OpenMP

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Introduction to OpenMP

OpenMP provides a way to write C++ code which uses multiple cores to run a program.

We control the number of threads using the global OMP_NUM_THREADS variable in a bash shell:

```
export OMP_NUM_THREADS=4
```

For example, the hello_openmp.cpp file contains:

```
#include <iostream>
int main(int argc, const char **argv)
{
    #pragma omp parallel
    {
        std::cout << "Hello OpenMP!\n";
    }
    return 0;
}</pre>
```

where the **#pragma omp parallel** line indicates the contents of the curly brackets below form a parallel section. Then the output depends on the value of the <code>OMP_NUM_THREADS</code> value.

We can now compile the code and set the thread variable as shown above, or specify it when running the code:

```
g++ -fopenmp hello_openmp.cpp -o hello_openmp

#setting parameter globally
export OMP_NUM_THREADS=4
./hello_openmp
```

```
## Hello OpenMP!
## Hello OpenMP!
## Hello OpenMP!
## Hello OpenMP!
```

```
OMP_NUM_THREADS=10 ./hello_openmp

## Hello OpenMP!
```

Directives

Hello OpenMP!
Hello OpenMP!

#setting parameter inline

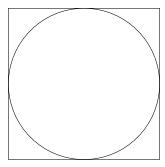
We use directives in OpenMP (e.g. the parallel directive above) to indicate how we want to use threads. Some basic examples include:

- sections: Used for code which can be run in parallel by different threads
- for: Does what sections does but for the contents of a loop
- parallel: Indicates a block to be executed by a team of threads.
- critical: Used when code can only be executed by one thread at a time.

We now use the information in this section to demonstrate how OpenMP can be useful in a concrete example.

Monte Carlo Example

A common algorithm for estimating π uses Monte Carlo methods and a unit circle (which has area π) inside a square with side length 2 (which has area 4):



Then if we generate points randomly inside the square, we expect the proportion of points inside the circle to be approximately $\frac{\pi}{4}$.

We can use a parallel loop to generate a large quantity of random numbers, determine which are in the circle and use the **reduce** directive to sum these counts.

We can generate random numbers using the <random> header. We can set the seed for the program by writing std::random_device before the parallel directive.

For each iteration we can then get a new generator and define a uniform distribution on the interval [-1,1] using these two lines of code:

```
std::default_random_engine generator(rd());
std::uniform_real_distribution random(-1.0, 1.0);
```

Then, if the count variables for points inside and outside the circle are called count_in and count_out respectively, the generating code under a for directive would be:

```
for (int i=0; i<1000000; i++){
   double x = random(generator);
   double y = random(generator);

if (r < 1.0){
    ++count_in;
}
else{
   ++count_out;
}</pre>
```

We can then use the reduction directive to combine the count_in and count_out variables from all the threads to estimate π .

Putting these parts together we obtain the solution:

```
#include <cmath>
#include <random>
#include <iostream>
int main()
{
   int n_inside = 0;
   int n_outside = 0;
   std::random_device rd;
    #pragma omp parallel reduction(+ : n_inside, n_outside)
        int count_in = 0;
        int count_out = 0;
        std::default_random_engine generator(rd());
        std::uniform_real_distribution<> random(-1.0, 1.0);
        #pragma omp for
        for (int i=0; i<1000000; ++i)
            double x = random(generator);
            double y = random(generator);
            double r = std::sqrt(x*x + y*y);
            if (r < 1.0)
            {
                ++count_in;
```

Now supposing the above is contained in a file called $\mathtt{pi.cpp}$ we can compile and run it in the terminal as follows:

```
g++ -fopenmp pi.cpp -o pi
./pi
```

pi is approximately 3.13989