

Why is it required?



Recall:

- Need to synchronise actions on shared variables.
- Need to ensure correct ordering of reads and writes.
- Need to protect updates to shared variables (not atomic by default)

BARRIER directive



- No thread can proceed past a barrier until all the other threads have arrived.
- Note that there is an implicit barrier at the end of DO/FOR, SECTIONS and SINGLE directives.

• Syntax:

Fortran: !\$OMP BARRIER

C/C++: #pragma omp barrier

• Either all threads or none must encounter the barrier: otherwise DFADLOCK!!

BARRIER directive (cont)



Example:

```
#pragma omp parallel private(myid,neighb) shared(a,b,c)
{
    myid = omp_get_thread_num();
    neighb = myid - 1;
    if (myid.eq.0) neighb = omp_get_num_threads()-1;
        ...
    a[myid] *= 3.5;
#pragma omp_barrier
    b[myid] = a[neighb] + c;
    ...
}
```

• Barrier required to force synchronisation on a





Example:

```
!$OMP PARALLEL PRIVATE (MYID, NEIGHB) SHARED (A,B,C)
    myid = omp_get_thread_num()
    neighb = myid - 1
    if (myid.eq.0) neighb = omp_get_num_threads()-1
    ...
    a(myid) = a(myid)*3.5
!$OMP BARRIER
    b(myid) = a(neighb) + c
    ...
!$OMP END PARALLEL
```

• Barrier required to force synchronisation on a

Critical sections



- A critical section is a block of code which can be executed by only one thread at a time.
- Can be used to protect updates to shared variables.

CRITICAL directive



• Syntax:

Fortran: !\$OMP CRITICAL

block

!\$OMP END CRITICAL

C/C++: #pragma omp critical

structured block

CRITICAL directive (cont)



Example: appending to a shared list

```
#pragma omp parallel for shared(list, N) private(newitem_p)
for (int i=0; i<N; i++) {
    newitem_p = createitem(i);
#pragma omp critical
    {
        append(&list,p_newitem);
    }
}</pre>
```

CRITICAL directive (cont)



Example: appending to a shared list

```
!$OMP PARALLEL DO SHARED(list,n) PRIVATE(newitem)
do i=1,n
   newitem = createitem(i)
!$OMP CRITICAL
   call append(list,newitem)
!$OMP END CRITICAL
end do
```





Example: pushing and popping a task stack

CRITICAL directive (cont)



Example: pushing and popping a task stack

```
!$OMP PARALLEL SHARED(stack), PRIVATE(next, new, done)
  do while (.not. done)
!$OMP CRITICAL
      next = pop(stack)
!$OMP END CRITICAL
      new = process(next)
!$OMP CRITICAL
      if (valid(new)) call push(new, stack)
      done = isempty(stack)
!$OMP END CRITICAL
      end do
!$OMP END PARALLEL
```

ATOMIC directive



- Used to protect a single update to a shared scalar variable of basic type.
- Applies only to a single statement.
- Syntax:

Fortran: ! \$OMP ATOMIC

statement

where statement must have one of these forms:

```
x = x op expr, x = exprop x, x = intr (x, expr) or x = intr (expr, x) op is one of +, *, -, /, .and., .or., .eqv., or .neqv. intr is one of MAX, MIN, IAND, IOR or IEOR
```

ATOMIC directive (cont)



where statement must have one of the forms:

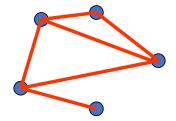
$$x \ binop = \ expr, x++, ++x, x--, \text{ or } --x$$
 and $binop$ is one of +, *, -, /, &, ^, <<, or >>

- Note that the evaluation of expr is not atomic.
- May be more efficient than using CRITICAL directives, e.g. if different array elements can be protected separately.
- No interaction with CRITICAL directives

ATOMIC directive (cont)



Example (compute degree of each vertex in a graph):



Lock routines



- Occasionally we may require more flexibility than is provided by CRITICAL directive.
- A lock is a special variable that may be set by a thread. No other thread may set the lock until the thread which set the lock has unset it.
- Setting a lock can either be blocking or non-blocking.
- A lock must be initialised before it is used, and may be destroyed when it is not longer required.
- Lock variables should not be used for any other purpose.

Lock routines - syntax



```
Fortran:
```

```
USE OMP_LIB

SUBROUTINE OMP_INIT_LOCK(OMP_LOCK_KIND var)

SUBROUTINE OMP_SET_LOCK(OMP_LOCK_KIND var)

LOGICAL FUNCTION OMP_TEST_LOCK(OMP_LOCK_KIND var)

SUBROUTINE OMP_UNSET_LOCK(OMP_LOCK_KIND var)

SUBROUTINE OMP_DESTROY_LOCK(OMP_LOCK_KIND var)
```

var should be an INTEGER of the same size as addresses (e.g. INTEGER*8 on a 64-bit machine)

OMP_LIB defines OMP_LOCK_KIND

Lock routines - syntax

epcc

C/C++:

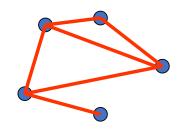
```
#include <omp.h>
  void omp_init_lock(omp_lock_t *lock);
  void omp_set_lock(omp_lock_t *lock);
  int omp_test_lock(omp_lock_t *lock);
  void omp_unset_lock(omp_lock_t *lock);
  void omp_destroy_lock(omp_lock_t *lock);
```

Lock example

Example (compute degree of each vertex in a graph):

```
omp lock t lockvar[nvertices];
for (i=0; i<nvertexes; i++) {</pre>
  omp_init_lock(&lockvar[i]);
#pragma omp parallel for
      for (j=0; j<nedges; j++) {</pre>
         omp_set_lock(&lockvar[edge[j].vertex1]);
           degree[edge[j].vertex1]++;
         omp unset lock(&lockvar[edge[j].vertex1]);
         omp set lock(&lockvar[edge[j].vertex2]);
           degree[edge[j].vertex2]++;
         omp_unset_lock(&lockvar[edge[j].vertex2]);
```





Exercise: Molecular dynamics



- The code supplied is a simple molecular dynamics simulation of the melting of solid argon.
- Computation is dominated by the calculation of force pairs in subroutine **forces**.
- Parallelise this routine using a DO/FOR directive and critical sections.
 - Watch out for PRIVATE and REDUCTION variables.
 - Choose a suitable loop schedule
- Extra exercise: can you improve the performance by using locks, or atomics, or by using a reduction array.

Reusing this material





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