Heterogeneous computing with performance modelling

GPU programming basics

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Why should I use GPUs?

Why should I use GPUs?









Why should I use GPUs?

- Nvidia says that when compared to CPUs, GPUs provides much higher
 - instruction throughput and
 - memory bandwidth.
- In particular, this performance is delivered in a relatively small price and power envelope.

What does this mean?









Why should I use GPUs? (flop rate, double precision)

Quad-core Intel Skylake CPU:

$$\sim$$
 200 GFlops

▶ 14-core Intel Xeon Gold 6132 CPU:

$$\sim$$
 1200 GFlops

Nvidia Tesla V100 GPU:

 \sim **7000** GFlops







Why should I use GPUs? (memory bandwidth)

Quad-core Intel Skylake CPU:

$$\sim 35~\text{GB/s}$$

▶ 14-core Intel Xeon Gold 6132 CPU:

$$\sim$$
 100 GB/s

Nvidia Tesla V100 GPU:

$$\sim$$
 900 GB/s







Why should I use GPUs? (power usage)

Quad-core Intel Skylake CPU:



▶ 14-core Intel Xeon Gold 6132 CPU:

$$\sim$$
 140 W

Nvidia Tesla V100 GPU:

 \sim 250 W







Why should I use GPUs? (price)

Quad-core Intel Skylake CPU:



▶ 14-core Intel Xeon Gold 6132 CPU:

$$\sim$$
 \$2 000

Nvidia Tesla V100 GPU:

 \sim \$10 000









Why should I use GPUs? (per Watt)

Quad-core Intel Skylake CPU:

 \sim 2 GFlops per Watt, \sim 0.4 GB/s per Watt

▶ 14-core Intel Xeon Gold 6132 CPU:

 \sim 9 GFlops per Watt, \sim 0.7 GB/s per Watt

Nvidia Tesla V100 GPU:

 \sim **28** GFlops per Watt, \sim **3.6** GB/s per Watt









Why should I use GPUs? (per dollar)

Quad-core Intel Skylake CPU:

$$\sim$$
 0.5 GFlops per \$, \sim 0.09 GB/s per \$

▶ 14-core Intel Xeon Gold 6132 CPU:

$$\sim$$
 0.6 GFlops per \$, \sim 0.05 GB/s per \$

Nvidia Tesla V100 GPU:

 \sim **0.7** GFlops per \$, \sim **0.09** GB/s per \$







Why should I use GPUs? (reduced precision)

Quad-core Intel Skylake CPU:

$$\sim$$
 400 GFlops (single precision)

▶ 14-core Intel Xeon Gold 6132 CPU:

$$\sim$$
 2400 GFlops (single precision)

Nvidia Tesla V100 GPU:

 \sim 14 000 GFlops (single precision) \sim **112 000** GFlops (half precision, tensor cores)









What is the catch?

What is the catch?

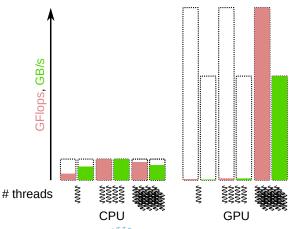






What is the catch?

- GPUs are highly parallel. Obtaining peak performance requires thousands of threads.
- ▶ The algorithms and the software must be highly parallel.











What is the catch?

- GPU programming is difficult.
 - As mentioned, requires a lot of parallelism.
 - Usually, the GPU-side code is written in "C++".
- The CUDA cores are not "true" cores.
 - The cores share resources such as schedulers and caches.
 - The tensor cores (half precision) are even more limited.
- Limited amount of memory.
 - Data must be moved between memories (RAM, VRAM, ...).
 - Certain memories are significantly faster than others.
 - Certain memories are fast only when used correctly.
- Not all algorithms are suitable for GPUs.
 - Some algorithms are just slightly faster on GPUs.
 - Does it make sense to pay the extra money?
- Not all GPUs are equally powerful.
 - Some GPUs have a very low double-precision flop rate.
 - Some GPUs do not have tensor cores.









CUDA basics

Lets go through some CUDA basics...









What is CUDA?

- Compute Unified Device Architecture (CUDA) is a parallel computing platform and an API created by Nvidia.
 - Only for Nvidia GPUs.
- Can be used
 - directly through CUDA C/C++ and CUDA Fortran,
 - directly through wrappers (Python, Perl, Fortran, Java, Ruby, Lua, etc),
 - indirectly through compiler directives (OpenACC), and
 - indirectly through other computational interfaces (OpenCL, DirectCompute, OpenGL, etc).
- During this training course, we will use CUDA C/C++.









What is CUDA?

- CUDA comes with its own compiler, nvcc.
 - GPU-specific code is compiled to PTX (Parallel Thread Execution) "assembly".
 - The PTX code is translated to binary code by the driver.
 - ▶ The rest is offloaded to the host compiler (e.g., g++).
- CUDA also comes with a large set of libraries:
 - CUDA Runtime library (CUDART), basic functionality.
 - CUDA Basic Linear Algebra Subroutines (cuBLAS).
 - CUDA Fast Fourier Transform library (cuFFT).
 - CUDA Sparse Matrix library(cuSPARSE).
 - CUDA Deep Neural Network library (cuDNN).
- Many 3rd party libraries are use these vendor libraries.
- In most cases, you want to use these libraries instead of implementing your own.









Hello world

A "Hello world" program (hello.cu) is a good place to start:

```
#include <stdlib.h>
#include <stdlib.h>
__global__ void say_hello()
{
    printf("GPU says, Hello world!\n");
}
int main()
{
    printf("Host says, Hello world!\n");
    say_hello <<<1,1>>>();
    cudaDeviceSynchronize();
    return EXIT_SUCCESS;
}
```







Hello world (compile and run)

- Load the correct toolchain:
 - \$ ml purge
 - \$ ml fosscuda/2019b buildenv
- Compile the source code with nvcc:
 - \$ nvcc -o hello hello.cu
- Queue a job:

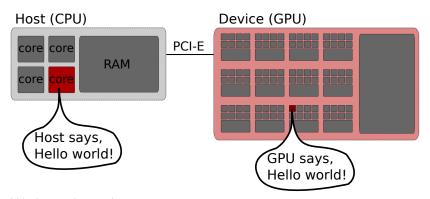
```
srun --account=SNIC2020-9-161 \
--reservation=snic2020-9-161-day1 \
--ntasks=1 --gres=gpu:v100:1,gpuexcl \
--time=00:02:00 ./hello
Host says, Hello world!
GPU says, Hello world!
```







Hello world (what is happening)



We have three objects:

Host CPUs + main memory (RAM)

Device A GPU consisting of several CUDA cores

PCI-E Fast interconnect between the host and the device









Hello world (kernels)

- ▶ The device-side code is written inside special functions called kernels.
- A kernel is declared with the __global__ keyword:

```
__global__ void say_hello()
{
    printf("GPU says, Hello world!\n");
}
```

- Kernel's return type is always void.
- All threads enter the kernel from the beginning of the body of the function.
 - Single Instruction, Multiple Thread (SIMT)
 - Threads are not spawned fork-join style (except when a kernel launches other kernel).









Hello world (kernel launch)

▶ The host **issues** the say_hello kernel as follows:

```
|| say_hello <<<1,1>>>();
```

- This places the kernel call into a queue known as stream.
 - Note that the kernel is not guaranteed to be executed!
- ▶ We will return to the <<< . , . >>> brackets later...
 - For now, you need to know that the number of threads is one.
- The cudaDeviceSynchronize call causes the host to wait until stream is empty, i.e., the kernel has finished:

```
|| cudaDeviceSynchronize();
```









Hello world (summary)

```
#include <stdlib.h>
#include <stdio.h>
__global__ void say_hello()
    printf("GPU says, Hello world!\n");
int main()
    // the host (CPU) executes these lines
    printf("Host says, Hello world!\n");
    // issue the say_hello kernel
    say_hello <<<1,1>>>();
    // wait until the kernel has finished
    cudaDeviceSynchronize();
    return EXIT SUCCESS:
```









AX example (scalar-vector multiplication)

Lets try something more complicated:

$$\alpha \in \mathbb{R}, \mathbf{x} \in \mathbb{R}^n$$

$$\mathbf{x} \leftarrow \alpha \mathbf{x}$$

▶ A host function would look something like this:

```
void ax(int n, double alpha, double *x)
{
   for (int i = 0; i < n; i++)
        x[i] = alpha * x[i];
}</pre>
```







AX example (kernel)

A matching kernel is still relatively simple:

- What are blockIdx.x, blockDim.x and threadIdx.x?
- ▶ Where is the for loop?
- Why is there a if block?

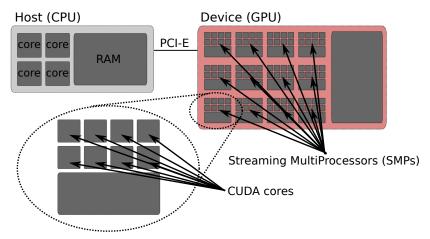






AX example (CUDA cores and SMPs)

Let's look at the earlier figure again...









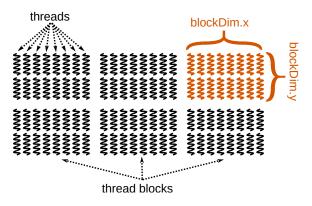


AX example (CUDA cores and SMPs)

- Each Streaming MultiProcessor (SMP) consists of several CUDA cores.
- Each CUDA core can execute several threads simultaneously.
 - The scheduler select the next instruction among a pool of active threads.
- Thus, the total number of threads can be in the millions.
- ▶ How do we decide what each thread should do?
- How do we manage all these threads?
 - Different problems sizes might require different number of threads.
 - Different GPUs might have different number of SMPs and CUDA cores.



The threads are divided into thread blocks:



blockDim.x and blockDim.y are the dimensions of the thread blocks.

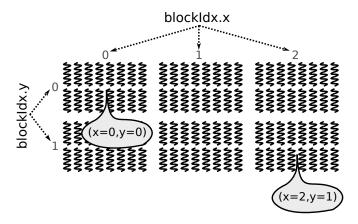








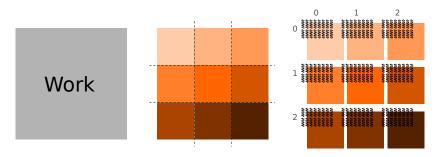
Each thread block gets an index number in a grid:



gridDim.x and gridDim.y are the dimensions of the grid.



- ▶ The overall idea is to **partition** the work into self-contained tasks.
- Each task is assign to one thread block.
 - The thread block indices are used to identify the task.

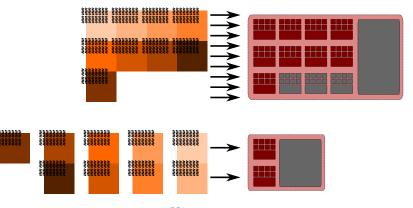








- The CUDA runtime is responsible for scheduling the thread blocks to SMPs.
- The execution order of the thread blocks is relaxed.
 - ► The code can therefore adapt to different GPUs:



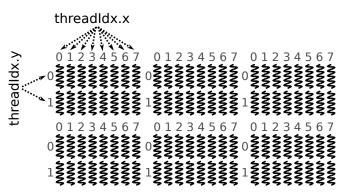








Each thread gets a **local** index number:







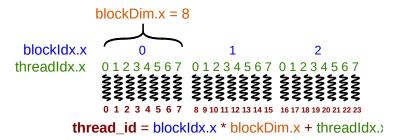




AX example (global indexing)

➤ A unique global index number can be calculated for each thread:

```
__global__ void ax_kernel(int n, double alpha, double *x)
{
    // query the global thread index
    int thread_id = blockIdx.x * blockDim.x + threadIdx.x;
    if (thread_id < n)
        x[thread_id] = alpha * x[thread_id];
}</pre>
```



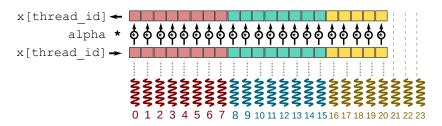


AX example (global indexing)

▶ The if block is used to filter out excess threads:

```
| __global__ void ax_kernel(int n, double alpha, double *x)
{
    // query the global thread index
    int thread_id = blockIdx.x * blockDim.x + threadIdx.x;

    // each thread updates one row
    if (thread_id < n)
        x[thread_id] = alpha * x[thread_id];
}</pre>
```



AX example (remarks)

- The thread blocks should be reasonably large.
 - Usually 32 threads is too small for practical use.
 - At the moment, the upper limit is 1024 thread.
 - My personal recommendation is to start from 256 threads, and tune the number if necessary.
 - Multiples of 32 are preferred (warps).
- The grid should be reasonably large.
 - Each thread block runs on a single SMP.
 - For optimal performance, each SMP should get a thread block.
 - Nvidia Tesla V100 GPU has 80 SMPs.
- ▶ Given a thread block of the size (Dx, Dy, Dz), the hardware indexes a thread of index (x, y, z) as (x + y*Dx + z*Dx*Dy).



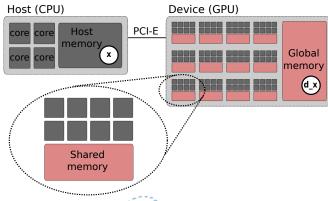




AX example (memory spaces)

► The host manages the memory:

```
double *x = (double *) malloc(n*sizeof(double));
for (int i = 0; i < n; i++)
x[i] = i;
cudaMalloc(&d_x, n*sizeof(double));
```







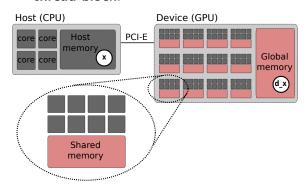




AX example (memory spaces)

Host memory is accessible by the host (and sometimes by all threads in all thread blocks).

Global memory is accessible by all threads in all thread blocks. Shared memory is accessible by threads that **belong to a same** thread block









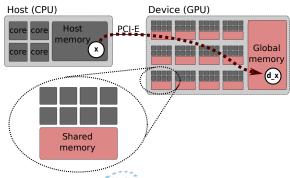


AX example (memory transfer to device)

► The host initializes a data transfer from the host memory to the global memory:

```
cudaMemcpy(d_x, x, n*sizeof(double), cudaMemcpyHostToDevice);
```

- The cudaMemcpy call is blocking.
 - ▶ The host waits until the transfer is ready.









AX example (kernel launch)

The host issues the ax_kernel kernel:

```
dim3 threads = 256;
dim3 blocks = (n+threads.x-1)/threads.x; // gridDim.x

ax_kernel <<< blocks, threads >>>(n, alpha, d_x);
```

▶ If the kernel used multi-dimensional threads blocks, then

```
| dim3 threads(Dx, Dy, Dz);
| dim3 blocks(Gx, Gy, Gz)
| ax_kernel<<<br/>blocks, threads>>>(n, alpha, d_x);
```

would create a $Gx \times Gy \times Gz$ grid of thread blocks, with Dx \times Dy \times Dz threads in each block.









AX example (kernel launch)

Alternatively, we could have written

```
| ax_kernel <<<(n+255)/256, 256>>>(n, alpha, d_x);
```

▶ (n+blockDim.x-1)/blockDim.x is simply a convenient way of making sure that

```
n < gridDim.x * blockDim.x.
```

I personally prefer the following approach:

```
// a function that returns the ceil of a/b. That is,
// DIVCEIL(5, 2) = ceil(5/2) = ceil(2.5) = 3.
static int DIVCEIL(int a, int b)
{
    return (a+b-1)/b;
}
```



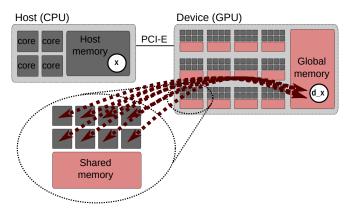






AX example (kernel launch)

▶ When the kernel finally starts, the threads access the data from the global memory:







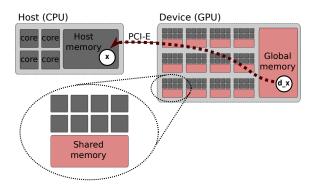




AX example (memory transfers from device)

▶ The host initializes a data transfer from the global memory to the host memory:

```
cudaMemcpy(x, d_x, n*sizeof(double), cudaMemcpyDeviceToHost)
```











AX example (cleanup)

Finally, we must free the allocated memory:

- All memory that has been allocated using the cudaMalloc function **must** the freed with the cudaFree function.
- The following will cause a segmentation fault:









AX example (compile and run)

- Load the correct toolchain:
 - \$ ml purge
 - \$ ml fosscuda/2019b buildenv
- Compile the source code with nvcc:
 - \$ nvcc -o ax ax.cu
- Queue a job:

```
$ srun --account=SNIC2020-9-161 \
--reservation=snic2020-9-161-day1 \
--ntasks=1 --gres=gpu:v100:1,gpuexcl \
--time=00:02:00 ./ax 10000
Residual = 0.000000e+00
```







Error handling (queries)

- Most CUDA functions return an error code of the type cudaError_t.
- A successful function call returns cudaSuccess. Other values indicate error.
- A kernel launch does not return anything.
 - Any errors must be gueried separately. See below.
- ▶ The previous error code can be checked and resetted with:

```
|| __host__ __device__ cudaError_t cudaGetLastError()
```







Error handling (queries)

▶ The previous error code can be checked without resetting:

```
| __host__ __device__ cudaError_t cudaPeekAtLastError()
```

An error code can be turned into a string:

```
| __host__ __device__ const char* cudaGetErrorName(cudaError_t error)
```

An error code can be turned into a longer description:

```
|| __host__ __device__ const char* cudaGetErrorString(cudaError_t error)
```





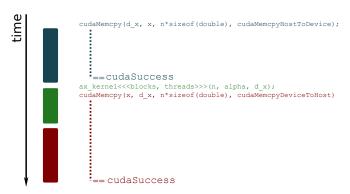




Error handling (some notes)

- Kernel launches and many other CUDA functions (*Async) are non-blocking / asynchronous.
 - The kernel or the function call is simply placed into a stream.

device host











Error handling (some notes)

It is possible that the returned error code is related to one of the earlier kernels or function calls!

device host

```
cudaMemcpy(d_x, x, n*sizeof(double), cudaMemcpyHostToDevice);
  ==cudaSuccess
ax kernel << blocks, threads>>> (n, alpha, d x);
cudaMemcpy(x, d_x, n*sizeof(double), cudaMemcpyDeviceToHost)
```









Error handling (some notes)

This can happen even when the outcome of the kernel launch is checked:

device host

```
cudaMemcpy(d x, x, n*sizeof(double), cudaMemcpyHostToDevice);
ax kernel << <blocks, threads>>> (n, alpha, d x);
cudaGetLastError() == cudaSuccess
cudaMemcpy(x, d x, n*sizeof(double), cudaMemcpyDeviceToHost)
```

Only errors that occurred during the kernel launch are returned by cudaGetLastError().









Hands-ons

- Three hands-ons under hands-ons/1.basics:
 - 1.threads Learn how to launch a kernel. Learn how to coordinate threads and thread blocks.
 - 2 errors. Learn how to detect and handle errors.
 - 3.memory Learn how to allocate global memory and transfer data to/from a device.
- Solutions can be found under solutions/1.basics.







