### MULTICORE PROGRAMMING WITH OPENMP

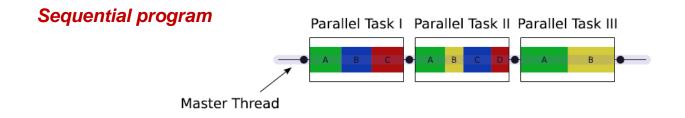
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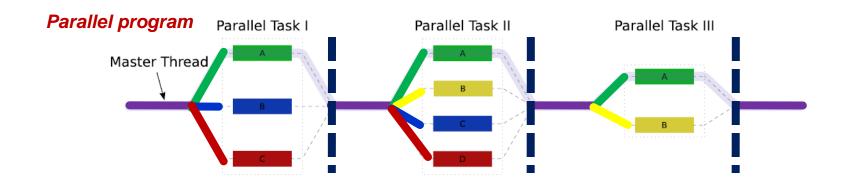
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#### Programming model: OpenMP

- De-facto standard for the shared memory programming model
- A collection of compiler directives, library routines and environment variables
- Easy to specify parallel execution within a serial code
- Requires special support in the compiler
- Generates calls to threading libraries (e.g. pthreads)
- Focus on loop-level parallel execution
- Popular in high-end embedded

### Fork/Join Parallelism





- Initially only master thread is active
- Master thread executes sequential code
- Fork: Master thread creates or awakens additional threads to execute parallel code
- Join: At the end of parallel code created threads are suspended upon barrier synchronization

#### Pragmas

- Pragma: a compiler directive in C or C++
- Stands for "pragmatic information"
- A way for the programmer to communicate with the compiler
- Compiler free to ignore pragmas: original sequential semantic is not altered
- Syntax:

#pragma omp <rest of pragma>

#### Components of OpenMP

#### **Directives**

- Parallel regions
  - #pragma omp parallel
- Work sharing
  - #pragma omp for
  - #pragma omp sections
- Synchronization
  - #pragma omp barrier
  - #pragma omp critical
  - #pragma omp atomic

#### **Runtime Library**

- Data scope attributes
  - private
  - shared
  - reduction
- Loop scheduling
  - static
  - dynamic
- Thread Forking/Joining

Clauses

- omp\_parallel\_start()
- omp\_parallel\_end()
- Loop scheduling
- Thread IDs
  - omp\_get\_thread\_num()
  - omp\_get\_num\_threads()

# Outlining parallelism The parallel directive

- Fundamental construct to outline parallel computation within a sequential program
- Code within its scope is replicated among threads
- Defers implementation of parallel execution to the runtime (machinespecific, e.g. pthread\_create)

### A sequential program.. .. is easily parallelized

```
int main()
{
    #pragma omp parallel
    {
       printf ("\nHello world!");
    }
}
```

```
int main()
{
   omp_parallel_start(&parfun, ...);
   parfun();
   omp_parallel_end();
}
int parfun(...)
{
   printf ("\nHello world!");
}
```

```
int main()
Code originally contained
within the scope of the
                                 omp_parallel_start(&parfun, ...);
pragma is outlined to a new
                                 parfun();
function within the compiler
                                 omp_parallel_end();
                               int parfun(...)
int main()
                                printf ("\nHello world!");
#pragma omp parallel
    printf ("\nHello world!");
```

```
int main()
The #pragma construct in
the main function is
                                 omp_parallel_start(&parfun, ...);
replaced with function calls
                                 parfun();
to the runtime library
                                 omp_parallel_end();
                               int parfun(...)
int main()
                                printf ("\nHello world!");
#pragma omp parallel
    printf ("\nHello world!");
```

First we call the runtime to fork new threads, and pass them a pointer to the function to execute in parallel

#pragma omp parallel

printf ("\nHello world!");

int main()

```
int main()
{
   omp_parallel_start(&parfun, ...);
   parrun();
   omp_parallel_end();
}
int parfun(...)
{
   printf ("\nHello world!");
}
```

```
int main()
Then the master itself calls
the parallel function
                                 omp_parallel_start(&parfun, ...);
                                 parfun();
                               int parfun(...)
int main()
                                printf ("\nHello world!");
#pragma omp parallel
    printf ("\nHello world!");
```

Finally we call the runtime to synchronize threads with a barrier and suspend them

#pragma omp parallel

int main()

```
int main()
                            omp_parallel_start(&parfun, ...);
                           parfun();
                            omp_parallel_end();
                          int parfun(...)
                          printf ("\nHello world!");
printf ("\nHello world!");
```

```
int main()
                                    Call runtime to get thread ID:
                                    Every thread sees a different value
  int id;
  int a = 5i
#pragma omp parallel
                                            Master and slave threads
                                             access the same variable a
    id = omp get thread num();
    if (id == 0)
      printf ("Master: a = %d.", a*2);
    else
      printf ("Slave: a = %d.", a
                    A slightly more complex example
```

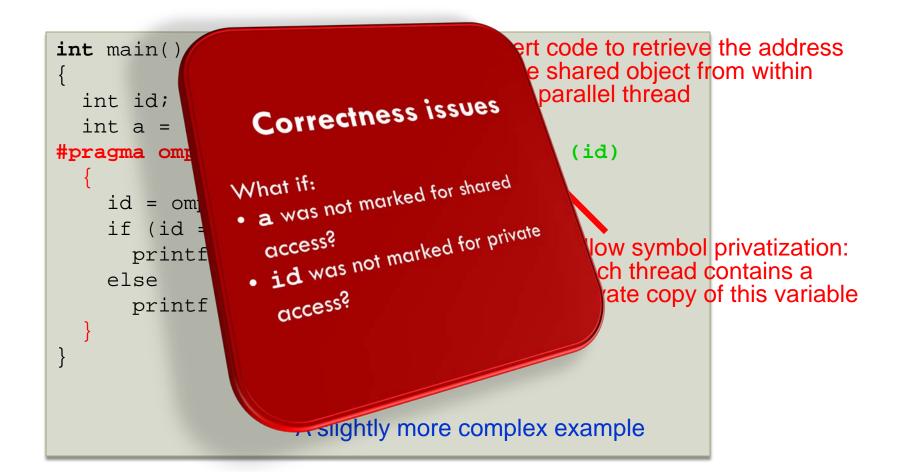
```
int main()
                                             Call runtime to get thread ID:
int id;
int a = 5;

#pragma omp parallel

#bread, num (ese different)

thread, num (ese different)
                                             Every thread sees a different value
     id = omp_get_thread_num()
behaviors?
                                                                 nd slave threads
      else
        printf ("Slave: a = %d.", a = %d)
                         A slightly more complex example
```

```
Insert code to retrieve the address
int main()
                                     of the shared object from within
                                     each parallel thread
  int id;
  int a = 5i
#pragma omp parallel shared (a) private (id)
    id = omp get thread num();
    if (id == 0)
                                            Allow symbol privatization:
      printf ("Master: a = %d.", a*2);
                                             Each thread contains a
    else
                                             private copy of this variable
     printf ("Slave: a = %d.", a);
                     A slightly more complex example
```



#### More data sharing clauses

- firstprivate
  - copyin, private storage
- lastprivate
  - Copyout, private storage

### Sharing work among threads The for directive

- The parallel pragma instructs every thread to execute all of the code inside the block
- If we encounter a for loop that we want to divide among threads, we use the for pragma

#pragma omp for

### #pragma omp for

The code of the **for** loop is moved inside the outlined function.

```
int main()
  omp_parallel_start(&parfun, ...);
 parfun();
  omp_parallel_end();
int parfun(...)
  int LB = \dots;
  int UB = ...;
  for (i=LB; i<UB; i++)
      a[i] = i;
```

### #pragma omp for

Every thread works on a different subset of the iteration space..

```
int main()
  omp_parallel_start(&parfun, ...);
  parfun();
  omp_parallel_end();
int parfun(...)
 int LB = ...;
```

### #pragma omp for

..since lower and upper boundaries (**LB**, **UB**) are computed locally

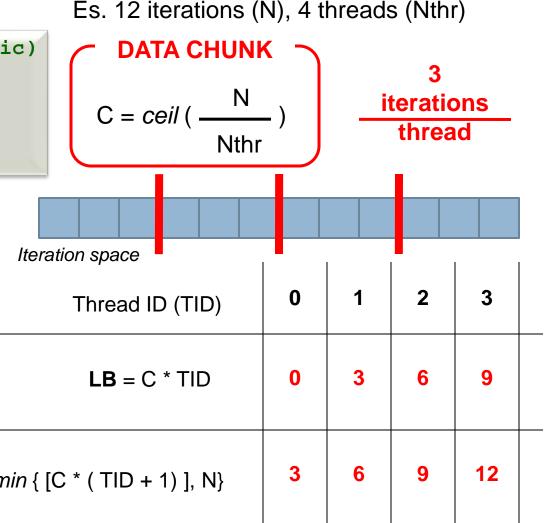
```
int main()
{
    #pragma omp parallel for
    {
       for (i=0; i<10; i++)
        a[i] = i;
    }
}</pre>
```

```
int main()
  omp_parallel_start(&parfun, ...);
  parfun();
  omp_parallel_end();
int parfun(...)
  int LB = \dots;
  int UB = ...;
  for (i=LB; i<UB; i++)
      a[i] = i;
```

#### #pragma omp for schedule(static) for (i=0; i<12; i++)a[i] = i;

#### Useful for:

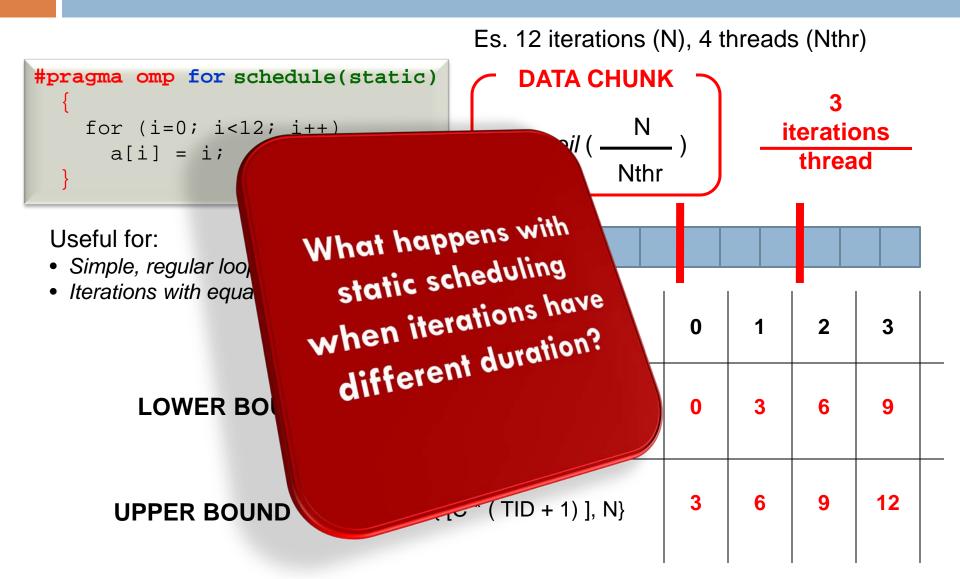
- Simple, regular loops
- Iterations with equal duration

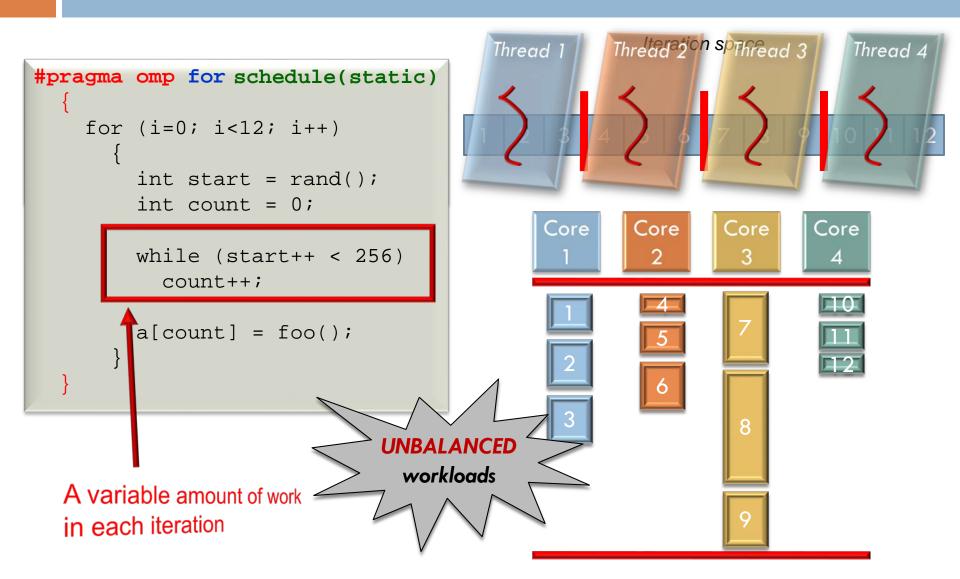


**UPPER BOUND** 

LOWER BOUND

$$UB = min \{ [C * (TID + 1)], N \}$$



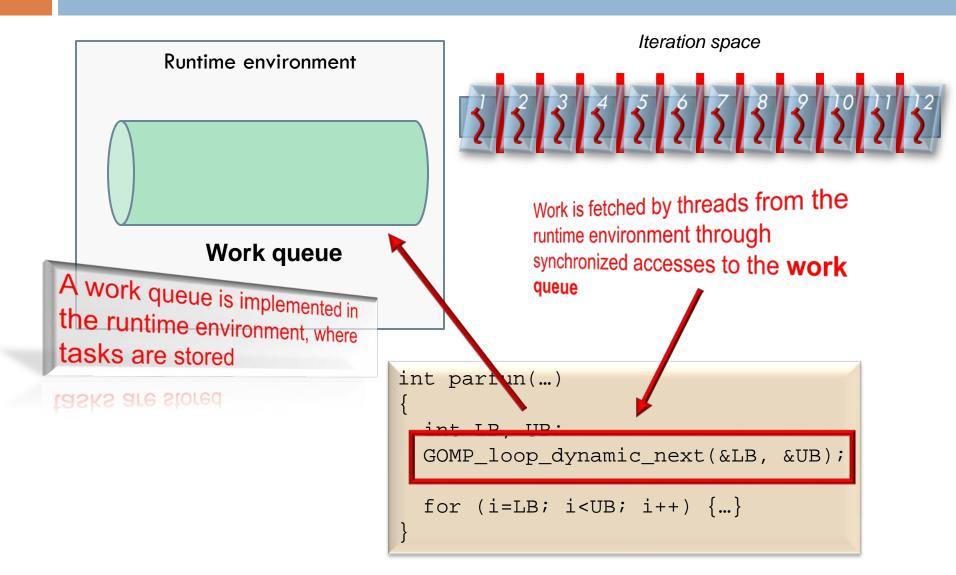


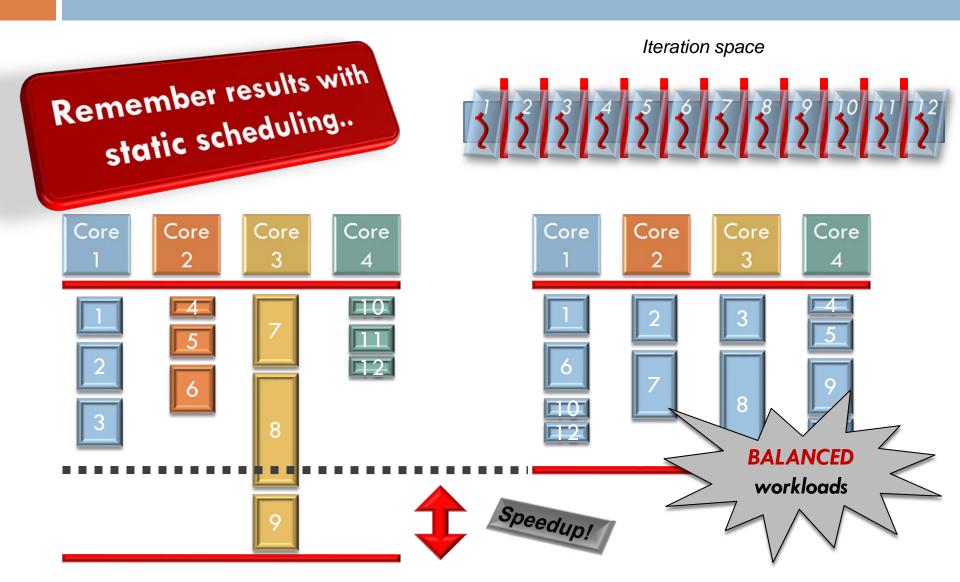
```
#pragma omp for schedule(dtatmcc)
    for (i=0; i<12; i++)
        int start = rand();
        int count = 0;
        while (start++ < 256)
          count++;
        a[count] = foo();
```

Iteration space



A thread is generated for every single iteration

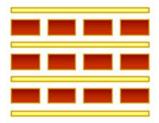




#### Parallelization granularity

#### Iteration chunking

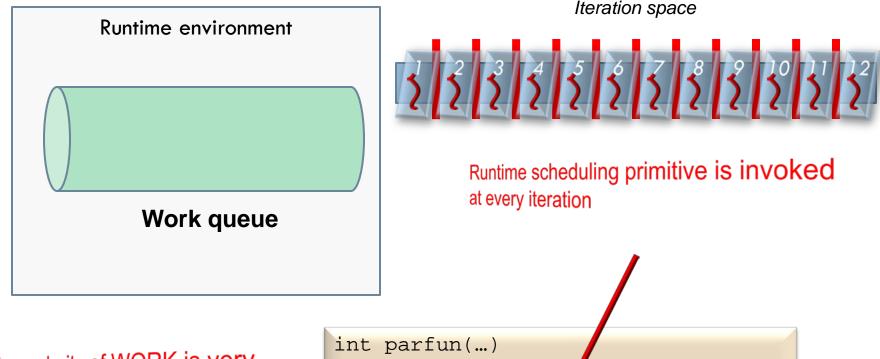
- Fine-grain Parallelism
  - Best opportunities for load balancing, but..
  - Small amounts of computational work between parallelism computation stages
  - Low computation to parallelization ratio → High parallelization overhead



- Coarse-grain Parallelism
  - Harder to load balance efficiently, but..
  - Large amounts of computational work between parallelism computation stages
  - High computation to parallelization ratio → Low parallelization overhead
  - Harder to load balance



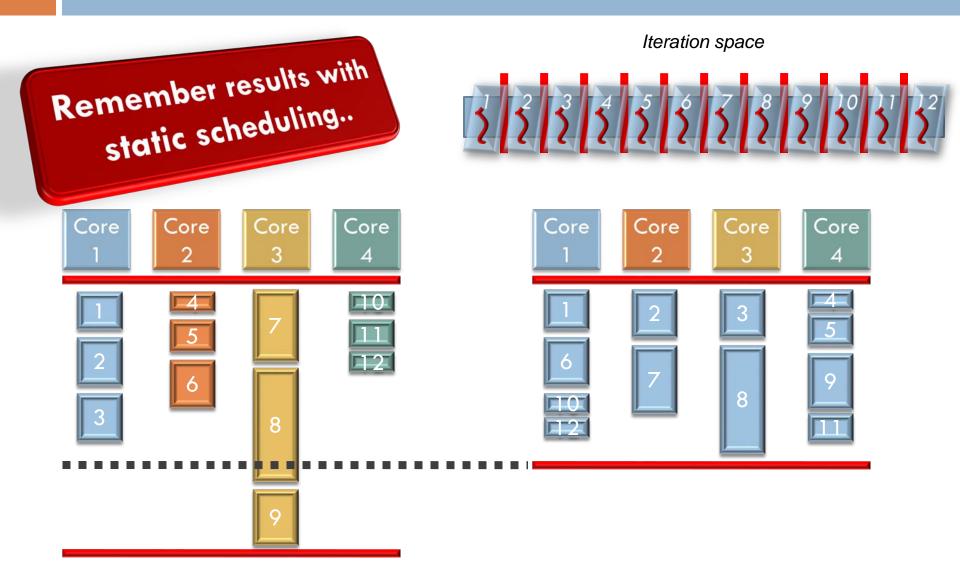
```
Iteration space
#pragma omp for schedule(dynamic, 1)
    for (i=0; i<12; i++)
                                  3 2 3 4 5 6 7 8 8 2 2 2 2
        int start = rand();
        int count = 0;
       while (start++ < 256)
          count++;
       a[count] = foo();
```

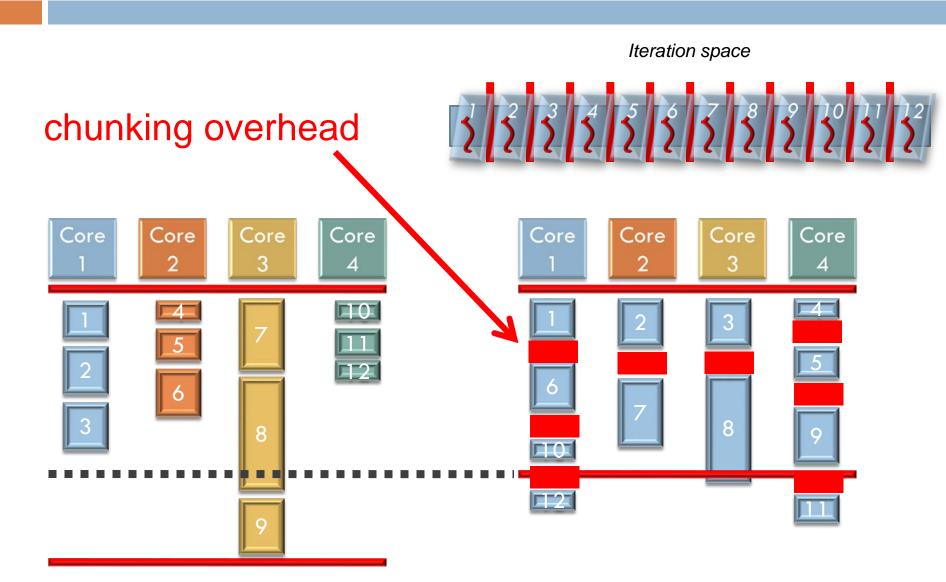


If granularity of WORK is very small the overhead for fine chunking is significant

```
int parfun(...)
{
    int IB, UB:
    GOMP_loop_dynamic_next(&LB, &UB);

for (i=LB; i<UB; i++) {WORK}
}</pre>
```



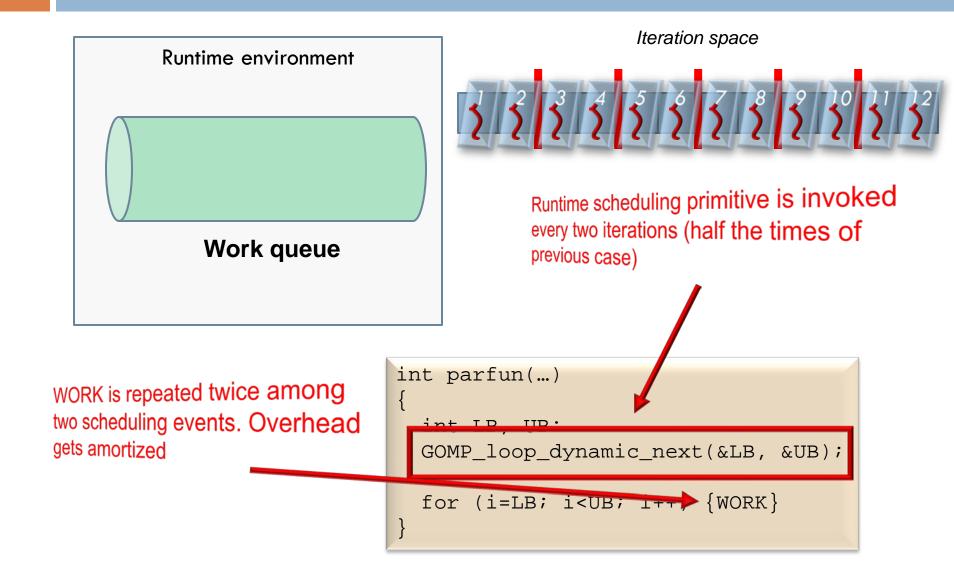


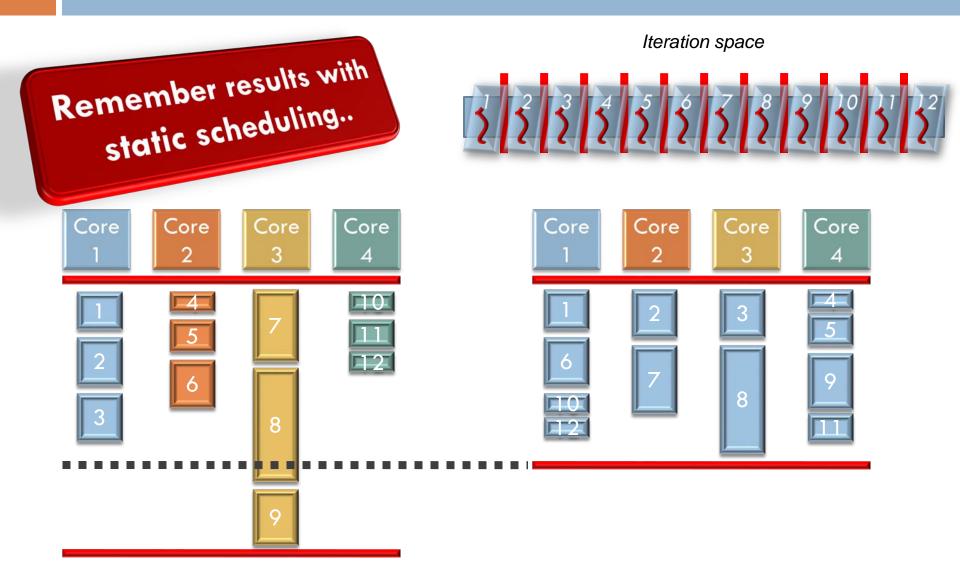
```
#pragma omp for schedule(dynamic, 2)

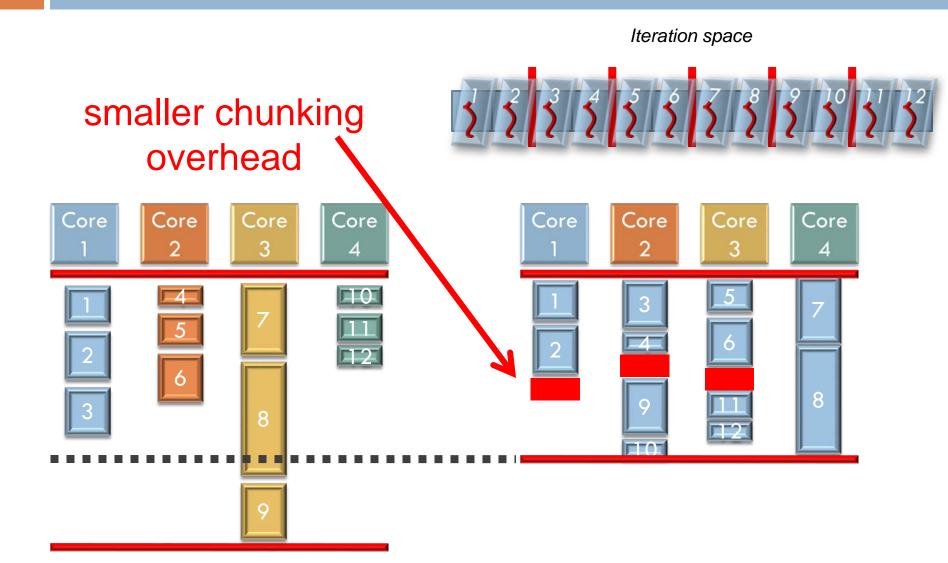
{
    for (i=0; i<12; i++)
        {
        int start = rand();
        int count = 0;

        while (start++ < 256)
        count++;

        a[count] = foo();
    }
```







#### More scheduling clauses

- schedule (guided[, chunk])
  - Threads dynamically grab blocks of iterations. The size of the block starts large and shrinks down to size "chunk" as the calculation proceeds.
- schedule (runtime[, chunk])
  - Schedule and chunk size taken from the OMP\_SCHEDULE environment variable (or the runtime library ... for OpenMP 3.0)

# Sharing work among threads The sections directive

- The for pragma allows to exploit data parallelism in loops
- OpenMP also provides directives to exploit task parallelism

#pragma omp sections

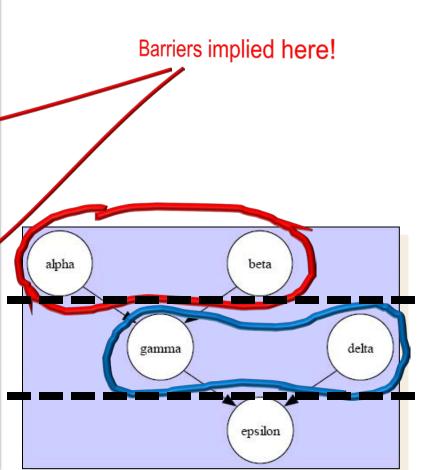
#### SPMD VS MPMD

#### Recall..

- SPMD (single program, multiple data)
  - Processors execute the same stream of instructions over different data
  - □ #pragma omp for
- MPMD (multiple program, multiple data)
  - Processors execute different streams of instructions over (possibly) different data
  - #pragma omp sections
  - □ #pragma omp task

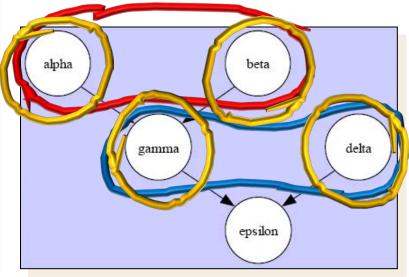
```
int main()
                                                      Identify independent nodes
                                                      in the task graph, and outline
       v = alpha();
                                                      parallel computation with the
                                      SOLUTION
                                                          sections directive
       w = beta();
       y = delta();
                                               alpha
                                                                   beta
       x = gamma(v, w);
       z = epsilon(x, y);
                                                                             delta
                                                         gamma
  printf ("%f\n'', z);
                                                                   epsilon
```

```
int main()
#pragma omp parallel sections {
     v = alpha();
     w = beta();
#pragma omp parallel sections {
     y = delta();
     x = gamma(v, w);
     z = epsilon(x, y);
 printf ("%f\n", z);
```



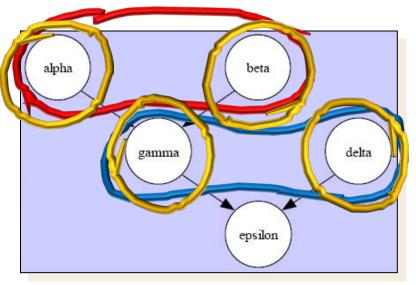
```
int main()
#pragma omp parallel sections {
     v = alpha();
     w = beta();
#pragma omp parallel sections {
     y = delta();
     x = gamma(v, w);
      z = epsilon(x, y);
 printf ("%f\n'', z);
```

Each parallel task within a sections block identifies a section



```
int main()
#pragma omp parallel sections {
 #pragma omp section
     v = alpha();
 #pragma omp section
     w = beta();
#pragma omp parallel sections {
 #pragma omp section
     y = delta();
 #pragma omp section
     x = gamma(v, w);
      z = epsilon(x, y);
 printf ("%f\n", z);
```

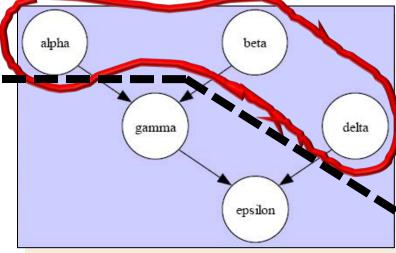
Each parallel task within a sections block identifies a section



```
int main()
                                                      Identify independent nodes
                                                      in the task graph, and outline
       v = alpha();
                                                      parallel computation with the
                                      SECOND
                                      SOLUTION
                                                          sections directive
       w = beta();
       y = delta();
                                               alpha
                                                                   beta
       x = gamma(v, w);
       z = epsilon(x, y);
                                                                             delta
                                                         gamma
  printf ("%f\n'', z);
                                                                  epsilon
```

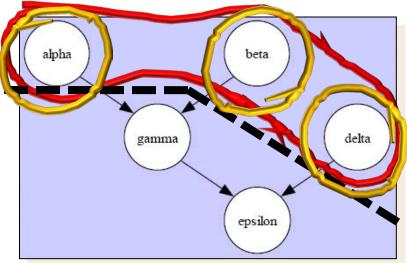
```
int main()
#pragma omp parallel sections {
     v = alpha();
     w = beta();
     y = delta();
     x = gamma(v, w);
     z = epsilon(x, y);
 printf ("%f\n", z);
```

Barrier implied here!



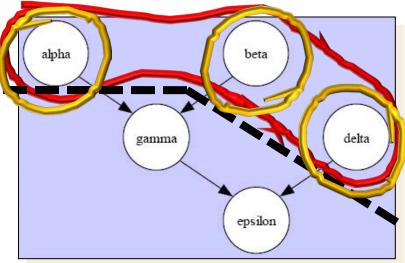
```
int main()
#pragma omp parallel sections {
     v = alpha();
     w = beta();
     y = delta();
     x = gamma(v, w);
     z = epsilon(x, y);
 printf ("%f\n", z);
```

Each parallel task within a sections block identifies a section



```
int main()
#pragma omp parallel sections {
   #pragma omp section
     v = alpha();
   #pragma omp section
      w = beta ();
   #pragma omp section
      y = delta();
      x = qamma(v, w);
      z = epsilon(x, y);
 printf ("%f\n'', z);
```

Each parallel task within a sections block identifies a section



## #pragma omp barrier

- Most important synchronization mechanism in shared memory fork/join parallel programming
- All threads participating in a parallel region wait until everybody has finished before computation flows on
- This prevents later stages of the program to work with inconsistent shared data
- It is implied at the end of parallel constructs, as well as for and sections (unless a nowait clause is specified)

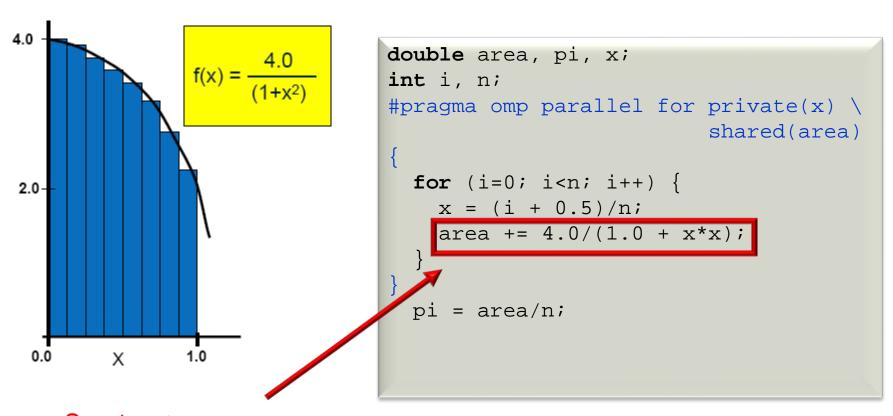
# #pragma omp critical

- Critical Section: a portion of code that only one thread at a time may execute
- □ We denote a critical section by putting the pragma

#pragma omp critical

in front of a block of C code

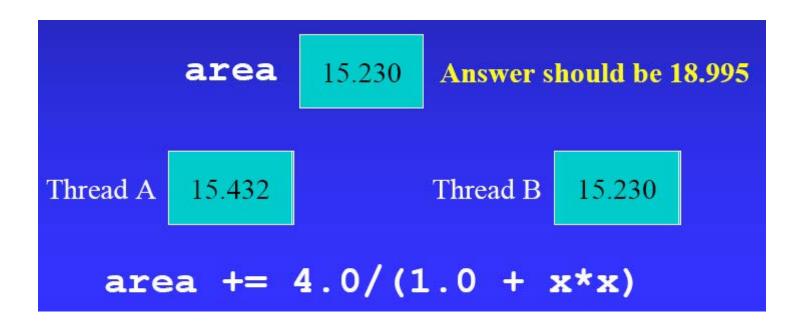
## $\pi$ -finding code example



Synchronize accesses to shared variable **area** to avoid inconsistent results

#### Race condition

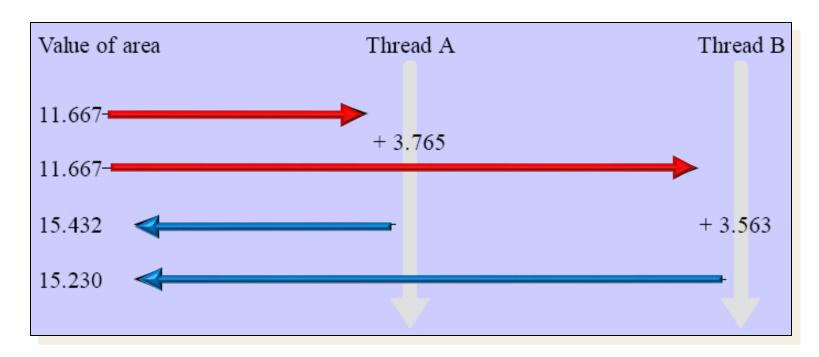
Ensure atomic updates of the shared variable area to avoid a race condition in which one process may "race ahead" of another and ignore changes



## Race condition (Cont'd)

- Thread A reads "11.667" into a local register
- Thread B reads "11.667" into a local register
- Thread A updates area with "11.667+3.765"
- Thread B ignores write from thread A and updates area with "11.667 + 3.563"





## $\pi$ -finding code example

```
double area, pi, x;
int i, n;
#pragma omp parallel for private(x) shared(area)
{
   for (i=0; i<n; i++) {
      x = (i +0.5)/n;
#pragma omp critical
      area += 4.0/(1.0 + x*x);
   }
}
pi = area/n;</pre>
```

**#pragma omp critical** protects the code within its scope by acquiring a lock before entering the critical section and releasing it after execution

### Correctness, not performance!

- As a matter of fact, using locks makes execution sequential
- To dim this effect we should try use fine grained locking (i.e. make critical sections as small as possible)
- A simple instruction to compute the value of area in the previous example is translated into many more simpler instructions within the compiler!
- The programmer is not aware of the real granularity of the critical section

```
_suif_start.omp_fn.0 (.omp_data_i)
<bb 2>:
  D.2605 = .omp_data_i - n;
  D.2615 = __builtin_omp_get_num_threads ();
   D.2616 = __builtin_omp_get_thread_num ();
   D.2623 = MIN_EXPR < D.2622, D.2605>;
   if (D.2621 >= D.2623) goto (L4); else goto (L2);
  (L4):;
    return;
  (12):
    D.2586 = (double) i;
     D.2587 = D.2586 + 5.0e-1;
     D.2605 = .omp_data_i \rightarrow n;
     D.2606 = D.2605;
     D.2588 = (double) D.2606;
      __builtin_GOMP_critical_start ();
     x = D.2587 / D.2588;
      D.2589 = x * x;
      D.2590 = D.2589 + 1.0e+0;
      D.2591 = 4.0e+0 / D.2590;
      D.2607 = .omp_data_i \rightarrow area;
      D.2608 = D.2607;
       D.2609 = D.2591 + D.2608;
       .onp_data_i->area = 1.2609;
       _builtin_GOMP_critical_end ();
        if (D.2627) goto (12); else goto (14);
```

#### mance!

kecution sequential

This is a dump of the intermediate representation of the program within the compiler

e.

3.0-1

```
_suif_start.omp_fn.0 (.omp_data_i)
<bb 2>:
  D.2605 = .omp_data_i - n;
   D.2615 = __builtin_omp_get_num_threads ();
   D.2616 = __builtin_omp_get_thread_num ();
   D.2623 = MIN_EXPR < D.2622, D.2605>;
   if (D.2621 >= D.2623) goto (L4); else goto (L2);
  (L4):;
    return;
  (12):
     D.2586 = (double) i;
     D.2587 = D.2586 + 5.0e-1;
     D.2605 = .omp_data_i \rightarrow n;
     D.2606 = D.2605;
     D.2588 = (double) D.2606;
      _builtin_GOMP_critical_start ();
     x = D.2587 / D.2588;
      D.2589 = x * x;
        2590 = D.2589 + 1.0e+0;
         2591 = 4.0e+0 / D.2590;
        2607 = .omp_data_i->area;
       D.2608 = D.2607;
       D.2609 = D.2591 + D.2608;
       .onp_data_i->area = D.2609;
       _builtin_GOMP_critical_end ();
        if (D.2627) goto (12); else goto (14);
```

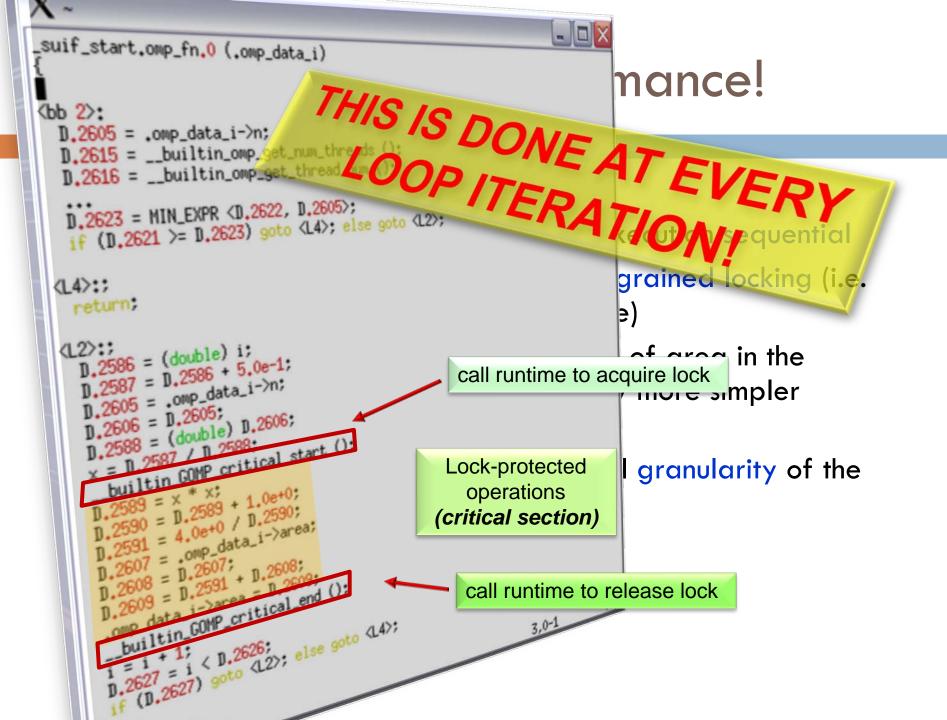
#### mance!

kecution sequential grained locking (i.e. a)
of area in the more simpler

granularity of the

This is how the compiler represents the instruction

area 
$$+= 4.0/(1.0 + x*x);$$



## $\pi$ -finding code example

```
double area, pi, x;
int i, n;
#pragma omp parallel for \
             private(x)
             shared(area)
  for (i=0; i<n; i++) {</pre>
    x = (i + 0.5)/n;
                          Parallel
  #pragma omp critical
    area += 4.0/(1.0 + x*x);
                        Sequential
  pi = area/n;
        Waiting for lock
```

## Correctness, not performance!

- A programming pattern such as area += 4.0/(1.0 + x\*x); in which we:
  - Fetch the value of an operand
  - Add a value to it
  - Store the updated value

is called a reduction, and is commonly supported by parallel programming APIs

 OpenMP takes care of storing partial results in private variables and combining partial results after the loop

## Correctness, not performance!

```
double area, pi, x;
int i, n;
#pragma omp parallel for private(x) shared(area) reduction(+:area)
{
  for (i=0; i<n; i++) {
    x = (i +0.5)/n;
    area += 4.0/(1.0 + x*x);
  }
}
pi = area/n;</pre>
```

The **reduction** clause instructs the compiler to create **private** copies of the **area** variable for every thread. At the end of the loop partial sums are combined on the shared **area** variable

```
_suif_start.omp_fn.0 (.omp_data_i)
        ⟨bb 2⟩:
          area = 0.0;
          D.2605 = .omp_data_i > n;
           D.2615 = __builtin_omp_get_num_threads ();
           D.2616 = __builtin_omp_get_thread_num ();
           D.2623 = MIN_EXPR <D.2622, D.2605>;
           if (D.2621 >= D.2623) goto (L4); else goto (L2);
doul
int
              _builtin_GOMP_atomic_start ();
          <L4>::
#pra
            D.2607 = &.omp_data_i-\rangle area;
             D.2608 = *D.2607;
   fo
             D.2609 = D.2608 + area;
             *D.2607 = D.2609;
             __builtin_GOMP_atomic_end ();
             return;
                   6 = (double) i;
                   87 = D.2586 + 5.0e-1;
                   05 = .omp_data_i - n;
   рi
                   06 = D.2605;
                   88 = (double) D.2606;
                   D.2587 / D.2588;
                  2590 = D.2589 + 1.0e+0;
               D.2591 = 4.0e+0 / D.2590;
              i = i + 1;

D.2626; else goto (14);

if (D.2627) goto (12); else goto (14);
  The r
  the ai
  comb
```

#### rformance!

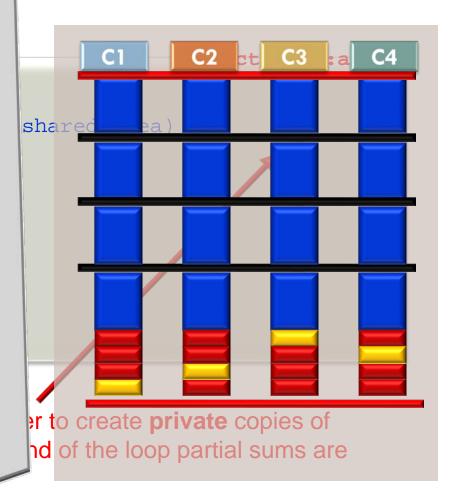
Shared variable is only updated at the end of the loop, when partial sums are computed shared(area) reduction(+:area) Each thread computes partial sums on private copies of the reduction variable

er to create **private** copies of and of the loop partial sums are

```
_suif_start.omp_fn.0 (.omp_data_i)
                                <br >
                                         area = 0.0;
                                          D.2605 = .omp_data_i->n;
                                           D.2615 = __builtin_omp_get_num_threads ();
                                            D.2616 = __builtin_omp_get_thread_num ();
                                             D.2623 = MIN_EXPR <D.2622, D.2605>;
                                             if (D.2621 >= D.2623) goto (L4); else goto (L2);
doul
int
                                                       _builtin_GOMP_atomic_start ();
                                       <L4>::
#pra
                                                 D.2607 = \&.omp_data_i \rightarrow area;
                                                  D.2608 = *D.2607;
             fo
                                                  D.2609 = D.2608 + area;
                                                   *D.2607 = D.2609;
                                                  __builtin_GOMP_atomic_end ();
                                                    return;
                             6
                                                              2586 = (double) i;
                                                                2587 = D.2586 + 5.0e-1;
                                                                   2605 = .omp_data_i - n;
             рi
                                                                 2588 = (double) D.2606;
                                                                    7606 = D.2605;
                                                                         D.2587 / D.2588;
                                                          D.2590 = D.2589 + 1.0e+0;
                                                           D.2591 = 4.0e+0 / D.2590;
       The r
                                                        i = i + 1; \ D.2626; else goto \((14\));

if \((10.2627)\) goto \((12\)); else goto \((14\));
        the ar
         comb
```

#### rformance!



#### More worksharing constructs

#### The master directive

- The master construct denotes a structured block that is only executed by the master thread.
- The other threads just skip it (no synchronization is implied).

```
#pragma omp parallel
{
   do_many_things();
   #pragma omp master
   { exchange_boundaries(); }

   #pragma omp barrier
   do_many_other_things();
}
```

#### More worksharing constructs

#### The single directive

- The single construct denotes a block of code that is executed by only one thread (not necessarily the master thread).
- A barrier is implied at the end of the single block (can remove the barrier with a nowait clause).

```
#pragma omp parallel
{
   do_many_things();
   #pragma omp single
   { exchange_boundaries(); }

   #pragma omp barrier
   do_many_other_things();
}
```

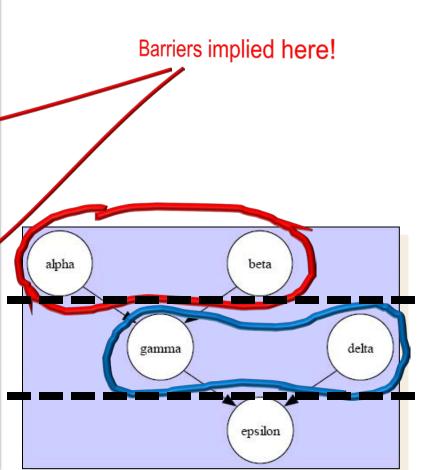
# Recap: TASK parallelism in OpenMP 2.5 The sections directive

- The for pragma allows to exploit data parallelism in loops
- OpenMP 2.5 also provides a directive to exploit task parallelism

#pragma omp sections

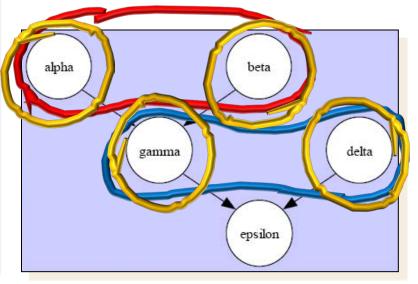
```
int main()
                                                      Identify independent nodes
                                                      in the task graph, and outline
       v = alpha();
                                                      parallel computation with the
                                      SOLUTION
                                                         sections directive
       w = beta();
       y = delta();
                                               alpha
                                                                   beta
       x = gamma(v, w);
       z = epsilon(x, y);
                                                                             delta
                                                         gamma
  printf ("%f\n", z);
                                                                  epsilon
```

```
int main()
#pragma omp parallel sections {
     v = alpha();
     w = beta();
#pragma omp parallel sections {
     y = delta();
     x = gamma(v, w);
     z = epsilon(x, y);
 printf ("%f\n", z);
```



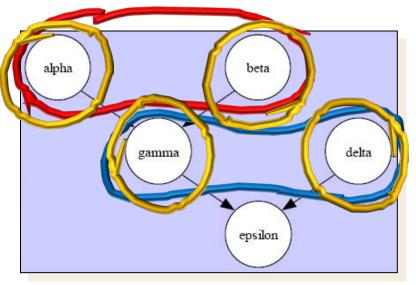
```
int main()
#pragma omp parallel sections {
     v = alpha();
     w = beta();
#pragma omp parallel sections {
     y = delta();
     x = gamma(v, w);
     z = epsilon(x, y);
 printf ("%f\n", z);
```

Each parallel task within a sections block identifies a section



```
int main()
#pragma omp parallel sections {
 #pragma omp section
     v = alpha();
 #pragma omp section
     w = beta();
#pragma omp parallel sections {
 #pragma omp section
     y = delta();
 #pragma omp section
     x = gamma(v, w);
      z = epsilon(x, y);
 printf ("%f\n", z);
```

Each parallel task within a sections block identifies a section



 The sections directive allows a very limited form of task parallelism

- All tasks must be statically outlined in the code
  - What if a functional loop (while) body is identified as a task?
    - Unrolling? Not feasible for high iteration count
  - What if recursion is used?

- □ Why?
  - Example: list traversal

#### EXAMPLE

```
void traverse_list (List 1)
{
   Element e ;
#pragma omp parallel private ( e )
   for ( e = e first; e; e = e next )
#pragma omp single nowait
     process ( e ) ;
}
```

#### OpenMP v2.5

- Awkward!
- Poor performance
- Not composable

- Why?
  - Example: tree traversal

#### **EXAMPLE**

```
void traverse tree (Tree *tree)
#pragma omp parallel sections
#pragma omp section
  if ( tree→left )
    traverse_tree ( tree > left );
#pragma omp section
  if ( tree→right)
    traverse_tree ( tree > right);
  process (tree);
```

#### OpenMP v2.5

- Too many parallel regions
  - Extra overheads
  - Extra synchronizations
  - Not always well supported

- Better solution for those problems
- Main addition to OpenMP 3.0a
- Allows to parallelize irregular problems
  - unbounded loops
  - recursive algorithms
  - producer/consumer schemes
  - **-** ...

- □ The OpenMP tasking model
  - Creating tasks
  - Data scoping
  - Syncronizing tasks
  - Execution model

# What is an OpenMP task?

- Tasks are work units which execution may be deferred
  - they can also be executed immediately!
- Tasks are composed of:
  - code to execute
  - data environment
    - Initialized at creation time
  - internal control variables (ICVs)

## Task directive

```
#pragma omp task [ clauses ] structured block
```

- Each encountering thread creates a task
  - Packages code and data environment
- Highly composable. Can be nested
  - inside parallel regions
  - inside other tasks
  - inside worksharing constructs (for, sections)

- □ Why?
  - Example: list traversal

#### **EXAMPLE**

```
void traverse_list (List 1)
{
   Element e ;

   for ( e = e first; e; e = e next )
#pragma omp task
     process ( e ) ;
}
What is 1
```

What is the scope of e?

# Task data scoping

- Data scoping clauses
  - □ shared(list)
  - private(list)
  - firstprivate(list)
    - data is captured at creation
  - default(shared | none)

## Task data scoping

when there are no clauses...

- ☐ If no clause
  - Implicit rules apply
    - e.g., global variables are shared
- Otherwise...
  - firstprivate
  - shared attribute is lexically inherited

#### **EXAMPLE**

```
int a ;
void foo ( ) {
int b , c ;
#pragma omp parallel shared(c)
  int d;
#pragma omp task
    int e ;
    a = shared
   b = firstprivate
   c = shared
    d = firstprivate
    e = private
```

### Tip

default(none) is your friend Use it if you do not see it clear

#### **EXAMPLE**

```
void traverse_list (List 1)
{
   Element e ;

   for ( e = e first; e; e = e next )
#pragma omp task
     process ( e ) ;
}
e is firstprivate
```

```
void traverse_list (List 1)
{
   Element e ;

   for ( e = e → first; e; e = e → next )
#pragma omp task
     process ( e ) ;
}
```

how we can guarantee here that the traversal is finished?

# Task synchronization

- Barriers (implicit or explicit)
  - All tasks created by any thread of the current team are guaranteed to be completed at barrier exit
- Task barrier

### #pragma omp taskwait

- Encountering task suspends until child tasks complete
  - Only direct childs, not descendants!

#### **EXAMPLE**

```
void traverse_list (List 1)
{
   Element e ;

   for ( e = e→first; e; e = e→next )
#pragma omp task
     process ( e ) ;

#pragma omp taskwait
}
```

All tasks guaranteed to be completed here

## Task execution model

- Task are executed by a thread of the team that generated it
  - Can be executed immediately by the same thread that creates it
- Parallel regions in 3.0 create tasks!
  - One implicit task is created for each thread
    - So all task-concepts have sense inside the parallel region
- Threads can suspend the execution of a task and start/resume another

- □ Why?
  - Example: list traversal

### **CAREFUL!**

• Multiple traversal of the same list

#### **EXAMPLE**

```
List 1;
#pragma omp parallel
traverse_list (1);
```

```
void traverse_list (List 1)
{
   Element e ;

   for ( e = e→first; e; e = e→next )
#pragma omp task
     process ( e ) ;
}
```

- Why?
  - Example: list traversal

#### **EXAMPLE**

```
#pragma omp parallel
#pragma omp single
traverse_list (1);
```

### Single traversal

- One thread enters single and creates all tasks
- All the team cooperates executing them

```
void traverse_list (List 1)
{
   Element e ;

   for ( e = e→first; e; e = e→next )
#pragma omp task
     process ( e ) ;
}
```

In case task is within a regular counted loop an alternative is to parallelize task creation among threads

#### **EXAMPLE**

```
/* A DIFFERENT EXAMPLE */
#pragma omp parallel
Myfunc ();
```

### **Multiple traversals**

- Multiple threads create tasks
- All the team cooperates executing them

```
void Myfunc ()
{
  int i;
#pragma omp for
  for (i=LB; i<UB; i++)
#pragma omp task
    process ( i ) ;
}</pre>
```

# Task scheduling

#### How it works?

- Tasks are tied by default
  - Tied tasks are executed always by the same thread
  - Tied tasks have scheduling restrictions
    - Deterministic scheduling points (creation, synchronization, ...)
    - Another constraint to avoid deadlock problems
  - Tied tasks may run into performance problems
- Programmer can use untied clause to lift all restrictions
  - Note: Mix very carefully with threadprivate, critical and thread-ids

## And last..

### The IF clause

- If the expression of a if clause is false
  - The encountering task is suspended
  - The new task is executed immediately
    - with its own data environment
    - different task with respect to synchronization
  - The parent task resumes when the task finishes
  - Allows implementations to optimize task creation

```
#pragma omp task if (...)
process ( e );
```

# EXERCISE