

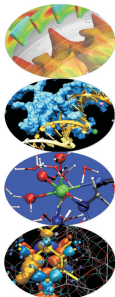
OpenMP Exercises

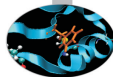
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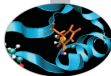
SuperComputing Applications and Innovation Department





Warm-up with OpenMP

- 1 Compile and run "Hello World" and experiment with the `OMP_NUM_THREADS` variable. If any errors occur, try to fix it.
- 2 Parallelize the MM (Matrix Multiplication) serial code acting only on the most important loop



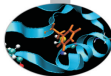
Hello World from C

C

```
#include <stdio.h>
#ifdef _OPENMP
#include<omp.h>
#endif
int main(int argc, char* argv[ ])
{
#ifdef _OPENMP
    int iam;
    #pragma omp parallel \
private(iam) /* the parallel block starts here */
    {
        iam=omp_get_thread_num();

        #pragma omp critical
        printf("Hello from %d\n",iam);

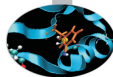
    } /* the parallel block ends here */
#else
    printf("Hello, this is a serial program.\n");
#endif
    return 0;
}
```



Hello World from Fortran

Fortran

```
Program Hello_from_Threads
#ifdef _OPENMP
    use omp_lib
#endif
    implicit none
    integer :: iam
#ifdef _OPENMP
    !$omp parallel &
    !$omp private(iam)
        iam=omp_get_thread_num()
        !$omp critical
            write( *,* ) 'Hello from', iam
        !$omp end critical
    !$omp end parallel
#else
    write( *,* ) 'Hello, this is a serial program'
#endif
end program Hello_from_Threads
```



Matrix Multiplication in C

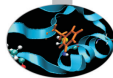
```

C
#include <stdio.h>
#include <stdlib.h>
#include <math.h>

int main(int argc, char **argv) {
    int n;
    int i, j, k;
    ...
    double ( *a ) [n] = malloc(sizeof(double[n][n]));
    double ( *b ) [n] = malloc(sizeof(double[n][n]));
    double ( *c ) [n] = malloc(sizeof(double[n][n]));
    ...

    for (i=0; i<n; i++)
        for (j=0; j<n; j++) {
            a[i][j] = ((double)rand())/((double)RAND_MAX);
            b[i][j] = ((double)rand())/((double)RAND_MAX);
            c[i][j] = 0.0;
        }

    #pragma omp parallel for private(j,k)
        for (i=0; i<n; ++i)
            for (k=0; k<n; k++)
                for (j=0; j<n; ++j)
                    c[i][j] += a[i][k]*b[k][j];
    ...
    return 0;
}
  
```

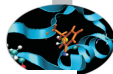


Matrix Multiplication in Fortran

Fortran

```
program mat_prod
  implicit none
  integer :: n
  real(kind(1.d0)), dimension(:,,:), allocatable :: a, b, c
  integer :: i, j, k
  ...
  allocate(a(n,n),b(n,n),c(n,n),stat=ierr)
  ...
  call random_number(a)
  call random_number(b)
  c = 0.d0

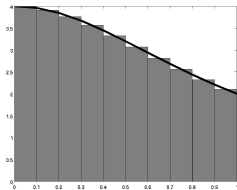
  !$omp parallel do
    do j=1, n
      do k=1, n
        do i=1, n
          c(i,j) = c(i,j) + a(i,k)*b(k,j)
        end do
      end do
    end do
  !$omp end parallel do
  ...
end program mat_prod
```



Let's play with OpenMP

- 3 Parallelize the serial code `Pi`. It computes the Reimann approximation of

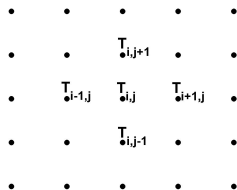
$$\int_0^1 \frac{4}{1+x^2} dx = 4 \arctan x \Big|_0^1 = \pi$$



- 4 Parallelize the serial code `Laplace`. It applies the iterative Jacobi method to a finite differences approximation of the Laplace equation with Dirichlet boundary condition:

$$T_{i,j}^{n+1} = \frac{1}{4}(T_{i+1,j}^n + T_{i-1,j}^n + T_{i,j-1}^n + T_{i,j+1}^n)$$

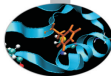
- start from the most computationally intensive loop
- then try to include the `while` loop in the parallel region





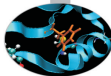
C

```
#include <omp.h>
...
double timel = omp_get_wtime();
sum = 0.0;
dx = 1.0 / (double) intervals;
#pragma omp parallel for private(x,f) reduction(+:sum)
for (i = 1; i <= n; i++) {
    x = dx * ((double) (i - 0.5));
    f = 4.0 / (1.0 + x*x);
    sum = sum + f;
}
pi = dx*sum;
time2 = omp_get_wtime() - timel;
...
```

Fortran

```
use omp_lib
...
time1 = omp_get_wtime()
sum=0.d0
dx=1.d0/intervals
!$omp parallel do private(x,f) reduction(+:sum)
do i=1,n
  x=dx*(i-0.5d0)
  f=4.d0/(1.d0+x*x)
  sum=sum+f
end do
!$omp end parallel do
pi=dx*sum
time2 = omp_get_wtime()
...
```

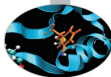


Laplace

```

C
...
while(var > tol && iter <= maxIter) {
  ++iter;
  var = 0.0;
  #pragma omp parallel for private(j) reduction(max:var)
  for (i=1; i<=n; ++i)
    for (j=1; j<=n; ++j) {
      Tnew[i*n2+j] = 0.25*(T[(i-1)*n2+j] + T[(i+1)*n2+j]
                          + T[i*n2+(j-1)] + T[i*n2+(j+1)]);
      var = fmax(var, fabs(Tnew[i*n2+j] - T[i*n2+j]));
    }
  Tmp=T; T=Tnew; Tnew=Tmp;
  if (iter%100 == 0)
    printf("iter: %8u, variation = %12.4lE\n", iter, var);
}
...

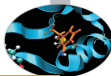
```



Laplace

Fortran

```
...  
do while (var > tol .and. iter <= maxIter)  
  iter = iter + 1  
  var = 0.d0  
  !$omp parallel do reduction(max:var)  
  do j = 1, n  
    do i = 1, n  
      Tnew(i,j)=0.25d0*(T(i-1,j)+T(i+1,j)+T(i,j-1)+T(i,j+1))  
      var = max(var, abs( Tnew(i,j) - T(i,j) ))  
    end do  
  end do  
  !$omp end parallel do  
  Tmp =>T; T =>Tnew; Tnew => Tmp;  
  if( mod(iter,100) == 0 ) ...  
end do  
...
```

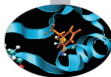


Lacks support for OpenMP 3.1?

C

```

while(var > tol && iter <= maxIter) {
  ++iter;
  var = 0.0;
  #pragma omp parallel
  {
    double pvar = 0.0;
    #pragma omp for private(j)
    for (i=1; i<=n; ++i)
      for (j=1; j<=n; ++j) {
        Tnew[i*n2+j] = 0.25*( T[(i-1)*n2+j] + T[(i+1)*n2+j]
                             + T[i*n2+(j-1)] + T[i*n2+(j+1)] );
        pvar = fmax(pvar , fabs(Tnew[i*n2+j] - T[i*n2+j]));
      }
    #pragma omp critical
    if (pvar > var) var = pvar;
  }
  Tmp=T; T=Tnew; Tnew=Tmp;
  if (iter%100 == 0) ...
}
  
```

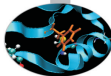


Laplace incremented

C

```
...
#pragma omp parallel
{
  while(var > tol && iter <= maxIter) {
    #pragma omp barrier
    #pragma omp single
    {
      ++iter;
      var = 0.0;
    }
    #pragma omp for private(j) reduction(max:var)
    ...
    #pragma omp single nowait
    {
      Tmp=T; T=Tnew; Tnew=Tmp;
      if (iter%100 == 0) ...
    }
  }
}
...

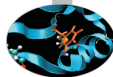
```



Laplace incremented

Fortran

```
...  
!$omp parallel  
  do while (var > tol .and. iter <= maxIter)  
    !$omp barrier  
    !$omp single  
      iter = iter + 1  
      var = 0.d0  
    !$omp end single  
    !$omp do reduction(max:var)  
      ...  
    !$omp end do  
    !$omp single  
      Tmp =>T; T =>Tnew; Tnew => Tmp;  
      if( mod(iter,100) == 0 ) ...  
    !$omp end single nowait  
  end do  
!$omp end parallel  
...
```

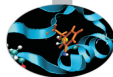


Still lacks OpenMP 3.1 support?

```

C
#pragma omp parallel
{
  while(var > tol && iter <= maxIter) {
    #pragma omp barrier
    #pragma omp single
    {
      ++iter;
      var = 0.0;
    }
    double pvar = 0.0;
    #pragma omp for nowait private(j)
    for (i=1; i<=n; ++i)
      for (j=1; j<=n; ++j) {
        Tnew[i*n2+j] = 0.25*( T[(i-1)*n2+j] + T[(i+1)*n2+j]
                             + T[i*n2+(j-1)] + T[i*n2+(j+1)] );
        pvar = fmax(pvar, fabs(Tnew[i*n2+j] - T[i*n2+j]));
      }
    #pragma omp critical
    if (pvar > var) var = pvar;

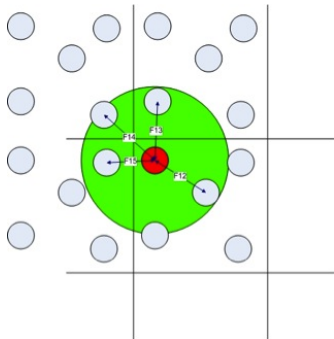
    #pragma omp barrier
    #pragma omp single nowait
    {
      Tmp=T; T=Tnew; Tnew=Tmp;
      if (iter%100 == 0) printf("iter: %8u, variation = %12.41E\n", iter, var);
    }
  }
}
  
```

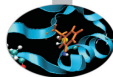


When the Going Gets Tough, ...

5 Parallelize the serial code **Nbody**. It computes the total energy and the forces of a system of N particles with potential $V = 1/r$ if r is less of a threshold and $V = 0$ otherwise.

- pay attention to the update of **forces**
 - try to update them atomically
 - try to reduce them
 - try different schedules and test their performance
-
- to compile use the preprocessing MACRO DIM=55000, for example
 - `gcc -O3 -DDIM=55000 Nbody.c -o nbody -lm`



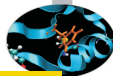


Nbody atomic

C

```

#pragma omp parallel for private(i,j,k,rij,d,d2,d3) reduction(+:ene) \
schedule(guided)
  for(i=0; i<nbodies; ++i)
    for(j=i+1; j<nbodies; ++j) {
      d2 = 0.0;
      for(k=0; k<3; ++k) {
        rij[k] = pos[i][k] - pos[j][k];
        d2 += rij[k]*rij[k];
      }
      if (d2 <= cut2) {
        d = sqrt(d2);
        d3 = d*d2;
        for(k=0; k<3; ++k) {
          double f = -rij[k]/d3;
          #pragma omp atomic
          forces[i][k] += f;
          #pragma omp atomic
          forces[j][k] -= f;
        }
        ene += -1.0/d;
      }
    }
  }
  
```

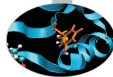


Nbody atomic

Fortran

```

!$omp parallel do private(i,j,k,rij,d,d2,f) reduction(+:ene) &
!$omp schedule(guided)
  do i = 1, DIM
    do j = i+1, DIM
      rij(:) = pos(:,i) - pos(:,j)
      d2 = 0.d0
      do k = 1, 3
        d2 = d2 + rij(k)**2
      end do
      if (d2 .le. cut2) then
        d = sqrt(d2)
        f(:) = - 1.d0 / d**3 * rij(:)
        do k=1, 3
          !$omp atomic
            forces(k,i) = forces(k,i) + f(k)
          !$omp atomic
            forces(k,j) = forces(k,j) - f(k)
        end do
        ene = ene + (-1.d0/d)
      end if
    end do
  end do
!$omp end parallel do
  
```



Nbody hand reduced in C

```

C ...
int tot_threads;
double ( *gforces ) [3];

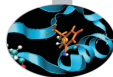
#pragma omp parallel private(i, j, k, rij, d, d2, d3)
{
#ifdef _OPENMP
  tot_threads = omp_get_num_threads();
#else
  tot_threads = 1;
#endif
  #pragma omp single
  gforces = calloc(nbodies*tot_threads, \
    sizeof( *gforces ));

  double ( *pforces ) [3];

#ifdef _OPENMP
  pforces = gforces + nbodies*omp_get_thread_num();
#else
  pforces = gforces;
#endif
}
  
```

```

C
#pragma omp for reduction(+:ene) schedule(guided)
for(i=0; i<nbodies; ++i)
  for(j=i+1; j<nbodies; ++j) {
    d2 = 0.0;
    for(k=0; k<3; ++k) {
      rij[k] = pos[i][k] - pos[j][k];
      d2 += rij[k]*rij[k];
    }
    if (d2 <= cut2) {
      d = sqrt(d2);
      d3 = d*d2;
      for(k=0; k<3; ++k) {
        double f = -rij[k]/d3;
        pforces[i][k] += f;
        pforces[j][k] -= f;
      }
      ene += -1.0/d;
    }
  }
#pragma omp for
for(i=0; i<nbodies; ++i)
  for(j=0; j<tot_threads; j++)
    for(k=0; k<3; ++k)
      forces[i][k] += gforces[i+j*nbodies][k];
}
  
```



Nbody reduction in Fortran

Fortran

```
!$omp parallel do private(i,j,k,rij,d,d2,f) &  
!$omp reduction(+:ene,forces) &  
!$omp schedule(guided)  
  do i = 1, DIM  
    do j = i+1, DIM  
      rij(:) = pos(:,i) - pos(:,j)  
      d2 = 0.d0  
      do k = 1, 3  
        d2 = d2 + rij(k)**2  
      end do  
      if (d2 .le. cut2) then  
        d = sqrt(d2)  
        f(:) = - 1.d0 / d**3 * rij(:)  
        forces(:,i) = forces(:,i) + f(:)  
        forces(:,j) = forces(:,j) - f(:)  
        ene = ene + (-1.d0/d)  
      end if  
    end do  
  end do  
!$omp end parallel do
```