

Notebook 1: #pragma omp parallel

The first step in OpenMP programming

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Introduction

We would like to optimize the computation of pi by numerical integration method.

A serial version of the algorithm is available on GitHub, alongside some useful material:



<https://github.com/gabrielefronze/iCSC/tree/master>

This version of the code is the sequential baseline to be used to measure speedup of parallel versions.

```
#include "StopWatch.h"
#include <iostream>

const long num_steps = 500000000; //number of x bins

int main(){
    Stopwatch stopWatch; //this object will be destroyed when out of
    scope

    double x, pi, sum = 0.0;
    step = 1.0/(double) num_steps; //x-step
    int n_threads=1;

    for (long i=1; i<=num_steps; i++) {
        x = (i - 0.5) * step; //computing the x value
        sum += 4.0 / (1.0 + x * x); //adding to the cumulus
    }

    pi = step * sum;

    printf("Pi value: %f\n
    Number of steps: %d\n
    Number of threads: %d\n",
    pi, num_steps, n_threads);
    return 0;
}
```

StopWatch

This class is useful for timing purposes.
It is an interface around:

omp_get_wtime()

This function returns a time measure which is safe
in a multithread environment.

The class is implemented as a default constructor
and a destructor: it is enough to create an object
inside the main function and to let it “die” outside
of the scope. The destructor prints on screen the
measured elapsed time.

```
#include <omp.h>
#include <iostream>

class Stopwatch{
public:
    Stopwatch(){
        start = omp_get_wtime();
    }
    ~StopWatch(){
        printf("Elapsed time: %f\n", omp_get_wtime()-start);
    }
private:
    double start;
}
```

Compilation

Compiling OpenMP code requires a compatible compiler, such as gcc, clang-omp or icc.

Each compiler has his own flag to enable OpenMP support. Typically the flag is `-fopenmp`.
For icc the flag is `-openmp`.

To compile a C/C++ file use the following command:

```
<compiler> -o <executable_name> -fopenmp <inputfile>
```

Ingredients

We want to use the pragma:

#pragma omp parallel

This pragma does a fork right after, running an exact copy of the following scope in each spawned thread.

The number of threads is defined and controlled by the environment variable:

OMP_NUM_THREADS

Accessible via the functions:

omp_set_num_threads()

omp_get_num_threads()

omp_get_thread_num()

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int main()
{
    Stopwatch stopWatch;

    double x, pi, sum = 0.0;
    step = 1.0/(double) num_steps; //x-step
    int n_threads=1;

    #pragma omp parallel
    {
        n_threads = omp_get_num_threads();

        // Some changes have to be made:
        // at the moment each thread performs the same operations
        // TIP: work on the for loop ranges
        for (long i=1; i<=num_steps; i++) {
            x = (i - 0.5) * step; //computing the x value
            sum += 4.0 / (1.0 + x * x); //adding to the cumulus
        }

        pi = step * sum;

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        }

        pi = step * sum;

        printf("Pi value: %f\n", pi);
        printf("Number of steps: %d\n", num_steps);
        printf("Number of threads: %d\n", n_threads);
        return 0;
    }
}
```

And now... Make the code *rain*!

Try to complete the exercise using `#pragma omp parallel`

