## ME 759

## High Performance Computing for Engineering Applications Assignment 3 Due Thursday 09/29/2022 at 9:00 PM

Submit responses to all tasks which don't specify a file name to Canvas in a file called assignment3. {txt, docx, pdf, rtf, odt} (choose one of these formats). Submit all plots (if any) on Canvas. Do not zip your Canvas submission.

All source files should be submitted in the HW03 subdirectory on the main branch of your GitLab repo. Please use the name HW03 exactly as shown here (both in terms of capitalization & name). The HW03 subdirectory should have no subdirectories. For this assignment, your HW03 folder should contain task1.cu, task2.cu, task3.cu, and vscale.cu.

All commands or code must work on *Euler* with only the nvidia/cuda module loaded. The commands may behave differently on your computer, so be sure to test on *Euler* before you submit.

Please submit clean code. Consider using a formatter like clang-format.

IMPORTANT: Before you begin, copy any provided files from Assignments/HW03 directory of the ME759 Resource Repo. Do not change any of the provided files since these files will be overwritten with clean, reference copies when grading.

- 1. Write a C++ program using CUDA in a file called task1.cu which computes the factorial of integers from 1 to 8, by launching a GPU kernel with 1 block and 8 threads. Inside the kernel, each thread should use std::printf to write out a!=b (followed by a newline), where a is one of the 8 integers, and b is the result of a!. (Follow your kernel call with a call to cudaDeviceSynchronize() so that the host waits for the kernel to finish printing before returning from main.)
  - Compile: nvcc task1.cu -Xcompiler -03 -Xcompiler -Wall -Xptxas -03 -std=c++17 -o task1
  - Run (on Euler, use Slurm sbatch!): ./task1
  - Expected output (showing only 4 out of the 8 lines expected; lines could be out of order):
    - 1!=1
    - 2!=2
    - 3!=6
    - 4!=24
  - It is ok to publish your sbatch script on Piazza. For this assignment, you will need to ask Slurm to execute your program on a node that has GPU cards

- 2. Write a C++ program using CUDA in a file called task2.cu which does the following:
  - From the host, allocates an array of 16 ints on the device called dA.
  - Launches a kernel with 2 blocks, each block having 8 threads.
  - Each thread computes ax+y and writes the result in one distinct entry of the dA array. Here,
    - x is the thread's threadIdx;
    - y is the thread's blockIdx;
    - a is an integer argument that the kernel takes (so all threads use the same a). You need to generate a randomly and then call the kernel with it. It is up to you how you generate this random number, one possible approach is described here BestPractice.
  - Copies back the data stored in the device array dA into a host array called hA.
  - Prints (from the host) the 16 values stored in the host array separated by a single space each.

How to go about it, and what the expected output looks like:

- Compile: nvcc task2.cu -Xcompiler -03 -Xcompiler -Wall -Xptxas -03 -std=c++17 -o task2
- Run (on Euler, use Slurm sbatch!): ./task2
- Expected output (followed by newline; yours could be different depending on the random number generation): 0 10 20 30 40 50 60 70 1 11 21 31 41 51 61 71

3. a) Implement in a file called vscale.cu, the vscale kernel function as declared and described in vscale.cuh. This function should take in two arrays, a and b, and do an element-wise multiplication of the two arrays:  $b_i = a_i \cdot b_i$ . In the process, b gets overwritten. Each thread should do at most one of the multiplication operations.

## Example:

$$\mathbf{a} = [-5.0, 2.0, 1.5], \quad \mathbf{b} = [0.8, 0.3, 0.6], \quad \mathbf{n} = 3$$

The resulting **b** array is:

$$\mathbf{b} = [-4.0, 0.6, 0.9].$$

- b) Write a file task3.cu which does the following:
  - Creates two arrays of length n filled by random numbers where n is read from the first command line argument. The range of values for array a is [-10.0, 10.0], whereas the range of values for array b is [0.0, 1.0].
  - Calls your vscale kernel with a 1D execution configuration that uses 512 threads per block.
  - Prints the amount of time taken to execute the kernel in *milliseconds* using CUDA events<sup>2</sup>.
  - Prints the first element of the resulting array.
  - Prints the last element of the resulting array.

How to go about it, and what the expected output looks like:

- Compile: nvcc task3.cu vscale.cu -Xcompiler -03 -Xcompiler -Wall -Xptxas -03 -std=c++17 -o task3
- Run (where n is a positive integer): ./task3 n
- Example expected output (followed by newline):
  - 0.012
  - 1.3
  - 2.3
- c) On an Euler *compute node*, run task3 for each value  $n = 2^{10}, 2^{11}, \dots, 2^{29}$  and generate a plot task3.pdf which plots the time taken by your vscale as a function of n. Overlay another plot which shows the scaling results when using 16 threads per block.

<sup>&</sup>lt;sup>1</sup>Details about random number generation can be found in random\_numbers.md.

<sup>&</sup>lt;sup>2</sup>Recall the GPU timing section of the document timing.md.