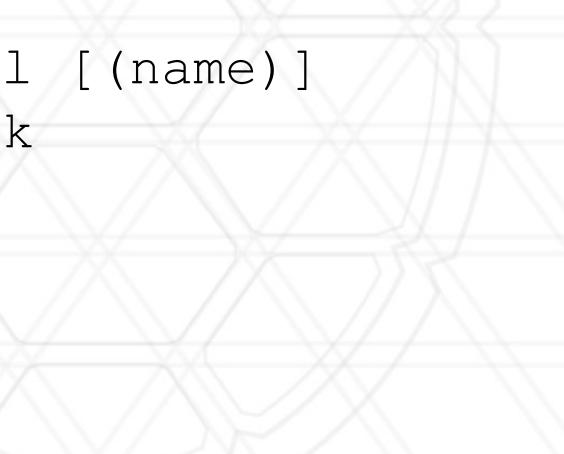


critical directive

Specifies that the code is only to be executed by one thread at a time

#pragma omp critical [(name)] structured block







atomic directive

Specifies that a memory location should be updated atomically

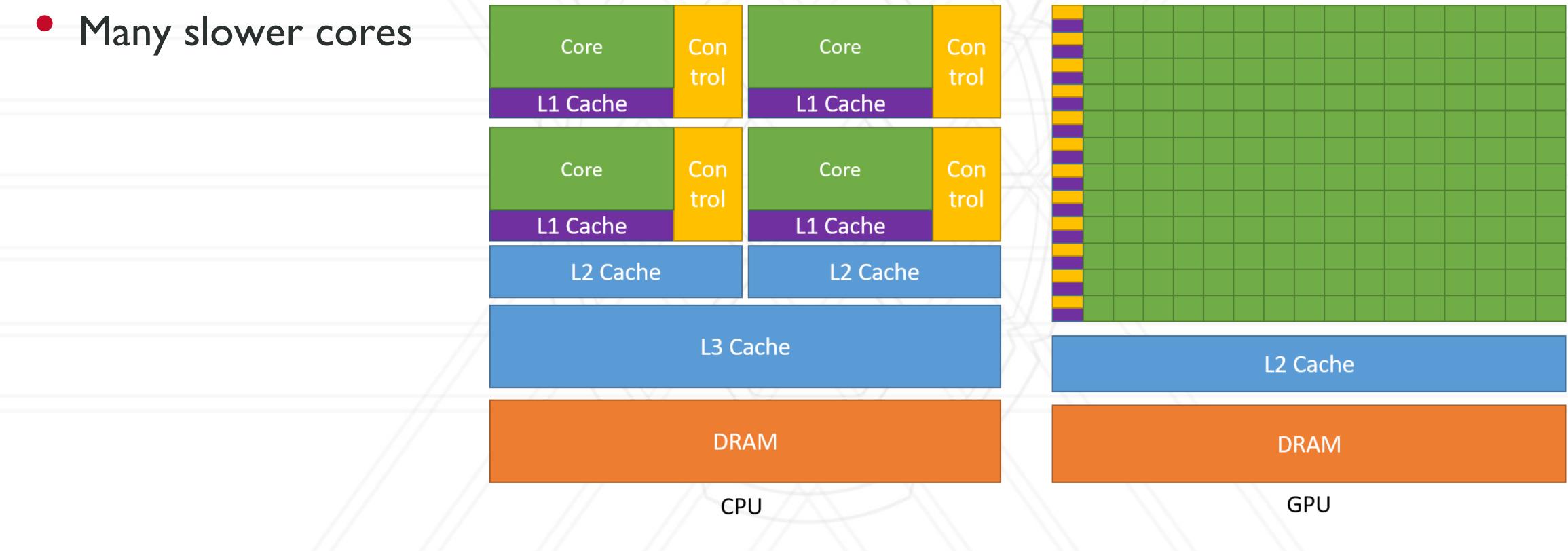
#pragma omp atomic expression





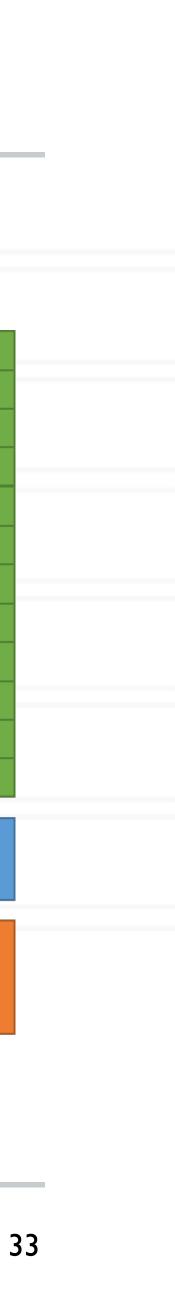
GPGPUs

GPGPU: General Purpose Graphical Processing Unit





https://docs.nvidia.com/cuda/cuda-c-programming-guide/index.html



OpenMP on GPUs

• *target*: run on accelerator / device

#pragma omp target teams distribute parallel for for (int i = 0; i < n; i++) { z[i] = a * x[i] + y[i];

• teams distribute: creates a team of worker threads and distributes work amongst them









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