Outline



- Introduction to OpenMP
- Creating Threads
- Synchronization
- Parallel Loops
- Data Environment
- Memory Model



- Irregular Parallelism and Tasks
- Recap
- Beyond the Common Core:
 - Worksharing Revisited
 - Synchronization Revisited: Options for Mutual exclusion
 - Thread Affinity and Data Locality
 - Thread Private Data
 - Memory Models and Point-to-Point Synchronization
 - Programming your GPU with OpenMP

Irregular Parallelism

- Let's call a problem "irregular" when one or both of the following hold:
 - Data Structures are sparse
 - Control structures are not basic for-loops
- Example: Traversing Linked lists:

```
p = listhead;
while (p) {
  process(p);
  p=p->next;
}
```

Using what we've learned so far, traversing a linked list in parallel using OpenMP is difficult.

Exercise: Traversing linked lists

- Consider the program linked.c
 - Traverses a linked list computing a sequence of Fibonacci numbers at each node.
- Parallelize this program selecting from the following list of constructs:

```
#pragma omp parallel
#pragma omp for
#pragma omp parallel for
#pragma omp for reduction(op:list)
#pragma omp critical
int omp_get_num_threads();
int omp_get_thread_num();
double omp_get_wtime();
schedule(static[,chunk]) or schedule(dynamic[,chunk])
private(), firstprivate(), default(none)
```

 Hint: Just worry about the while loop that is timed inside main(). You don't need to make any changes to the "list functions"

Linked Lists with OpenMP (without tasks)

See the file solutions/linked_notasks.c

```
while (p != NULL) {
   p = p-next;
                                                            Count number of items in the linked list
   count++;
struct node *parr = (struct node*) malloc(count*sizeof(struct node));
p = head;
for(i=0; i<count; i++) {
                                                            Copy pointer to each node into an array
   parr[i] = p;
   p = p - next;
#pragma omp parallel
   #pragma omp for schedule(static,1)
                                                            Process nodes in parallel with a for loop
   for(i=0; i<count; i++)
                                                             Number of
                                                                                      Schedule
     processwork(parr[i]);
                                                             threads
                                                                                Default
                                                                                              Static, 1
                                                                           48 seconds
                                                                                              45 seconds
```

28 seconds

39 seconds

Linked Lists with OpenMP (without tasks)

See the file solutions/linked_notasks.c

```
while (p != NULL) {
   p = p-next;
    count++;
struct node *parr = (struct node*) malloc(count*sizeof(struct node)):
p = head:
for(i=0; i<count; i++) {
    parr[i] = p;
    p = p - next;
#pragma omp parallel
   #pragma omp for schedule(static,1)
   for(i=0; i<count; i++)
     processwork(parr[i]);
```

Count number of items in the linked list

With so much code to add and three passes through the data, this is really ugly.

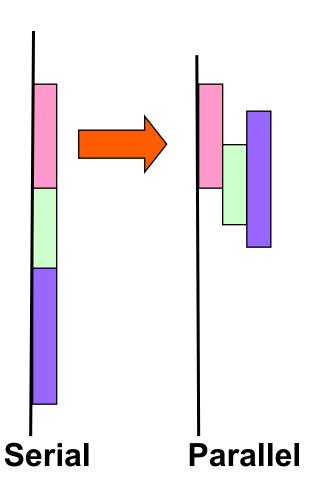
There has got to be a better way to do this

Process nodes in parallel with a for loop

Number of threads	Schedule	
	Default	Static,1
1	48 seconds	45 seconds
2	39 seconds	28 seconds

What are Tasks?

- Tasks are independent units of work
- Tasks are composed of:
 - code to execute
 - data to compute with
- Threads are assigned to perform the work of each task.
 - The thread that encounters the task construct may execute the task immediately.
 - The threads may defer execution until later

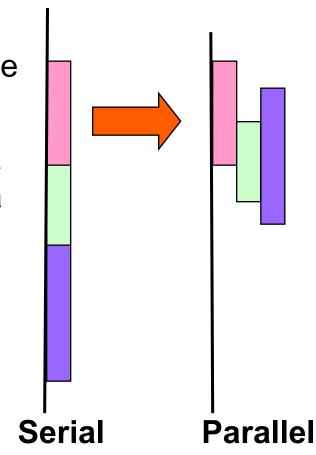


What are Tasks?

The task construct includes a structured block of code

 Inside a parallel region, a thread encountering a task construct will package up the code block and its data for execution

 Tasks can be nested: i.e. a task may itself generate tasks.



A common Pattern is to have one thread create the tasks while the other threads wait at a barrier and execute the tasks

Single Worksharing Construct

- The single construct denotes a block of code that is executed by only one thread (not necessarily the primary* 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(); }
          do_many_other_things();
}
```

^{*}This used to be called the "master thread". The term "master" has been deprecated in OpenMP 5.1 and replaced with the term "primary".

Task Directive

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

```
Create some threads
#pragma omp parallel ← — —
  #pragma omp single __
                                        One Thread
                                        packages tasks
      #pragma omp task
          fred();
      #pragma omp task
                                        Tasks executed by
          daisy();
                                        some thread in some
      #pragma omp task
                                        order
         billy();
              All tasks complete before this barrier is released
```

Exercise: Simple tasks

- Write a program using tasks that will "randomly" generate one of two strings:
 - "I think " "race" "car" "s are fun"
 - "I think " "car" "race" "s are fun"
- Hint: use tasks to print the indeterminate part of the output (i.e. the "race" or "car" parts).
- This is called a "Race Condition". It occurs when the result of a program depends on how the OS schedules the threads.
- NOTE: A "data race" is when threads "race to update a shared variable". They produce race conditions. Programs containing data races are undefined (in OpenMP but also ANSI standards C++'11 and beyond).

```
#pragma omp parallel#pragma omp task#pragma omp single
```

Racey Cars: Solution

```
#include <stdio.h>
#include <omp.h>
int main()
{ printf("I think");
 #pragma omp parallel
   #pragma omp single
     #pragma omp task
       printf(" car");
     #pragma omp task
       printf(" race");
 printf("s");
 printf(" are fun!\n");
```

Data Scoping with Tasks

- Variables can be shared, private or firstprivate with respect to task
- These concepts are a little bit different compared with threads:
 - If a variable is shared on a task construct, the references to it inside the construct are to the storage with that name at the point where the task was encountered
 - If a variable is private on a task construct, the references to it inside the construct are to new uninitialized storage that is created when the task is executed
 - If a variable is firstprivate on a construct, the references to it inside the construct are
 to new storage that is created and initialized with the value of the existing storage of
 that name when the task is encountered

Data Scoping Defaults

- The behavior you want for tasks is usually firstprivate, because the task may not be executed until later (and variables may have gone out of scope)
 - Variables that are private when the task construct is encountered are firstprivate by default
- Variables that are shared in all constructs starting from the innermost enclosing parallel construct are shared by default

Exercise: Traversing linked lists

- Consider the program linked.c
 - Traverses a linked list computing a sequence of Fibonacci numbers at each node.
- Parallelize this program selecting from the following list of constructs:

```
#pragma omp parallel
#pragma omp single
#pragma omp task
int omp_get_num_threads();
int omp_get_thread_num();
double omp_get_wtime();
private(), firstprivate()
```

Hint: Just worry about the contents of main(). You don't need to make any changes to the "list functions"

Parallel Linked List Traversal

```
Only one thread
                                        packages tasks
#pragma omp parallel
  #pragma omp single*
    p = listhead ;
    while (p) {
        #pragma omp task firstprivate(p)
                 process (p);
        p=next (p) ;
                                     makes a copy of p
                                     when the task is
                                      packaged
```

When/Where are Tasks Complete?

- At thread barriers (explicit or implicit)
 - all tasks generated inside a region must complete at the next barrier encountered by the threads in that region. Common examples:
 - Tasks generated inside a single construct: all tasks complete before exiting the barrier on the single.
 - Tasks generated inside a parallel region: all tasks complete before exiting the barrier at the end of the parallel region.

At taskwait directive

i.e. Wait until all tasks defined in the current task have completed.

```
#pragma omp taskwait
```

- Note: applies only to tasks generated in the current task, not to "descendants".

Example

```
#pragma omp parallel
  #pragma omp single
                                        fred() and daisy()
      #pragma omp task
                                        must complete before
          fred();
                                        billy() starts, but
      #pragma omp task
                                        this does not include
          daisy();
                                        tasks created inside
      #pragma omp taskwait
                                        fred() and daisy()
      #pragma omp task
          billy();
                          All tasks including those created
                          inside fred() and daisy() must
                          complete before exiting this barrier
```

Example

```
#pragma omp parallel
  #pragma omp single nowait
     #pragma omp task
        fred();
     #pragma omp task
        daisy();
     #pragma omp taskwait
     #pragma omp task
        billy();
```

The barrier at the end of the single is expensive and not needed since you get the barrier at the end of the parallel region. So use nowait to turn it off.

All tasks including those created inside fred() and daisy() must complete before exiting this barrier

Example: Fibonacci numbers

```
int fib (int n)
 int x,y;
 if (n < 2) return n;
 x = fib(n-1);
  y = fib (n-2);
  return (x+y);
Int main()
  int NW = 5000;
 fib(NW);
```

- $F_n = F_{n-1} + F_{n-2}$
- Inefficient O(n²) recursive implementation!

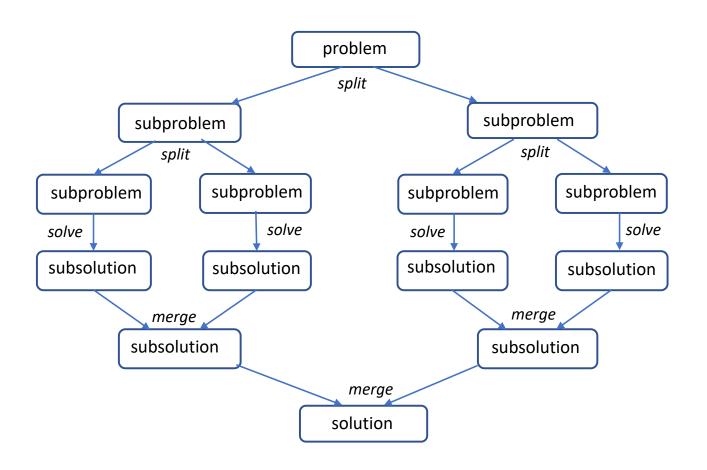
Parallel Fibonacci

```
int fib (int n)
{ int x,y;
 if (n < 2) return n;
#pragma omp task shared(x)
 x = fib(n-1);
#pragma omp task shared(y)
 y = fib (n-2);
#pragma omp taskwait
 return (x+y);
Int main()
\{ \text{ int NW} = 5000; 
 #pragma omp parallel
    #pragma omp single
        fib(NW);
```

- Binary tree of tasks
- Traversed using a recursive function
- A task cannot complete until all tasks below it in the tree are complete (enforced with taskwait)
- x,y are local, and so by default they are private to current task
 - must be shared on child tasks so they don't create their own firstprivate copies at this level!

Divide and Conquer

 Split the problem into smaller sub-problems; continue until the sub-problems can be solved directly



- 3 Options for parallelism:
 - □ Do work as you split into sub-problems
 - □ Do work only at the leaves
 - Do work as you recombine

Exercise: PI with tasks

- Go back to the original pi.c program
 - Parallelize this program using OpenMP tasks

```
#pragma omp parallel
#pragma omp task
#pragma omp taskwait
#pragma omp single
double omp_get_wtime()
int omp_get_thread_num();
int omp_get_num_threads();
```

Hint: first create a recursive pi program and verify that it works. Think about the
computation you want to do at the leaves. If you go all the way down to one
iteration per leaf-node, won't you just swamp the system with tasks?

Program: OpenMP tasks

```
include <omp.h>
static long num_steps = 100000000;
#define MIN BLK 10000000
double pi comp(int Nstart,int Nfinish,double step)
{ int i,iblk;
  double x, sum = 0.0,sum1, sum2;
  if (Nfinish-Nstart < MIN_BLK){</pre>
    for (i=Nstart;i< Nfinish; i++){</pre>
      x = (i+0.5)*step;
      sum = sum + 4.0/(1.0+x*x);
  else{
    iblk = Nfinish-Nstart;
    #pragma omp task shared(sum1)
       sum1 = pi comp(Nstart,
                                    Nfinish-iblk/2,step);
    #pragma omp task shared(sum2)
        sum2 = pi_comp(Nfinish-iblk/2, Nfinish,
                                                  step);
    #pragma omp taskwait
      sum = sum1 + sum2;
 }return sum;
```

```
int main ()
 int i;
 double step, pi, sum;
 step = 1.0/(double) num_steps;
 #pragma omp parallel
    #pragma omp single
      sum =
         pi_comp(0,num_steps,step);
   pi = step * sum;
```

Results*: Pi with tasks

threads	1 st SPMD	SPMD critical	PI Loop	Pi tasks
1	1.86	1.87	1.91	1.87
2	1.03	1.00	1.02	1.00
3	1.08	0.68	0.80	0.76
4	0.97	0.53	0.68	0.52

^{*}Intel compiler (icpc) with no optimization on Apple OS X 10.7.3 with a dual core (four HW thread) Intel® CoreTM i5 processor at 1.7 Ghz and 4 Gbyte DDR3 memory at 1.333 Ghz.

Using Tasks

- Don't use tasks for things already well supported by OpenMP
 - -e.g. standard do/for loops
 - the overhead of using tasks is greater

- Don't expect miracles from the runtime
 - best results usually obtained where the user controls the number and granularity of tasks