

Homework 4—due Friday May 10, 11 PM

Total number of points: 100 (+ 15 bonus). Late day policy: 2 late days with a 10% grade penalty.

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.

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 grid $G_{i,j}$, $1 \le i \le n_x$, $1 \le j \le n_y$, 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 gridpoints.

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

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

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

The $C^{(x)}$ (xfcl in the code) and $C^{(y)}$ (yfcl 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 a maximum value given by the Courant-Friedrichs-Lewy condition.³ You do not have to worry about this, because the starter code already takes care of picking the temporal discretization step as 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 / 2). This way, you do not have to worry about the size of the stencil as you approach the wall.⁴

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

²For instance, if we use a 4th order scheme, we will express the derivative with respect to x of T at point (i, j)

read more about this at http://en.wikipedia.org/wiki/ Courant-Friedrichs-Lewy_condition.

⁴A 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 gpuStencil): 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 gpuStencilLoop): 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 $n_x \times n_y$ /numYPerStep threads total.
- (Extra Credit) Shared (function gpuShared): 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.⁵ 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 be 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.
- *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.
- *create_vm.sh—This script is used to initialize a Google Cloud Platform VM with a NVIDIA K80 GPU, with name hw4. See the course website for further instructions.

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

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:

\$./main [-g] [-b] [-s]

where -g stands for global, -b for block, and -s for shared. Note that you can run several implementations at the same time (for instance ./main -g -b to run the global and block implementations). The output produced by the program will contain:

- The time and bandwidth for the CPU implementation, as well as for the implementations that you specified when 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_{∞} 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.

Typical ranges for the errors (and for the parameters that you will be using) are:

- $[10^{-5}, 10^{-3}]$ for L_{∞} norm error.
- $[10^{-6}, 10^{-4}]$ for L₂ norm error.

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

Question 1

(30 points) Implement the function gpuStencil that runs the solver using global memory, and create 3D surface plots of temperature on a 256 × 256 grid at iteration 0, 1000 and 2000 respectively, with 8th order. You must also fill in the function gpuComputation. 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 gpuStencilLoop that runs the solver using global memory but where each thread computes numYPerStep points on the grid. You must also fill in the function gpuComputationLoop. You should use 2D blocks and grid to implement the blocking strategy we talked about in the class. You need to decide your block and grid dimensions, as well as numYPerStep to optimize 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: 256×256 ; 512×512 ; 1024×1024 ; 2048×2048 ; 4096×4096 . You can choose number of iterations to run in order 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.

Question 5

(Extra credit 15 points) Implement the function gpuShared 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 in order 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 Google Cloud VM 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:
- \$ /usr/bin/python /usr/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 Google Cloud VM. We provide a script <code>create_vm.sh</code> in your starter code to set up the correct machine for this homework. You can also reuse your hw3 machine, which has the same configurations. You have all the necessary libraries and tools pre-installed on the VM.

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 shared L2 cache, which has 32-byte cache lines).
- Make sure your implementations are able to deal with square grids as well as rectangular ones.