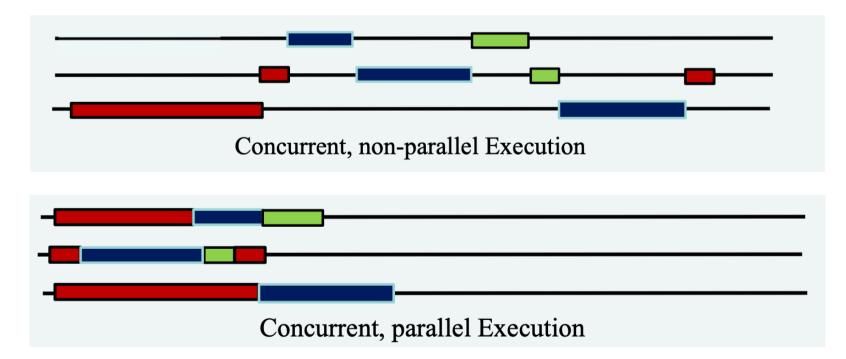
- Some useful resources:
 - Tim Mattson OpenMP notes (https://www.openmp.org/wp-content/uploads/Intro_To_OpenMP_Mattson.pdf) and videos (https://www.youtube.com/playlist?list=PLLX-Q6B8xqZ8n8bwjGdzBJ25X2utwnoEG)
 - ARCHER2 Youtube courses (https://www.archer2.ac.uk/training/courses/201006-openmp/)
 - Many other online lecture courses

- Around 20 years ago, the clock rate for CPUs stalled we can no longer speed up code simply by running on newer chips
- We must now make use of multiple chips to parallelise our code
- This means to provide access for one program to run on many cores
- We first concentrate on *threads* independent instances of code within a process which *share the same memory*

- An important distinction:
 - Concurrency multiple tasks are logically active at the same time but not running at the same time
 - E.g. checking emails while watching Netflix
 - E.g. C++ <thread> library
 - Concurrency can reduce wasted clock cycles (reduce latency) for example, while the program, is fetching something from
 memory, a program can implement some other part of the
 program
 - However it is not always associated with speedup

Threads and OpenMP

- Parallelism multiple tasks are actually active at the same time
 - Implemented to speed up execution



https://www.openmp.org/wp-content/uploads/Intro_To_OpenMP_Mattson.pdf

Threads and OpenMP

- Another important concept in parallelism is scaling this is how efficiently the program scales to using more cores
- There are two types of scaling:
 - Strong scaling how does the solution time vary with core number for a fixed problem
 - Eg. When we want to speed up an existing serial program
 - Weak scaling how does the solution time vary with core number for a fixed problem size per core
 - Eg. If we want to run a larger problem in the same time
- You will use these terms frequently if you do any kind of HPC development

- OpenMP is an API that supports shared-memory parallelisation via multithreading, developed in ~1990s
- The computational workload is divided up between threads that can run at the same time
- Practically:
 - Directives start with #pragma omp
 - We usually need to add #include <omp.h>

Example to demonstrate speedup:

- Set up the Docker container and `connect to running container'
- Compile and run openmp.cpp as normal, then time the program with time
- Uncomment the line containing #pragma
- Compile again, but this time with the flag -fopenmp
- Run and time again also try with different numbers of threads, set using export OMP NUM THREADS=number
- (But note: are we getting the right answer..?)

Threads and OpenMP

Exercise: Write a multithreaded program that prints "hello world"

```
#include <stdio.h>
int main() {
  int ID = 0;
  printf("hello(%d)", ID);
  printf("world(%d)\n", ID);
}
```

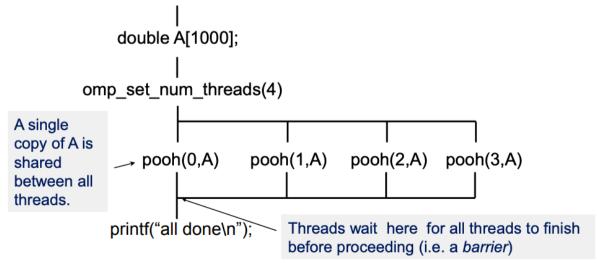
Exercise: Edit to run on multiple threads (remember the compiler flag -fopenmp)

```
#include <stdio.h>
int main() {
#pragma omp parallel
  int ID = 0;
  printf("hello(%d)", ID);
  printf("world(%d)\n", ID);
```

Exercise: Edit so we know which output comes from which thread

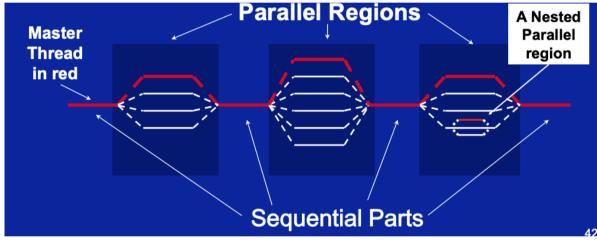
```
#include <stdio.h>
#include <omp.h>
int main() {
#pragma omp parallel
  int ID = omp_get_thread_num();
  printf("hello(%d)", ID);
  printf("world(%d)\n", ID);
```

- Here we have the desired output, but it is very muddled
- As we know, OpenMP threads share an address space
- This can lead to race conditions this is where the output of the program changes due to the threads running/finishing at different times



Threads and OpenMP

 We have seen how OpenMP creates a team of threads from the main thread using #pragma omp parallel



https://www.openmp.org/wp-content/uploads/Intro_To_OpenMP_Mattson.pdf

- If we want to set the number of threads in the code instead of using environment variables, we use omp set num threads (num) before the pragma
- OR we can edit the pragma to #pragma omp parallel num threads (num)

Threads and OpenMP

Exercise: Write a multithreaded program to perform a numerical integration of:

$$\int_0^1 \frac{4}{1+x^2} \mathrm{d}x$$

- We start with the serial version numerical_integration.cpp
- How do we want to split this up?
- Solution in numerical_integration_solution.cpp
 - Important: pay attention to the local and global variables
- Try timing both of these with more steps

Threads and OpenMP

- numerical_integration_solution.cpp is an example of a Single
 Program Multiple Data (SPMD) algorithm uses the thread ID to control which tasks to run
- If we time this example, we currently don't get good scaling
- This could be due to *false sharing* this occurs if independent data elements happen to sit on the *same cache line*
 - This is common in OpenMP when we promote a scalar to an array to index by the thread number
 - The array elements are contiguous in memory, so share cache lines poor scalability
- One way around this is to pad the arrays so that we are on distinct cache lines - note though that this is architecture dependent
- Another way is to make use of private variables and restructure so there is less sharing of data (see later...)

Threads and OpenMP

- As we saw briefly at the beginning, we can parallelise for loops using #pragma omp parallel for (or in Fortran, #pragma omp parallel do)
- This splits up the loop across threads in a team
- Significantly improves the readability of the code over indexing by thread number

```
#pragma omp parallel
                               int id, i, Nthrds, istart, iend;
OpenMP parallel
                               id = omp get thread num();
region
                               Nthrds = omp get num threads();
                               istart = id * N / Nthrds:
                               iend = (id+1) * N / Nthrds;
                               if (id == Nthrds-1)iend = N;
                               for(i=istart;i<iend;i++) { a[i] = a[i] + b[i];}
OpenMP parallel
                       #pragma omp parallel
region and a
                       #pragma omp for
worksharing for
                               for(i=0;i<N;i++) {a[i] = a[i] + b[i];}
construct
```

- Must make sure that the loop iterations are independent
- We can also use this for *nested loops*

```
#pragma omp parallel for collapse(2)
```

where here, the number of loops is 2

- If we do have a loop where we are gathering values into one variable, such as eg. sum += value[i], we can use a reduction
- In this case, a local copy of each list variable is made and updates, then local copies are reduced to a single value and combined

Example:

```
double sum=0., average, values[max];
int i;
for(i=0; i<max; i++)
{
   sum += values[max];
}
average = sum/max;</pre>
```

Example:

```
double sum=0., average, values[max];
int i;
#pragma omp parallel for reduction (+:sum)
for(i=0; i<max; i++)

{
   sum += values[max];
}
average = sum/max;</pre>
```

Exercise for outside class - parallelise the integration program with a loop construct