

ECE 759
High Performance Computing for Engineering Applications
Assignment 5
Due Friday 11/08/2024 at 23:59 PM

Submit responses to all tasks which don't specify a file name to Canvas in a file called `assignment5.pdf`. Submit all plots (if any) on Canvas. Do not zip your Canvas submission.

All *source files* should be submitted in the `HW05` subdirectory on the `main` branch of your GitLab repo. Please use the name `HW05` exactly as shown here (both in terms of capitalization & name). The `HW05` subdirectory should have no subdirectories. For this assignment, your `HW05` 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.

- `#SBATCH --gpus-per-task=1` should be added, which requests one node with 1 GPU.

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

IMPORTANT: Before you begin, copy any provided files from `Assignments/HW05` directory of the [ECE 759 Resource Repo](#). Do not change any of the provided files since these files will be overwritten with clean, reference copies when grading.

The GitHub link to your code folder should be: <https://github.com/YourGitHubName/repo759/HW05>

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 -O3 -Xcompiler -Wall -Xptxas -O3 -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 -O3 -Xcompiler -Wall -Xptxas -O3 -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:

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

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 numbers¹ 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 events².
 - 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 -O3 -Xcompiler -Wall -Xptxas -O3 -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 = 210, 211, ..., 229` 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.

¹Details about random number generation can be found in [random_numbers.md](#).

²Recall the GPU timing section of the document [timing.md](#).