

CPSC 4240/5240: Parallel Programming Techniques

Lecture 6: OpenMP

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Race Conditions in Parallel Instructions

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

int main() {
    #pragma omp parallel num_threads(4)
    {
        std::cout << "Hello from thread " <<
omp_get_thread_num() << std::endl;
    }
    return 0;
}
```

Each << is a **separate instruction**; thread can interleave arbitrarily in between

E.g., Hello from thread
Hello from thread 2 3
...

Why are Dependencies So Bad?

- Two threads writing to the same memory location causes **race conditions**
- Threads are “independent” within a parallel region and can **execute in any order**
- When threads are racing to modify the same memory address, we call this a **race condition**
- Depends on which threads execute first, can get **different results**
- Generally not what we want in programming (getting different results)!

Simple OpenMP Race Condition Example

What could happen here?

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

int main() {
    int sum = 0;

    #pragma omp parallel for
    for (int i = 0; i < 10000; i++) {
        // Race condition: no synchronization!
        sum += 1;
    }

    std::cout << "Final sum = " << sum << std::endl;
    return 0;
}
```

Simple OpenMP Race Condition Example

sum is a shared variable

**Two threads
can add to sum
simultaneously**

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

int main() {
    int sum = 0;

    #pragma omp parallel for
    for (int i = 0; i < 10000; i++) {
        // Race condition: no synchronization!
        sum += 1;
    }

    std::cout << "Final sum = " << sum << std::endl;
    return 0;
}
```

Simple OpenMP Race Condition Example

**Two threads
can add to sum
simultaneously**

**Won't give
10000 with
high likelihood**

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

int main() {
    int sum = 0;

    #pragma omp parallel for
    for (int i = 0; i < 10000; i++) {
        // Race condition: no synchronization!
        sum += 1;
    }

    std::cout << "Final sum = " << sum << std::endl;
    return 0;
}
```

Simple OpenMP Race Condition Example

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

```
int main() {
    int sum = 0;
```

```
    #pragma omp parallel for
    for (int i = 0; i < 10000; i++) {
        // Race condition: no synchronization!
        sum += 1;
    }
```

```
    std::cout << "Final sum = " << sum << std::endl;
    return 0;
}
```

Final sum = 9992
Final sum = 9975
Final sum = 9941
Final sum = 10000
Final sum = 9987
Final sum = 9823

How do we fix it?

- Use a **reduction clause**
 - `#pragma omp parallel for reduction(+: sum)`
 - `reduction([operation]: [variable])`
 - **Performs operation on variable safely**
 - Under the hood: uses a **parallel reduce primitive**
 - Parallel primitive algorithm like those we discussed last week
 - Calculates partial sums and aggregate them

How do we fix it?

- Use a **reduction clause**
 - `#pragma omp parallel for reduction(+: sum)`

```
#pragma omp parallel for reduction(+: sum)
for (int i = 0; i < 10000; i++) {
    sum += 1;
}
```

What is allowed in reduction?

Arithmetic operators: `+`, `-`, `*`

Bitwise operators: `&`, `|`, `^`

Logical operators: `&&`, `||`

`min`, `max`

How do we fix it?

- Use a **reduction clause**
 - `#pragma omp parallel for reduction(+: sum)`

```
#pragma omp parallel for reduction(+: sum)
for (int i = 0; i < 10000; i++) {
    sum += 1;
}
```

What variables are allowed in reduction?

scalar variables: int, float, double, bool

Custom reductions (OpenMP 4.0+)

Subranges (OpenMP 4.5+)

Another Way to Fix

- Use an **atomic variable**
 - #pragma omp atomic

```
#pragma omp parallel for
for (int i = 0; i < 10000; i++) {
    #pragma omp atomic
    sum += 1;
}
```

Another Way to Fix: Atomic Variables

- What are **atomic variables**?
- Atomic variables are variables that can be accessed (read, written modified) without interference from multiple threads
 - Prevents race conditions
 - Operations on atomic variables happen as a single, individual (atomic) operation
- Under the hood: generate a hardware atomic instruction (eg. LOCK XADD) or internal (software) lock if hardware lock isn't available
- Locks are expensive! Can lead to serialization due to contention

Another Way to Fix: Atomic Variables

- What are **atomic variables**?
- Atomic variables are variables that can be modified without interference from multiple processors
 - Prevents race conditions
 - Operations on atomic variables happen as a single, individual (atomic) operation
- Under the hood: generate a hardware atomic instruction (eg. LOCK XADD) or internal (software) lock if hardware lock isn't available
- Locks are expensive! Can lead to serialization due to contention

Use to (quickly) ensure safety when you have infrequent updates to the same variable!

Dealing with Loop Dependencies

- Last class we discussed issues with data race conditions
- Two threads writing to the same memory can cause incorrect outputs because they are competing to write
- OpenMP will likely **not warn** you about these race conditions
- Best way to resolve this is to restructure the algorithm/code to avoid race conditions
- Let's take a look at some more examples

OpenMP **nowait**

- By default, many work-sharing constructs contain a barrier at the end
 - for
 - sections
 - single
 - **Does not work for:** parallel

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

Will return an error

OpenMP **nowait**

- By default, many work-sharing constructs contain a barrier at the end
 - for
 - sections
 - single
 - **Does not work for:** parallel
- Recall that barriers act as joins, all threads complete tasks before going to the next parts of the code

OpenMP **nowait**

- OpenMP **nowait** construct modifies default synchronization behavior of OpenMP constructs
- Default OpenMP synchronization results in unnecessary waiting sometimes when not all threads need to finish before moving on to the next computation
- Let's take a look at an example

OpenMP **nowait**

Is this code correct?

```
#pragma omp parallel
{
    // First loop: Compute a[i] based on b[i+1] and b[i-1]
    #pragma omp for nowait
    for (int i = 1; i < N - 1; ++i) {
        a[i] = (b[i + 1] - b[i - 1]) / (2.0 * h);
    }

    // Second loop: Compute b[i] based on a[i+1] and a[i-1]
    #pragma omp for
    for (int i = N - 2; i > 0; --i) {
        b[i] = (a[i + 1] - a[i - 1]) / (2.0 * h);
    }
} // End of parallel region
```

OpenMP **nowait**

No! Why is it not correct?

```
#pragma omp parallel
{
    // First loop: Compute a[i] based on b[i+1] and b[i-1]
    #pragma omp for nowait
    for (int i = 1; i < N - 1; ++i) {
        a[i] = (b[i + 1] - b[i - 1]) / (2.0 * h);
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    // Second loop: Compute b[i] based on a[i+1] and a[i-1]
    #pragma omp for
    for (int i = N - 2; i > 0; --i) {
        b[i] = (a[i + 1] - a[i - 1]) / (2.0 * h);
    }
} // End of parallel region
```

OpenMP **nowait**

Results will be incorrect
because second loop
depends on first

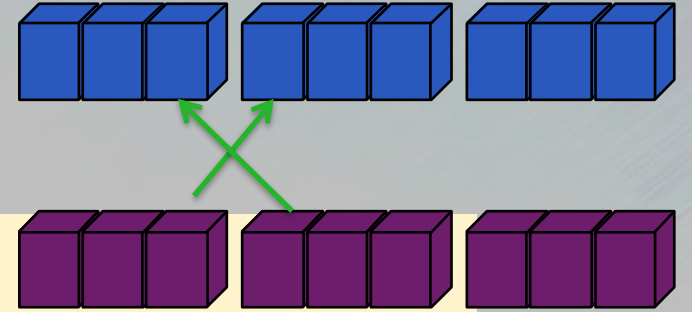
```
#pragma omp parallel
{
    // First loop: Compute a[i] based on b[i+1] and b[i-1]
    #pragma omp for nowait
    for (int i = 1; i < N - 1; ++i) {
        a[i] = (b[i + 1] - b[i - 1]) / (2.0 * h);
    }

    // Second loop: Compute b[i] based on a[i+1] and a[i-1]
    #pragma omp for
    for (int i = N - 2; i > 0; --i) {
        b[i] = (a[i + 1] - a[i - 1]) / (2.0 * h);
    }
} // End of parallel region
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OpenMP nowait

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    for (int i = N - 2; i > 0; --i) {
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OpenMP nowait

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        a[i] = (b[i + 1] - b[i - 1]) / (2.0 * h);
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    #pragma omp for
    for (int i = N - 2; i > 0; --i) {
        b[i] = (a[i + 1] - a[i - 1]) / (2.0 * h);
    }
} // End of parallel region
```

Suppose $a[i] = 0$ and $b[i] = i$ at initialization

OpenMP nowait

Suppose $a[i] = 0$ and $b[i] = i$ at initialization

```
#pragma omp parallel
```

```
{
```

```
// First loop: Compute a[i] based on b[i+1] and b[i-1]
```

```
#pragma omp for nowait
```

```
for (int i = 1; i < N - 1; ++i) {
```

```
     $a[i] = (b[i + 1] - b[i - 1]) / (2.0 * h);$ 
```

```
}
```

```
// Second loop: Compute b[i] based on a[i]
```

```
#pragma omp for
```

```
for (int i = N - 2; i > 0; --i) {
```

```
     $b[i] = (a[i + 1] - a[i - 1]) / (2.0 * h);$ 
```

```
}
```

```
} // End of parallel region
```

Sample output:

Last 10 elements 'a':

a[990] = 0.00000

a[991] = 0.00000

a[992] = 0.00000

a[993] = 0.00000

a[994] = 0.00000

a[995] = 0.00000

a[996] = 0.00000

a[997] = 0.00000

a[998] = 49950.00000

a[999] = 0.00000

Last 10 elements 'b':

b[990] = 0.00000

b[991] = 0.00000

b[992] = 0.00000

b[993] = 0.00000

b[994] = 0.00000

b[995] = 0.00000

b[996] = 0.00000

b[997] = 0.00000

b[998] = 0.00000

b[999] = 999.00000

OpenMP nowait

```
#pragma omp parallel
{
    // First loop: Compute a[i] based on b[i+1] and b[i-1]
    #pragma omp for nowait
    for (int i = 1; i < N - 1; ++i) {
        a[i] = (b[i + 1] - b[i - 1]) / (2.0 * h);
    }

    // Second loop: Compute b[i] based on a[i+1] and a[i-1]
    #pragma omp for
    for (int i = N - 2; i > 0; --i) {
        b[i] = (a[i + 1] - a[i - 1]) / (2.0 * h);
    }
} // End of parallel region
```

Is this correct?

Why did this happen?

Sample output:

Last 10 elements 'a':

a[990] = 0.00000
a[991] = 0.00000
a[992] = 0.00000
a[993] = 0.00000
a[994] = 0.00000
a[995] = 0.00000
a[996] = 0.00000
a[997] = 0.00000
a[998] = 49950.00000
a[999] = 0.00000

Last 10 elements 'b':

b[990] = 0.00000
b[991] = 0.00000
b[992] = 0.00000
b[993] = 0.00000
b[994] = 0.00000
b[995] = 0.00000
b[996] = 0.00000
b[997] = 0.00000
b[998] = 0.00000
b[999] = 999.00000

OpenMP nowait

```
#pragma omp parallel
{
    // First loop: Compute a[i] based on b
    #pragma omp for nowait
    for (int i = 1; i < N - 1; ++i) {
        a[i] = (b[i + 1] - b[i - 1]) / (2.0 * h);
    }

    // Second loop: Compute b[i] based on a
    #pragma omp for
    for (int i = N - 2; i > 0; --i) {
        b[i] = (a[i + 1] - a[i - 1]) / (2.0 * h);
    }
} // End of parallel region
```

The $b[i]$'s **computed their values with the default values of $a[i]$ before the $a[i]$'s computed their values!**

Sample output:

Last 10 elements 'a':

a[990] = 0.00000
a[991] = 0.00000
a[992] = 0.00000
a[993] = 0.00000
a[994] = 0.00000
a[995] = 0.00000
a[996] = 0.00000
a[997] = 0.00000
a[998] = 49950.00000
a[999] = 0.00000

Last 10 elements 'b':

b[990] = 0.00000
b[991] = 0.00000
b[992] = 0.00000
b[993] = 0.00000
b[994] = 0.00000
b[995] = 0.00000
b[996] = 0.00000
b[997] = 0.00000
b[998] = 0.00000
b[999] = 999.00000

OpenMP **nowait**

```
#pragma omp parallel
{
    // First loop: Compute a[i] based on b[i+1] and b[i-1]
    #pragma omp for
    for (int i = 1; i < N - 1; ++i) {
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    }
} // End of parallel region
```

Now **we remove nowait** so we allow for the default behavior with the for loop (with the barrier)

OpenMP **nowait**

```
#pragma omp parallel
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    // First loop: Compute a[i] based on b[i+1] and b[i-1]
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Now **we remove nowait** so we allow for the default behavior with the for loop (with the barrier)

Sample output:

Last 10 elements 'a':

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a[990] = 100.00000
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a[995] = 100.00000
a[996] = 100.00000
a[997] = 100.00000
a[998] = 100.00000
a[999] = 0.00000
```

Last 10 elements 'b':

```
b[990] = 0.00000
b[991] = 0.00000
b[992] = 0.00000
b[993] = 0.00000
b[994] = 0.00000
b[995] = 0.00000
b[996] = 0.00000
b[997] = 0.00000
b[998] = -5000.00000
b[999] = 999.00000
```

OpenMP nowait

```
#pragma omp parallel
{
    // First loop: Compute a[i] based on b[i+1] and b[i-1]
    #pragma omp for
    for (int i = 1; i < N - 1; ++i) {
        a[i] = (b[i + 1] - b[i - 1]) / (2.0 * h);
    }

    // Second loop: Compute b[i] based on a[i]
    #pragma omp for
    for (int i = N - 2; i > 0; --i) {
        b[i] = (a[i + 1] - a[i - 1]) / (2.0 * h);
    }
} // End of parallel region
```

This is now the correct behavior!

Questions?

Sample output:

Last 10 elements 'a':

```
a[990] = 100.00000
a[991] = 100.00000
a[992] = 100.00000
a[993] = 100.00000
a[994] = 100.00000
a[995] = 100.00000
a[996] = 100.00000
a[997] = 100.00000
a[998] = 100.00000
a[999] = 0.00000
```

Last 10 elements 'b':

```
b[990] = 0.00000
b[991] = 0.00000
b[992] = 0.00000
b[993] = 0.00000
b[994] = 0.00000
b[995] = 0.00000
b[996] = 0.00000
b[997] = 0.00000
b[998] = -5000.00000
b[999] = 999.00000
```

OpenMP **nowait**

Can also add an
explicit barrier

```
#pragma omp parallel
{
    // First loop: Compute a[i] based on b[i+1] and b[i-1]
    #pragma omp for nowait
    for (int i = 1; i < N - 1; ++i) {
        a[i] = (b[i + 1] - b[i - 1]) / (2.0 * h);
    }

    #pragma omp barrier

    // Second loop: Compute b[i] based on a[i+1] and a[i-1]
    #pragma omp for
    for (int i = N - 2; i > 0; --i) {
        b[i] = (a[i + 1] - a[i - 1]) / (2.0 * h);
    }
} // End of parallel region
```

OpenMP **nowait**

```
#pragma omp parallel
{
    // First loop: Compute a[i] based on b[i+1] and b[i-1]
    #pragma omp for nowait
    for (int i = 1; i < N - 1; ++i) {
        a[i] = (b[i + 1] - b[i - 1]) / (2.0 * h);
    }

    #pragma omp barrier

    // Second loop: Compute b[i] based on a[i]
    #pragma omp for
    for (int i = N - 2; i > 0; --i) {
        b[i] = (a[i + 1] - a[i - 1]) / (2.0 * h);
    }
} // End of parallel region
```

Can also add an
explicit barrier

Explicit or implicit
barrier: **no thread
may pass the
barrier until all
threads** have arrived
at the barrier

OpenMP **nowait**

```
#pragma omp parallel
{
    // First loop: Compute a[i] based on b[i+1]
    #pragma omp for nowait
    for (int i = 1; i < N - 1; ++i) {
        a[i] = (b[i + 1] - b[i - 1]) / (2.0 * h);
    }

    #pragma omp barrier

    // Second loop: Compute b[i] based on a[i]
    #pragma omp for
    for (int i = N - 2; i > 0; --i) {
        b[i] = (a[i + 1] - a[i - 1]) / (2.0 * h);
    }
} // End of parallel region
```

In this case, no writes will occur in b until all threads have written in a

Explicit or implicit barrier: **no thread may pass the barrier until all threads** have arrived at the barrier

More Dependency Examples

- Some loops cannot be parallelized at all!
- Need to be careful about which loops you can or cannot parallelize
- Let's look at some examples

More Dependency Examples

- Which of the following loops can OpenMP parallelize?

A)

```
for (int i = 1; i < N; i++)  
    A[i] = A[i] + B[i-1];
```

B)

```
for (int i = 0; i < N - 1; i++)  
    A[i + 1] = A[i] + 1;
```

C)

```
for (int i = 1; i < N; i++)  
    A[i] = A[0] + A[i];
```

D)

```
for (int i = 0; i < N-1; i++) {  
    A[i] = B[i];  
    C[i] = A[i] + B[i];  
    E[i] = C[i+1];  
}
```

More Dependency Examples

- Which of the following loops can OpenMP parallelize?

A)

```
for (int i = 1; i < N; i++)  
    A[i] = A[i] + B[i-1];
```

C)

```
for (int i = 1; i < N; i++)  
    A[i] = A[0] + A[i];
```

B)

```
for (int i = 0; i < N - 1; i++)  
    A[i + 1] = A[i] + 1;
```

D)

```
for (int i = 0; i < N-1; i++) {  
    A[i] = B[i];  
    C[i] = A[i] + B[i];  
    E[i] = C[i+1];  
}
```

More Dependency Examples

Let's go through these examples one by one!

- Which of the following loops can OpenMP parallelize?

A)

```
for (int i = 1; i < N; i++)  
    A[i] = A[i] + B[i-1];
```

C)

```
for (int i = 1; i < N; i++)  
    A[i] = A[0] + A[i];
```

B)

```
for (int i = 0; i < N - 1; i++)  
    A[i + 1] = A[i] + 1;
```

D)

```
for (int i = 0; i < N-1; i++) {  
    A[i] = B[i];  
    C[i] = A[i] + B[i];  
    E[i] = C[i+1];  
}
```

Let's go through these examples one by one!

More Dependency Examples

A)

```
for (int i = 1; i < N; i++)  
    A[i] = A[i] + B[i-1];
```

- Each $A[i]$ is calculated **independently of other $A[i]$'s**
- Each $A[i]$ depends only on itself and $B[i-1]$ neither of which are modified by other threads

Let's go through these examples one by one!

More Dependency Examples

- Here, each $A[i]$ depends on the previous value of $A[i]$
- What can go wrong here?
- **$A[i + 1]$ calculated with old value of $A[i]$** before $A[i]$ is calculated!

B)

```
for (int i = 0; i < N - 1; i++)  
    A[i + 1] = A[i] + 1;
```

Let's go through these examples one by one!

More Dependency Examples

C)

```
for (int i = 1; i < N; i++)  
    A[i] = A[0] + A[i];
```

- $A[i]$ only depends on itself and $A[0]$
- Iteration starts at 1 and doesn't recompute $A[0]$
- Hence, $A[0]$ is a static value and each $A[i]$ computation is independent of other $A[i]$'s

Let's go through these examples one by one!

More Dependency Examples

- There are several dependencies in this computation
- What can go wrong here?
Hint: **threads can interleave!**
 - Meaning **a thread can perform some computation in the middle of another thread's computation**

D)

```
for (int i = 0; i < N-1; i++) {  
    A[i] = B[i];  
    C[i] = A[i] + B[i];  
    E[i] = C[i+1];  
}
```

Let's go through these examples one by one!

More Dependency Examples

- There are several dependencies in this computation
- What can go wrong here?
Hint: **threads can interleave!**
 - Meaning **a thread can perform some computation in the middle of another thread's computation**

- One thread can compute $E[i]$ **after $C[i+1]$ computed by another thread**

D)

```
for (int i = 0; i < N-1; i++) {  
    A[i] = B[i];  
    C[i] = A[i] + B[i];  
    E[i] = C[i+1];  
}
```

How would you change the algorithm to parallelize loop B?

Parallelized B)

```
for (int i = 0; i < N - 1; i++)  
    A[i + 1] = A[0] + i;
```

B)

```
for (int i = 0; i < N - 1; i++)  
    A[i + 1] = A[i] + 1;
```

Another example for nowait

- To ensure correctness of the following code, where must we remove the nowait clause(s)?

Another example for nowait

```
#pragma omp parallel for shared(c) private(i) nowait  
    for (i=0; i<N; i++)  
        c[i] = (double) i
```

```
#pragma omp parallel for shared(c) private(i) nowait  
    for (i=1; i<N; i+=2)  
        c[i] = c[i] + c[i-1]
```

```
#pragma omp parallel for shared(c, d) private(i) nowait  
    for (i=1; i<N; i+=2)  
        d[i] = c[i] + c[i-1]
```

```
#pragma omp parallel for shared(c, e) private(i) nowait  
    for (i=2; i<N; i+=2)  
        e[i] = c[i] + c[i-1]
```

Another example for nowait

```
#pragma omp parallel for shared(c) private(i) nowait  
    for (i=0; i<N; i++)  
        c[i] = (double) i
```

```
#pragma omp parallel for shared(c) private(i) nowait  
    for (i=1; i<N; i+=2)  
        c[i] = c[i] + c[i-1]
```

```
#pragma omp parallel for shared(c, e) private(i) nowait  
    for (i=1; i<N; i+=2)  
        d[i] = e[i] + e[i-1]
```

```
#pragma omp parallel for shared(c, e) private(i) nowait  
    for (i=2; i<N; i+=2)  
        e[i] = c[i] + c[i-1]
```

Another example for nowait

```
#pragma omp parallel for shared(c) private(i) nowait  
    for (i=0; i<N; i++)  
        c[i] = (double) i
```

```
#pragma omp parallel for shared(c) private(i)  
    for (i=1; i<N; i+=2)  
        c[i] = c[i] + c[i-1]
```

```
#pragma omp parallel for shared(c, e) private(i)  
    for (i=1; i<N; i+=2)  
        d[i] = e[i] + e[i-1]
```

```
#pragma omp parallel for shared(c, e) private(i) nowait  
    for (i=2; i<N; i+=2)  
        e[i] = c[i] + c[i-1]
```

Now **need to remove the nowait on the third loop** because we cannot use the new value of e!

Removing Data Dependencies More Examples

- How can we split the following loop into two loops so that we can parallelize each loop and the result is correct?

B initially: 0 1 2 3 4 5 6 7

B on 1 thread: 7 7 7 7 11 12 13 14

```
#pragma omp parallel for shared (N,B)
for (i = 0; i < N; i++)
    B[i] += B[N-1-i];
```

B[0] += B[7],	B[1] += B[6],	B[2] += B[5]
B[3] += B[4],	B[4] += B[3],	B[5] += B[2]
B[6] += B[1],	B[7] += B[0]	

Removing Data Dependencies More Examples

- How can we split the following loop into two loops so that we can parallelize each loop and the result is correct?

B initially: 0 1 2 3 4 5 6 7
B on 10 threads: 7 7 9 7 11 7 13 7

```
#pragma omp parallel for shared (N,B)
for (i = 0; i < N; i++)
    B[i] += B[N-1-i];
```

B[0] += B[7],	B[1] += B[6],	B[2] += B[5]
B[3] += B[4],	B[4] += B[3],	B[5] += B[2]
B[6] += B[1],	B[7] += B[0]	

Removing Data Dependencies More Examples

- How can we split the following loop into two loops so that we can parallelize each loop and the result is correct?

B initially: 0 1 2 3 4 5 6 7
B on 10 threads: 7 7 9 7 11 7 13 7

```
#pragma omp parallel for shared (N,B)
for (i = 0; i < N; i++)
    B[i] += B[N-1-i];
```

What is happening here?

Removing Data Dependencies More Examples

- How can we split the following loop into two loops so that we can parallelize each loop and the result is correct?

B initially: 0 1 2 3 4 5 6 7
B on 10 threads: 7 7 9 7 11 7 13 7

```
#pragma omp parallel for shared (N,B)
for (i = 0; i < N; i++)
    B[i] += B[N-1-i];
```

B[7] updated before B[0]!
B[5] updated before B[2]!
B[2] updated after B[5]!

Removing Data Dependencies More Examples

- How can we split the following loop into two loops so that we can parallelize each loop and the result is correct?

B initially: 0 1 2 3 4 5 6 7
B on 10 threads: 7 7 9 7 11 7 13 7

```
#pragma omp parallel for shared (N,B)
for (i = 0; i < N; i++)
    B[i] += B[N-1-i];
```

How do we fix this?

Removing Data Dependencies More Examples

- How can we split the following loop into two loops so that we can parallelize each loop and the result is correct?

B initially: 0 1 2 3 4 5 6 7
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```
#pragma omp parallel for shared (N,B)
for (i = 0; i < N; i++)
    B[i] += B[N-1-i];
```

How do we fix this? Hint: **think about splitting the loop at some point**

Removing Data Dependencies More Examples

```
#pragma omp parallel for shared (N,B)  
for (i = 0; i < int(N/2 - 0.5); i++)  
    B[i] += B[N-1-i];
```

```
#pragma omp parallel for shared (N,B)  
for (i = int(N/2 + 0.5); i < N; i++)  
    B[i] += B[N-1-i];
```

How do we fix this? Hint: **think about splitting the loop at some point**

Removing Data Dependencies More Examples

```
#pragma omp parallel for shared (N,B)
for (i = 0; i < int(N/2 - 0.5); i++)
    B[i] += B[N-1-i];
```

```
#pragma omp parallel for shared (N,B)
for (i = int(N/2 + 0.5); i < N; i++)
    B[i] += B[N-1-i];
```

Why? Because only the second half depends on the first half

Removing Data Dependencies More Examples

Is this still correct now?

```
#pragma omp parallel shared(N, B)
{
    #pragma omp for nowait
    for(int i = 0; i <= int(N/2 - 0.5); ++i) {
        B[i] += B[N-1-i];
    }

    #pragma omp for
    for(int i = int(N/2 + 0.5); i < N; ++i) {
        B[i] += B[N-1-i];
    }
}
```

Removing Data Dependencies More

Example No! Revert back to the previous situation

```
#pragma omp parallel shared(N, B)
{
    #pragma omp for nowait
    for(int i = 0; i <= int(N/2 - 0.5); ++i) {
        B[i] += B[N-1-i];
    }

    #pragma omp for
    for(int i = int(N/2 + 0.5); i < N; ++i) {
        B[i] += B[N-1-i];
    }
}
```

OpenMP Loop Scheduling

- OpenMP has more fine-grained control over **loop scheduling**
- **Loop scheduling** determines how iterations of a parallel loop are divided among available threads
- Helps with balancing workloads, reducing latency (idle time), maximize CPU utilization, in some cases, helps with correctness
- `#pragma omp for schedule(...)`

OpenMP Scheduling: A Quick Aside

- **OpenMP Default Scheduler for loops**

- Allows compilers to define their default scheduling policy
- Can vary between different compilers and different versions of the same compiler
- Static scheduling is the default across most compilers: GCC, Clang, Intel's ICC
- Loop iterations divided into equal-sized chunks, each thread assigned a specific chunk before loop begins
- Other types of schedulers: **work-stealing** (discussed in later lectures!)

OpenMP Loop Scheduling

- Types of schedules: `#pragma omp for schedule(...)`
- **Static scheduling**
 - Divides loop iterations into chunks of a fixed size and assigns them to threads in round-robin fashion before loop starts
 - Hence it is static

```
#pragma omp for schedule(static)
```

```
#pragma omp for schedule(static, chunk_size)
```

OpenMP Loop Scheduling

- Types of schedules: `#pragma omp for schedule(...)`

- **Static scheduling**

- Divides loop iterations into chunks of a fixed size and assigns them to threads in round-robin fashion before loop starts
 - Hence it is static

Default chunk size when no
chunk_size passed in

```
#pragma omp for schedule(static)  
#pragma omp for schedule(static, chunk_size)
```

OpenMP Loop Scheduling

- Types of schedules: `#pragma omp for schedule(...)`

- **Static scheduling**

- Divides loop iterations into chunks of a fixed size and assigns them to threads in round-robin fashion before loop starts
 - Hence it is static

```
#pragma omp for schedule(static,  
#pragma omp for schedule(static,
```

Default chunk size when no
chunk_size passed in

Default chunk size is:

$$\frac{\text{Total Iters}}{\text{Num Threads}}$$

OpenMP Loop Scheduling

- **Static scheduling**

- **Pros:**

- Predictable work distribution: each thread receives roughly equal number of iterations
 - Low overhead for runtime decision-making
 - Better cache performance

- **Cons:**

- Potential load imbalance for irregular workloads

Best for regular workloads!

OpenMP Loop Scheduling

- **Static scheduling**

- **Pros:**

- Predictable work distribution: each thread receives roughly equal number of iterations
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Best for regular workloads! Typically you don't have to assign the `chunk_size` yourself

OpenMP Loop Scheduling

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- **Pros:**

- Predictable work distribution: each thread receives roughly equal number of iterations
 - Low overhead for runtime decision-making
 - Better cache performance

- **Cons:**

- Potential load imbalance for irregular workloads

Best for regular workloads! Typically you don't have to assign the `chunk_size` yourself

OpenMP Loop Scheduling

- **Dynamic Scheduling:**

- Divides loop iterations into chunks of a fixed size and assigns them to threads **on-the-fly** as the threads become available
- Hence it is dynamic

```
#pragma omp for schedule(dynamic)  
#pragma omp for schedule(dynamic, chunk_size)
```

OpenMP Loop Scheduling

- **Dynamic Scheduling:**

- Divides loop iterations into chunks of a fixed size and assigns them to threads **on-the-fly** as the threads become available
- Hence it is dynamic

```
#pragma omp for schedule(dy  
#pragma omp for schedule(dy
```

Default chunk size is determined by compiler for dynamic scheduling

Default for most compilers is 1

OpenMP Loop Scheduling

- **Dynamic Scheduling:**

- Divides loop iterations into chunks of a fixed size and assigns them to threads **on-the-fly** as the threads become available
- Hence
 - Default chunk size is determined by compiler for dynamic scheduling

#pragma omp for

#pragma omp for

Default for most compilers is 1

Why is it different? One is **extreme for uniform workloads**
the other **extreme for irregular workloads**

OpenMP Loop Scheduling

- **Dynamic Scheduling:**

- **Pros:**

- Better load balancing for irregular workloads
 - Reduced risk of load imbalance

- **Cons:**

- Increased overhead
 - Nondeterministic execution order
 - Potential memory overhead for task queues and scheduling structures

Lecture 5 Quiz

