

GPU BOOTCAMP MINI CHALLENGE

OpenACC
More Science, Less Programming



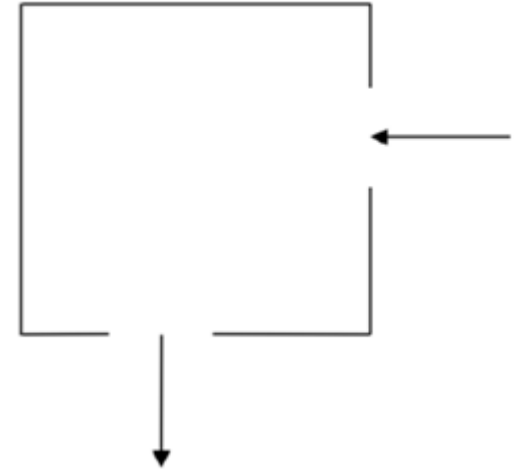
APPLICATION

Simple 2D regular-grid CFD simulation

```
set the boundary values for  $\Psi$ 
while (convergence == FALSE) do
  for each interior grid point do
    update  $\Psi$  by averaging with its 4 nearest neighbours
  end do

  check for convergence
end do

for each interior grid point do
  calculate  $u_x$  calculate  $u_y$ 
end do
```



The objective of this exercise is not to dwell into the maths but to make use of different approaches to GPU programming to parallelize and improve the performance.

Pseudo Code

```
int main(int argc, char **argv) {
```

```
    initialization loop
```

```
    boundary calculation loop
```

```
    Jacobi loop
```

```
    swap array loop
```

```
}
```

- [cfd.cpp](#)

- [boundary.cpp](#)

- [jacobi.cpp](#)

MORE ABOUT CODE

- Uses Makefile
- To run the code `./cfd 64 500`
 - Where `./cfd` is application name
 - 64 is size of scaling
 - 500 is number of max iteration

Output:

... finished

After 500 iterations, the error is **0.00211211** -> Check this value to confirm your porting

Time for 500 iterations was 18.8579 seconds

Each iteration took 0.0377159 seconds

HINTS

- Divide different methods to port among team members
- Use profiler to check the hotspots and bottlenecks in your code
- Make use of compiler flag to cross check if indeed parallelization was done e.g. `-Minfo`
- Key files to look out having maximum loops:
 - `cfid.cpp`
 - `jacobi.cpp`
- Download and take backup

Acknowledgment

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