# A Very Quick Introduction to CUDA 

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## Hardware Basics

CPU


GPU


## DRAM

- CPUs are latency oriented (minimize execution of serial code)
- GPUs are throughput oriented (maximize number of floating point operations)


## CPU vs GPU threads <br> core 1 <br> core 2



- If the CPU has $n$ cores, each core processes $1 / n$ elements
- Launching, scheduling threads adds overhead

- GPUs process one element per thread
- Scheduled by GPU hardware, not by OS


## CUDA C

- Compute Unified Device Architecture
- NVIDIA GPUs can be programmed by CUDA, extension of C language (CUDA Fortran is also available)
- CUDA C is compiled with nvcc
- Numerical libraries: cuBLAS, cuFFT, Magma, ...
- Host $->$ CPU; Device $->$ GPU (They do not share memory!)
- The HOST launches a kernel that execute on the DEVICE
- A kernel is a data-parallel computation, executed by many threads.
- The number of threads are very large ( $\sim 1000$ or more)

Thread Organization

Block 0


Block 1


Block 2


Block n-1


## CUDA C

- Threads are grouped into blocks.
- Each block shares memory.

Eg. Vector addition:

```
int main(void) {
    vecAdd<<< blocksPerGrid, THREADS_PER_BLOCK}>>> (d_A, d_B, d_C)
}
__global__ static void vecAdd (float *a, float * b, float * c){
    }
```

The __global__ qualifer alerts the compiler that the code block will run on the DEVICE, but can be called from the HOST.

## CUDA C

- Grids and threads can also be arranged in 2d arrays (useful for image processing)
dim3 blocks(2,2)
dim3 threads $(16,16)$
kernel <<< blocks, threads >>>( );



## Code Example - 1

## Hello World!

\#include <stdio.h>
__device__ const char *STR = "HELLO WORLD!";
const int STR_LENGTH = 12;
global__ void hello()
printf("\%c $\backslash$ ", STR[threadId.x \% STR_LENGTH]);
\}
int main(void)\{
int threads_per_block = STR_LENGHT;
int blocks_per_grid = l;
hello <<< blocks_per_grid, threads_per_block >>> ();
cudaDeviceSynchronize();
return 0 ;
\}


Halt host thread execution on CPU until the device has finished processing all previously requested tasks.

## Code Example - 2

## Vector Addition (Very large vectors)

```
__global__ void add( int *a, int *b, int *c){
    int tid = threadIdx.x + blockIdx.x * blockDim.x ; // handle the data at this index
    while (tid < N) {
        c[tid] = a[tid] + b[tid];
        tid += blockDim.x * gridDim.x;
    }
}
```

e.g.: blockDim $=4$, gridDim $=4$

$$
\begin{aligned}
\text { tid } & =\text { th.id + blk.id * blk.dim } \\
& =1+1 * 4 \\
& =5
\end{aligned}
$$

|  | th 0 | th 1 | th 2 | th 3 |
| :---: | :---: | :---: | :---: | :---: |
| block 0 |  |  |  |  |
| block 1 |  |  |  |  |
| block 2 |  |  |  |  |
| block 3 |  |  |  |  |

## Code Example - 2

## Vector Addition (Very large vectors)

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__global__ void add( int *a, int *b, int *c){
    int tid = threadIdx.x + blockIdx.x * blockDim.x ; // handle the data at this index
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        c[tid] = a[tid] + b[tid];
        tid += blockDim.x * gridDim.x;
    }
}
```

e.g.: $\mathbf{N}=256$, blockDim $=2$, gridDim $=2 \longrightarrow$ offset $=$ blockDim * gridDim
a

blockDim * gridDim
$+$
b


C


## Code Example - 2

- Define arrays to be used on the HOST, and allocate memory.

```
int a[N], b[N], c[N];
int *dev_a, *dev_b, *dev_c;
// Allocate memory on the GPU
cudaMalloc( (void**)&dev_a, N * sizeof(int) );
cudaMalloc( (void**)&dev_b, N * sizeof(int) );
cudaMalloc( (void**)&dev_c, N * sizeof(int) );
```

- Copy arrays to the DEVICE

```
//Copy the arrays 'a' and 'b' to the GPU
cudaMemcpy( dev_a, a, N * sizeof(int), cudaMemcpyHostToDevice );
cudaMemcpy( dev_b, b, N * sizeof(int), cudaMemcpyHostToDevice );
```

- Launch the kernel, then copy result from DEVICE to HOST

```
add<<<128,128>>>>( dev_a, dev_b, dev_c) ; // Launch N=128 blocks each containing M=128 threads
//Copy the array 'c' back from the GPU to the CPU
cudaMemcpy( c, dev_c, N * sizeof(int), cudaMemcpyDeviceToHost );
```

- Free memory

```
//Free memory
cudaFree(dev_a);
cudaFree(dev_b);
cudaFree(dev_c);
```


## Code Example - 3

## Dot product



- Recall, each Block shares memory!
- Each block will have a its own copy of cahce [], i.e. a partial result.
- Final step is reduction, i.e. summing all the partial results in cahce[] to obtain a final answer.


## Code Example - 3

```
// Reduction (even number of threads assumed)
int i = blockDim.x/2;
while ( i != 0 ){
    if (cacheIndex < i)
        cache[cacheIndex] += cache[cacheIndex + i];
    _syncthreads();
    i /= 2;
}
```

```
// Write to c using the threadId 0
if (cacheIndex == 0)
    c[blockIdx.x] = cache[0];
```

Parallel reduction

Finally, write the final answer, with one thread (serial).

Parallel reduction: (Not the best one!)


## GPUs on Comet

- 1944 Standard compute nodes
- 36 GPU Nodes:
- Intel Xeon E5-2680v3
- NVIDIA K80 GPUs (11GB)


GPU Examples:
/share/apps/examples/GPU

## GPUs on Comet

\$ module load cuda
\$ nvcc -o hello_cuda.x hello_cuda.cu
cuda.job

```
#!/bin/bash
#SBATCH -p gpu-shared
#SBATCH -gres=gpu:1
#SBATCH -job-name="hellocuda"
#SBATCH -output="hellocuda.%j.%N.out"
#SBATCH -t 00:01:00
#SBATCH -A TG-SEE150004
cd ~/Working_directory
./hello_cuda.x
```

\$ sbatch cuda.job

## Exercise

Examine and run the code add_vec_times.cu and compare it with add_vec_spu_thd-blk.cu and answer the following questions:

- Vary THREADS_PER_BLOCK: $1,2,4,8,16,32,64,128,256$
- Record the time printed

1. How many blocks are launched for each case?
2. Until what value the timing decreases linearly?
3. What is the explanation of the loss of the linear behavior after this value? (Hint: search for "warps")
