

## Homework 4

Total number of points: 100 (+ 15 bonus).

This assignment builds on the previous assignment's theme of examining memory access patterns. You will implement a finite difference solver for the 2D heat diffusion equation in different ways to examine the performance characteristics of different implementations. Please refer to the class and lecture slides on this homework.

**Background on the heat diffusion PDE.** The heat diffusion PDE that we will be solving can be written

$$\frac{\partial T}{\partial t} = \frac{\partial^2 T}{\partial x^2} + \frac{\partial^2 T}{\partial y^2}$$

To solve this PDE, we are going to discretize both the temporal and the spatial derivatives. To do this, we define a two-dimensional<sup>1</sup> grid  $G_{i,j}$ ,  $1 \leq i \leq n_x$ ,  $1 \leq j \leq n_y$ , where we denote by  $n_x$  (resp.  $n_y$ ) the number of points on the  $x$ -axis (resp.  $y$ -axis). At each time step, we will evaluate the temperature and its derivatives at these grid points.

While we will consistently use a first-order discretization scheme for the temporal derivative, we will use 2<sup>nd</sup>, 4<sup>th</sup> or 8<sup>th</sup> order discretization of the spatial derivative.<sup>2</sup>

If we denote by  $T_{i,j}^t$  the temperature at time  $t$  at point  $(i, j)$  of the grid, the 2<sup>nd</sup> order spatial discretization scheme can be written as:

$$T_{i,j}^{t+1} = T_{i,j}^t + C^{(x)} (T_{i+1,j}^t - 2T_{i,j}^t + T_{i-1,j}^t) + C^{(y)} (T_{i,j+1}^t - 2T_{i,j}^t + T_{i,j-1}^t)$$

The  $C^{(x)}$  (x`fc1` in the code) and  $C^{(y)}$  (y`fc1` in the code) constants are called Courant numbers. They depend on the temporal discretization step, as well as the spatial discretization step. To ensure the stability of the discretization scheme, they have to be less than the maximum value given by the Courant-Friedrichs-Lewy condition.<sup>3</sup> You do not have to worry about this, because the starter code already takes care of picking the temporal discretization step to maximize the Courant numbers while ensuring stability.

The starter code also contains host and device functions named `stencil` which contain the coefficients that go into the update equation. Therefore, you do not need to figure out how to implement the different order updates. You only need to understand how this function works and pass in the arguments correctly.

**Boundary conditions.** The starter code contains the functions that will update the boundary conditions for you (see file `BC.h`, in particular, the function `gpuUpdateBCsOnly`) and the points that are in the border (which has a size of order  $\sqrt{2}$ ). This way, you do not have to worry about the size of the stencil as you approach the wall.<sup>4</sup>

<sup>1</sup>In fact, the size of the grid is  $g_x \times g_y$  but we will only update the interior region, which size is  $n_x \times n_y$ .

<sup>2</sup>For instance, if we use a 4<sup>th</sup> order scheme, we will express the derivative with respect to  $x$  of  $T$  at point  $(i, j)$  using  $T_{i-2,j}^t$ ,  $T_{i-1,j}^t$ ,  $T_{i,j}^t$ ,  $T_{i+1,j}^t$  and  $T_{i+2,j}^t$ .

<sup>3</sup>If you are interested, you can read more about this at [http://en.wikipedia.org/wiki/Courant-Friedrichs-Lewy\\_condition](http://en.wikipedia.org/wiki/Courant-Friedrichs-Lewy_condition).

<sup>4</sup>A general way of dealing with this problem is to change the size of the stencil as you approach the wall. This is complicated and we simplified the process for this homework.

**Various implementations.** In this programming assignment, we are going to implement 2 different kernels (and you can do a third one for extra credit):

- Global (function `gpuStencilGlobal`): this kernel will use global memory and each thread will update exactly one point of the mesh. You should use a 1D Grid and 1D Blocks with  $n_x \times n_y$  threads total.
- Block (function `gpuStencilBlock`): this kernel will also use global memory. Each thread will update `numYPerStep` points of the mesh (these points form a vertical line). You should use a 2D grid with  $n_x \times n_y / \text{numYPerStep}$  threads total.
- (Extra Credit) Shared (function `gpuStencilShared`): this kernel will use shared memory. A group of threads must load a piece of the mesh in shared memory and then compute the new values of the temperatures on the mesh.<sup>5</sup> Each thread will load and update several elements.

**Parameter file.** The parameters used in the computation are read from the file `params.in`. You will need to modify *some parameters* (see description of the starter code) in this file to answer some of the questions. But this file will not be submitted through the submission script.

Here is a list of parameters that are used in the order they appear in the file:

```
int nx_, ny_;      // number of grid points in each dimension
double lx_, ly_;   // extent of physical domain in each dimension
int iters_;        // number of iterations to do
int order_;        // order of discretization
```

**Starter code** The starter code is composed of the following files (\* means the file will *not* be submitted by our script):

- `*main.cu` — This is the CUDA driver file. Do not modify this file.
- `gpuStencil.cu` — This is the file containing the kernels. You will need to modify this file.
- `Makefile` — `make` will build the binary. `make clean` will remove build files as well as debug output. Do not modify this file unless you are changing the `GTESTROOT` variable or are changing the `fmad` flag. More about this flag will be discussed below
- `*params.in` — This file contains a basic set of parameters. For debugging, performance testing, and to answer the questions, you will need to modify this file. The only parameters you should modify are `nx`, `ny` (line 1), and `order` (line 4). This file however will not get submitted through the script.
- `*simParams.h` and `*simParams.cpp` — These files contain the data structure necessary to handle the parameters of the problem. Do not modify these files.
- `*Grid.h` and `*Grid.cu` — These files contain the data structure that models the grid used to solve the PDE. Do not modify this file.
- `*BC.h` — This file contains the class `boundary_conditions` that will allow us to update the boundary conditions during the simulation. Do not modify this file.
- `*CPUComputation.h` — This file contains the functions related to the CPU implementation of the solver. Do not modify this file

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<sup>5</sup>Note, however, that the threads that loaded data on the borders of the small piece will stay idle during the computation step.

- `*Errors.h` — This file contains the functions we will use to test your code, i.e., to check the errors of the output and print out any errors to different file streams.
- `*hw4.sh` — This script is used to submit jobs to the queue (see Appendix A).

**Note** The files in the starter code contain some additional information about the implementation in the form of comments. Additionally, the CPU implementation should provide a clearer picture of the method and should aid your GPU work.

**Running the program** Type `make` to compile the code. Once this is done, you can use the command:  
`$ srun --partition=CME --gres=gpu:1 ./main` or just run `$ sbatch hw4.sh` which has all of the different param combinations we will ask about

When you run this program, the `main` function will be called and will initialize some global variables so that we may reuse the starting grid for all three of our tests `global`, `block`, and `shared`. Once the initialization is done in `main`, then the script will execute the tests in order. Note that if you aren't doing the extra credit (i.e. implementing the shared kernel), then your code will always fail the last test. This is expected. The output produced by the program will contain:

- The time and bandwidth for the CPU implementation, as well as the GPU implementations running the program.
- A report of the errors for the GPU implementations. Namely, the program will output:
  - The  $L_2$  norm of the final solution from the CPU implementation (i.e., the reference).
  - The  $L_\infty$  norm of the relative error between the reference solution and the GPU solution.
  - The  $L_2$  norm of the error normalized by the  $L_2$  norm of the CPU implementation.

The output of `hw4.sh` script can be found in the `result.txt` file that will be created when you run the script. With `--fmad=false`, the errors should be equal to 0. We will test your code with this flag, so make sure that your code passes all of our tests with this flag set. You may set the flag to true when calculating the bandwidth scores.

Without this option, typical ranges for roundoff errors are:

- $[10^{-8}, 10^{-7}]$  for  $L_\infty$  norm error.
- $[10^{-9}, 10^{-8}]$  for  $L_2$  norm error.

Search for `TODD` in `gpuStencil.cu` to see where you need to implement code.

### Question 1

(30 points) Implement the function `gpuStencilGlobal` that runs the solver using global memory, and creates 3D surface plots of temperature on a  $256 \times 256$  grid at iteration 0, 1000, and 2000 respectively, with 8th order. You must also fill in the function `gpuComputationGlobal`. The difference (in terms of the norms) between your implementation and the reference should be in the expected range. Grid class implements member function `saveStateToFile` to dump all the data of the grid to a CSV file. You can choose your own tools (Python, MATLAB, etc.) to generate those plots and include them in your writeup.

## Question 2

(35 points) Implement the function `gpuStencilBlock` that runs the solver using global memory but where each thread computes `numYPerStep` points on the grid. You must also fill in the function `gpuComputationBlock`. You should use 2D blocks and grids to implement the blocking strategy we talked about in class. You need to decide your block and grid dimensions, as well as `numYPerStep` to optimize the performance of your code. The difference (in terms of the norms) between your implementation and the reference should be in the expected range.

## Question 3

(15 points) Plot the bandwidth (GB/s) as a function of the grid size (in units of MegaPoints) for the following grid sizes:  $256 \times 256$ ;  $512 \times 512$ ;  $1024 \times 1024$ ;  $2048 \times 2048$ ;  $4096 \times 4096$ . You can choose the number of iterations to run to generate your results. You should run enough iterations so results are stable and meaningful (clearly 1 is not enough).

You must have 2 plots (or 3 plots if you choose to do the extra credit) as follows:

1. For order = 4, plot the bandwidth for the 2 (or 3) different algorithms.
2. For the block implementation, plot the bandwidth for the 3 different orders.
3. If you implemented the shared algorithm, plot the bandwidth for the 3 different orders.

## Question 4

(20 points) Which kernel (global, block, or shared) and which order gives the best performance (your answer may depend on the grid size)? Explain the performance results you got in Question 3. Specifically, explain performance differences (i) among kernels, (ii) from varying order, and (iii) from varying problem sizes.

## Question 5

(Extra credit 15 points) Implement the function `gpuStencilShared` that runs the solver using shared memory. You should also fill in the function `gpuComputationShared`. Note that you have to answer the questions related to shared memory implementation in Questions 3 and 4 to get the full extra credit.

# A Submission instructions

To submit:

1. For all questions that require explanations and answers besides source code, put those explanations and answers in a separate PDF file and upload this file on Gradescope.
2. Make sure your code compiles on `icme-gpu` and runs. To check your code, we will run:  

```
$ make
```

  
This should produce one executable: `main`
3. The homework should be submitted using a submission script on `cardinal`. The submission script must be run on `cardinal.stanford.edu`.
4. Copy your submission files to `cardinal.stanford.edu`. The script `submit.py` will copy only the files below to a directory accessible to the CME 213 staff. Only these files will be copied. Any other required files (e.g., `Makefile`) will be copied by us. Therefore, make sure you make changes only to the files below. You are free to change other files for your own debugging purposes, but make sure

you test it with the default test files before submitting. Also, do not use external libraries, additional header files, etc, that would prevent the teaching staff from compiling the code successfully. Here is the list of files we are expecting and that will be copied:

`gpuStencil.cu`

5. To submit, type:

```
$ /afs/ir.stanford.edu/class/cme213/script/submit.py hw4 <directory with your submission files>
```

6. You can submit at most 10 times before the deadline; each submission will replace the previous one.

## Machine

We will be using the `icme-gpu` cluster with NVIDIA Quadro RTX 6000 GPU. We have provided a script called `hw4.sh` in the starter code to help you submit a batch job with proper compute node configurations. You should run `module load cuda` once you log into `icme-gpu`, and you can submit a batch job using `sbatch hw4.sh`, which compiles your CUDA code, runs the main executable, and writes the output to `slurm.sh.out`. You can modify `hw4.sh` to run with different arguments.

## B Advice and hints

- Make sure you understand what are the parameters of the problem. In particular, the difference between `nx` and `gx` is important to understand.
- Spend some time looking at the `simParams` class as it contains the useful parameters to solve the problem.
- Keep in mind where your data are when trying to decide kernel parameters. You will be reading from global memory using Read-Only cache (L1, 128-byte cache lines). Writes to global memory can not be cached in L1, therefore have to go to the shared L2 cache, which has 32-byte cache lines.
- Make sure your implementations can deal with square grids as well as rectangular ones.