

## Drug Design

In the folder `~/CSinParallel/RaspberryPiBasics/drugdesign/` you will find three implementations of a program that attempts to locate small molecules called *ligands* that are good candidates for use as drugs. When a protein associated with a disease of interest is identified, its three-dimensional structure is found either experimentally or through a molecular modeling computation. A collection of ligands is tested against the protein. A score is set based on these binding properties, and the best scores are flagged, identifying ligands that would make good drug candidates. For simplicity, the programs we implement use strings to represent proteins and ligands. The three parallel implementations can be summarized as follows:

- `drugdesign-static`: Uses OpenMP to assign chunks of ligands at compile time to cores.
- `drugdesign-dynamic`: Uses OpenMP to assign ligands to cores dynamically (at run-time).
- `drugdesign-threads`: Uses C++11 threads (instead of OpenMP threads) to solve the problem.

To compile each version, `cd` into each individual directory and type `make`. Your goal is to determine which program is the fastest.

## Measure Run-Time

To measure the running time of each implementation, use `time -p`. For example:

```
> time -p ./drugdesign-static 1
```

Runs the `drugdesign-dynamic` executable on a single thread. Fill out the table below (use `real` for time):

Time (s)	1 Thread	2 Threads	3 Threads	4 Threads
<code>drugdesign-static</code>				
<code>drugdesign-dynamic</code>				
<code>drugdesign-threads</code>				

## Compute the Speedup

*Speedup* measures how much faster the parallel execution of a program is compared to the serial (single threaded) execution. The Speedup of a  $n$  core approach ( $S_n$ ) is computed by the following equation:

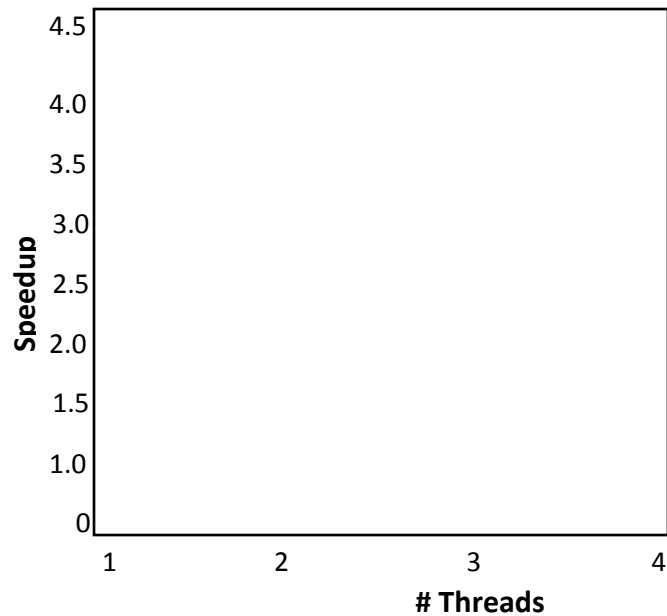
$$S_n = \frac{\text{Time}(1 \text{ thread})}{\text{Time}(n \text{ threads})}$$

Hint: Consider using the calculator application on the Raspberry Pi 2 to help you with the calculation!

Speedup	2 Threads	3 Threads	4 Threads
<code>drugdesign-static</code>			
<code>drugdesign-dynamic</code>			
<code>drugdesign-threads</code>			

## Plot Your Results

*Problem 2c:* Plot your speedup results in the area below. Use different lines for each of the different applications.



## Discussion Questions

1. Which approach is the fastest?
2. Determine the number of lines in each file (use `wc -l`). How does the C++11 implementation compare to the OpenMP implementations?
3. Increase the number of threads to 5 threads. What is the run time for each?
4. Increase the maximum ligand length to 7, and rerun each program. Are your speedup results similar?