INTRODUCTION TO PROGRAMMING IN CUDA C

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BASIC C PROGRAM

```
#include <iostream>

int main ( void ) {
  printf( "Hello, World!\n" );
  return 0;
}
```

BASIC CUDA C PROGRAM

```
#include <iostream>

__global__ void kernel ( void ) {
}

int main ( void ) {
  kernel<<<1,1>>>();
  printf( "Hello, World!\n" );
  return 0;
}
```

```
#include <iostream>

__global__ void kernel ( void ) {
}
```

Prefix that says: "only run this function on the GPU"

```
int main ( void ) {
  kernel<<<1,1>>>();
  printf( "Hello, World!\n" );
  return 0;
}
```

__global__: Call from CPU and run only on GPU.

__device__: Call from GPU and run only on GPU.
(More specifically, call only from within
a __global__ or another __device__ function.)

__host__: Call from CPU and run only on CPU.
(i.e., a traditional C function.)

	Executed on the:	Only callable from the:
__device__ float DeviceFunc()	device	device
__global__ void KernelFunc()	device	host
__host__ float HostFunc()	host	host

```
__device__ int dev1( void ){
}

__device__ int dev2( void ){
}

__global__ void kernel ( void ) {
  dev1();
  dev2();
}

int main ( void ) {
  kernel<<<1, 1>>>();
  printf( "Hello, World!\n" );
  return 0;
}
```

```
#include <iostream>

__global__ void kernel ( void ) {
}

int main ( void ) {
  kernel<<<1,1>>>();
  printf( "Hello, World!\n" );
  return 0;
}
```

Started on CPU

Sent to GPU

Done on CPU

Done on CPU

```
#include <iostream>

__global__ void kernel ( void ) {
}

int main ( void ) {
  kernel<<<1,1>>>(); ?
  printf( "Hello, World!\n" );
  return 0;
}
```

REVIEW OF PARALLELIZATION

Parallelization: Running different calculations simultaneously.

Kernel: An instruction set executed on the GPU. (All others are executed on the CPU.)

In CUDA C, a kernel is any function prefixed with the keyword, `__global__`. For example:

```
__global__ void kernel ( void ) {
}
```

PARALLELIZATION AND THE SIMD PARADIGM

SIMD: Single Instruction Multiple Data. The paradigm works like this:

1. The CPU sends a kernel (a SINGLE INSTRUCTION set) to the GPU.
2. The GPU executes the kernel MULTIPLE times simultaneously in **PARALLEL**. Each such execution of the kernel is called a **thread**, and each thread is executed on different parts of a giant data set.

PARALLELIZING A WORKLOAD ON A GPU

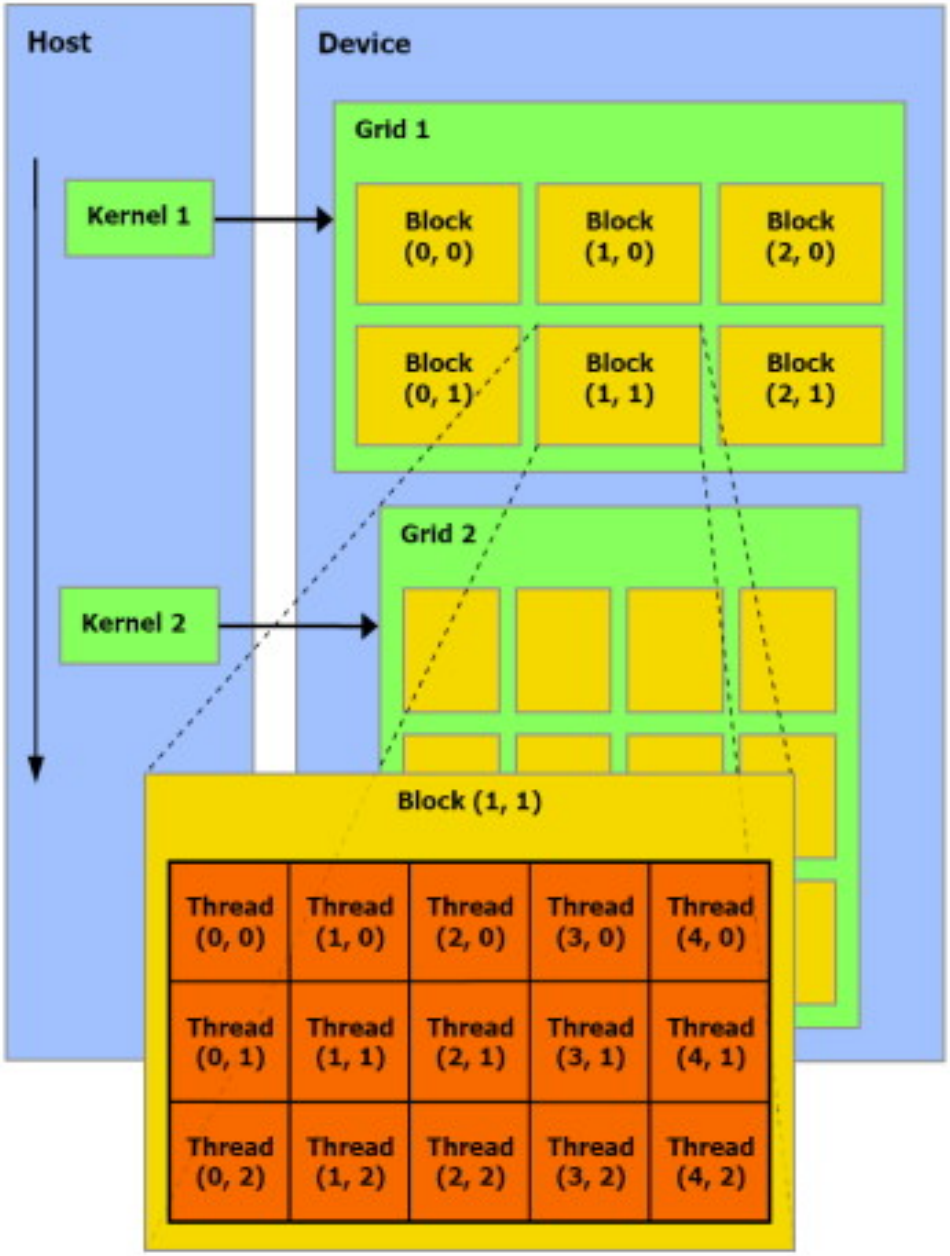
1. The CPU sends a CPU-to-GPU command called a **kernel** to a single GPU core.

2. The GPU core multitasks to execute the command:

   a. The GPU makes $N \cdot M$ copies of the kernel's code, and then runs all those copies simultaneously. Those parallel copies are called **threads**.

   b. The $N \cdot M$ threads are partitioned into N groups, called **blocks**, of M threads each.

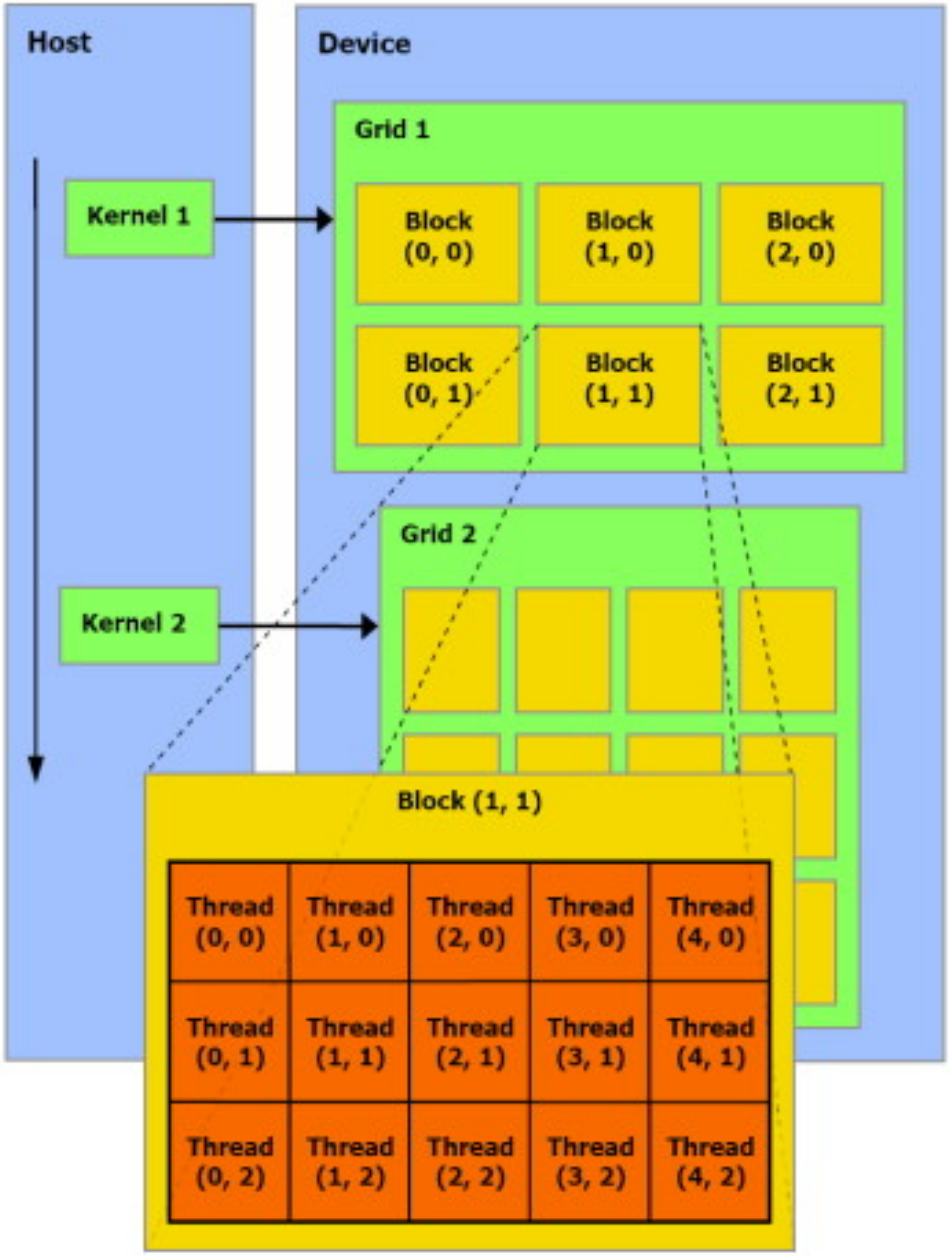
   c. The sum total of all the threads from a kernel call is a **grid**.


Host
Device
Grid 1
Kernel 1
Block (0, 0)
Block (1, 0)
Block (2, 0)
Block (0, 1)
Block (1, 1)
Block (2, 1)
Grid 2
Kernel 2
Block (1, 1)
Thread (0, 0)
Thread (1, 0)
Thread (2, 0)
Thread (3, 0)
Thread (4, 0)
Thread (0, 1)
Thread (1, 1)
Thread (2, 1)
Thread (3, 1)
Thread (4, 1)
Thread (0, 2)
Thread (1, 2)
Thread (2, 2)
Thread (3, 2)
Thread (4, 2)

ORGANIZATION OF THREADS

Grid; The collection of all the threads that are spawned when the CPU sends a kernel to the GPU.

Block: A collection of threads within a grid that share memory quickly and easily.


Host
Device
Grid 1
Kernel 1
Block (0, 0)
Block (1, 0)
Block (2, 0)
Block (0, 1)
Block (1, 1)
Block (2, 1)
Grid 2
Kernel 2
Block (1, 1)
Thread (0, 0)
Thread (1, 0)
Thread (2, 0)
Thread (3, 0)
Thread (4, 0)
Thread (0, 1)
Thread (1, 1)
Thread (2, 1)
Thread (3, 1)
Thread (4, 1)
Thread (0, 2)
Thread (1, 2)
Thread (2, 2)
Thread (3, 2)
Thread (4, 2)

NOW BACK TO THE STUFF IN ANGLE BRACKTES...

```
#include <iostream>

__global__ void kernel ( void ) {
}

int main ( void ) {
  kernel<<<1,1>>>();
  printf( "Hello, World!\n" );
  return 0;
}
```

Uses one grid.

Number of blocks

Number of threads per block

Here, the GPU runs kernel() one time:

```
#include <iostream>

__global__ void kernel ( void ) {
}

int main ( void ) {
  kernel<<<1,1>>>();
  printf( "Hello, World!" \n" )
  return 0;
}
```

Here, the GPU runs kernel() 5 times:

```
#include <iostream>

__global__ void kernel ( void ) {
}

int main ( void ) {
  kernel<<<5,1>>>();
  printf( "Hello, World!" \n" )
  return 0;
}
```

Here, the GPU runs kernel() 5 times:

```
#include <iostream>

__global__ void kernel ( void ) {
}

int main ( void ) {
  kernel<<<1,5>>>();
  printf( "Hello, World!" \n" )
  return 0;
}
```

Here, the GPU runs kernel() 20 times:

```
#include <iostream>

__global__ void kernel ( void ) {
}

int main ( void ) {
  kernel<<<4,5>>>();
  printf( "Hello, World!" \n" )
  return 0;
}
```

PASSING DATA TO AND FROM THE GPU: simple1.cu

```
#include <stdio.h>
#include <stdlib.h>

__global__ void colonel(int *dev_a){
  *dev_a = 1;
}

int main(){

  // Declare variables and allocate memory on the GPU.
  int a[1], *dev_a;
  cudaMalloc((void**) &dev_a, sizeof(int));

  // Execute kernel and copy the result to CPU memory.
  colonel<<<1,1>>>(dev_a);
  cudaMemcpy(a, dev_a, sizeof(int), cudaMemcpyDeviceToHost);

  // Print result and free dynamically allocated memory.
  printf("a[0] = %d\n", a[0]); // REMEMBER: INDEXING IN C STARTS FROM 0.
  cudaFree(dev_a);

}
```

```
[landau@impact1 simple]$ nvcc simple1.cu -o simple1.out
[landau@impact1 simple]$ ./simple1.out
a[0] = 1
[landau@impact1 simple]$
```

PASSING DATA TO AND FROM THE GPU: simple2.cu

```
#include <stdio.h>
#include <stdlib.h>

__global__ void colonel(int *dev_a){
  *dev_a = *dev_a + 1;
}

int main(){
  // Declare variables and allocate memory on the GPU.
  int a[1], *dev_a;
  cudaMalloc((void**) &dev_a, sizeof(int));

  // Intitialize argument a, executed kernel, and store result back in a.
  a[0] = 1; // REMEMBER: INDEXING IN C STARTS FROM 0.
  cudaMemcpy(dev_a, a, sizeof(int), cudaMemcpyHostToDevice);
  colonel<<<1,1>>>(dev_a);
  cudaMemcpy(a, dev_a, sizeof(int), cudaMemcpyDeviceToHost);

  // Print result and free dynamically allocated memory.
  printf("a[0] = %d\n", a[0]); // REMEMBER: INDEXING IN C STARTS FROM 0.
  cudaFree(dev_a);
}
```

```
[landau@impact1 simple]$ nvcc simple2.cu -o simple2.out
[landau@impact1 simple]$ ./simple2.out
b[0] = 2
[landau@impact1 simple]$
```

threadIdx.x AND blockIdx.x

Say the CPU sends a kernel to the GPU with B blocks and T threads per block.

The following CUDA C variables are available for use in the kernel:

blockIdx.x: the block ID corresponding to the current thread, an integer from 0 to $B - 1$ inclusive.

threadIdx.x: the thread ID of the current thread within its block, an integer from 0 to $T - 1$ inclusive.

NOTE: there are also variables like blockIdx.y and threadIdx.y, but those are only useful if you deliberately organize your blocks and threads in two-dimensional arrays (which can be done).

USING blockIdx.x AND threadIdx.x: identify.cu

```
#include <stdio.h>
#include <stdlib.h>

__global__ void isExecuted(int *dev_a, int blockid, int threadid){

  if(blockIdx.x == blockid && threadIdx.x == threadid)
    *dev_a = 1;

}

int main(){

  // Declare variables and allocate memory on the GPU.
  int init[1], a[1], *dev_a;
```

```
  cudaMalloc((void**) &dev_a, sizeof(int));

  // Initialize dev_a, execute kernel, and copy the result to CPU memory.
  init[1] = 0;
  cudaMemcpy(dev_a, init, sizeof(int), cudaMemcpyHostToDevice);
  isExecuted<<<100,100>>>(dev_a, 2, 4);
        // NOTE: INDEXING OF THREADS AND BLOCKS STARTS FROM 0.
  cudaMemcpy(a, dev_a, sizeof(int), cudaMemcpyDeviceToHost);

  // Print result and free dynamically allocated memory.
  printf("a[0] = %d\n", a[0]); // REMEMBER: INDEXING IN C STARTS FROM 0.
  cudaFree(dev_a);

}
```

```
[landau@impact1 simple]$ nvcc identify.cu -o identify.out
[landau@impact1 simple]$ ./identify.out
a[0] = 1
[landau@impact1 simple]$
```


isExecuted<<<2,3>>>(dev_a, blockid = 1, threadid = 1);
isExecuted(dev_a, blockid = 1, threadid = 1, blockIdx.x = 0, threadIdx.x = 0, ...);
*dev_a is unchanged.
isExecuted(dev_a, blockid = 1, threadid = 1, blockIdx.x = 0, threadIdx.x = 1, ...);
*dev_a is unchanged.
isExecuted(dev_a, blockid = 1, threadid = 1, blockIdx.x = 0, threadIdx.x = 2, ...);
*dev_a is unchanged.
isExecuted(dev_a, blockid = 1, threadid = 1, blockIdx.x = 1, threadIdx.x = 0, ...);
*dev_a is unchanged.
isExecuted(dev_a, blockid = 1, threadid = 1, blockIdx.x = 1, threadIdx.x = 1, ...);
*dev_a is set to 1.
isExecuted(dev_a, blockid = 1, threadid = 1, blockIdx.x = 1, threadIdx.x = 2, ...);
*dev_a is unchanged.

You can verify that when the line,

```
isExecuted<<<1,1>>>(dev_a, 0, 0);
```

is changed to:

```
isExecuted<<<1,1>>>(dev_a, 2, 4);
```

the output of the program:

```
a[0] = 0
```

which makes sense because there was no block number 2 or thread number 4.

ASIDE: AVOIDING PITFALLS IN PARALLELIZATION

What problem would arise if I used the function, isExectued_goofup(), instead of Executed():

```
__global__ void isExecuted_goofup(int *dev_a, int blockid,
                int threadid){

  if(blockIdx.x == blockid && threadIdx.x == threadid)
    *dev_a = 1;
  else
    *dev_a = 0;

}
```

If block 1 thread 1 finishes last, then *dev_a = 1.
If block 1 thread 1 does not finish last, *dev_a = 0.

We can't know or control which thread finishes last!

LECTURE SERIES MATERIALS

These lecture slides, a tentative syllabus for the whole lecture series, and code are available at:

https://github.com/wlandau/gpu.

After logging into you home directory on impact1, type:

```
git clone https://github.com/wlandau/gpu
```

into the command line to download all the course materials.

REFERENCES

David B. Kirk and Wen-mei W. Hwu. "Programming Massively Parallel Processors: a Hands-on Approach." Morgan Kaufman, 2010.

J. Sanders and E. Kandrot. *CUDA by Example.* Addison-Wesley, 2010.