

# 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:
<code>__device__ float DeviceFunc()</code>	device	device
<code>__global__ void KernelFunc()</code>	device	host
<code>__host__ float HostFunc()</code>	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 ) {  Started on CPU
    kernel<<<1,1>>>();  Sent to GPU
    printf( "Hello, World!\n" );  Done
    return 0;  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

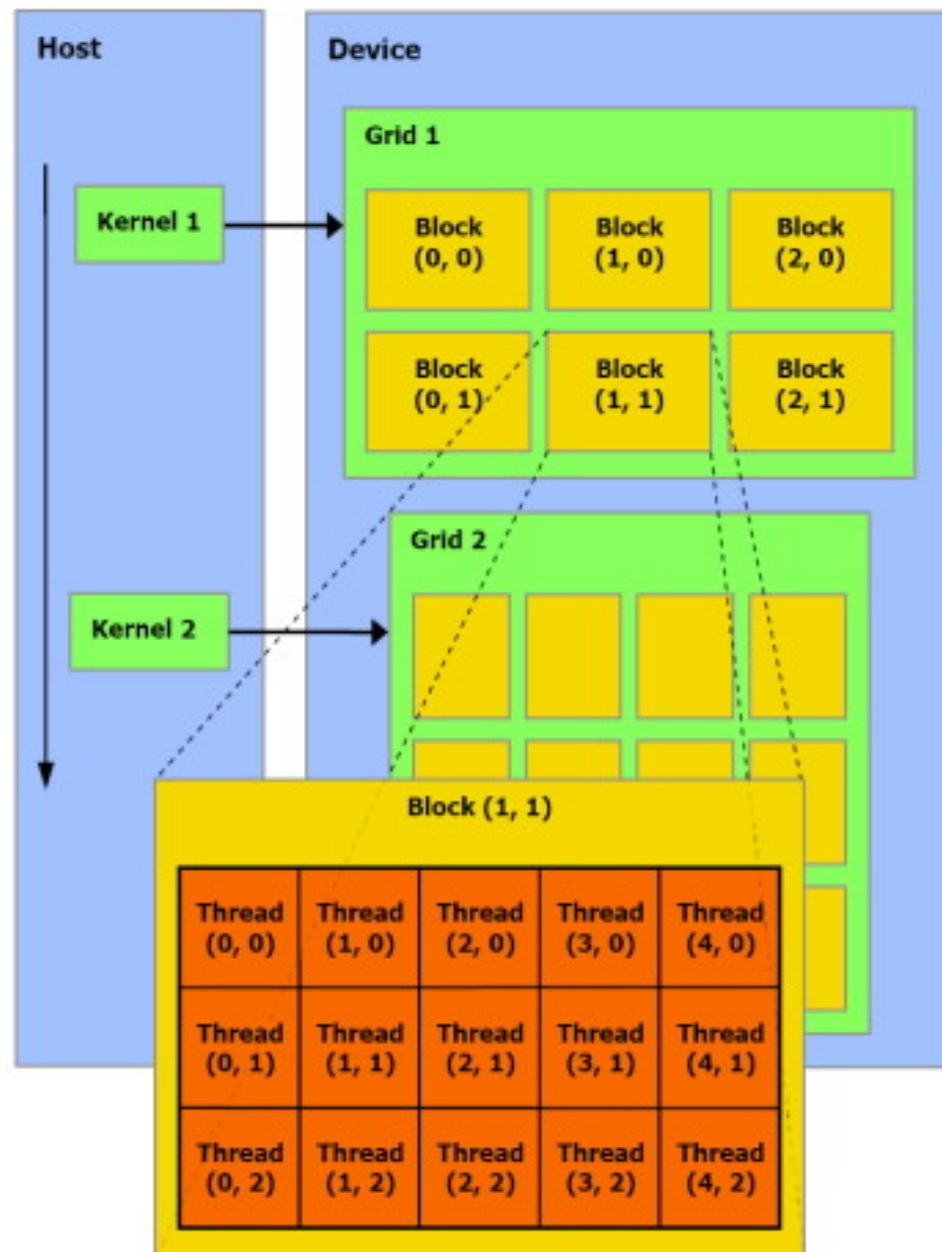
**SIMD**: Single Instruction Multiple Data. The paradigm works like this:

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.

# 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.



# NOW BACK TO THE STUFF IN ANGLE BRACKETS...

```
#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));

    // Initialize 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]$ |
```





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:

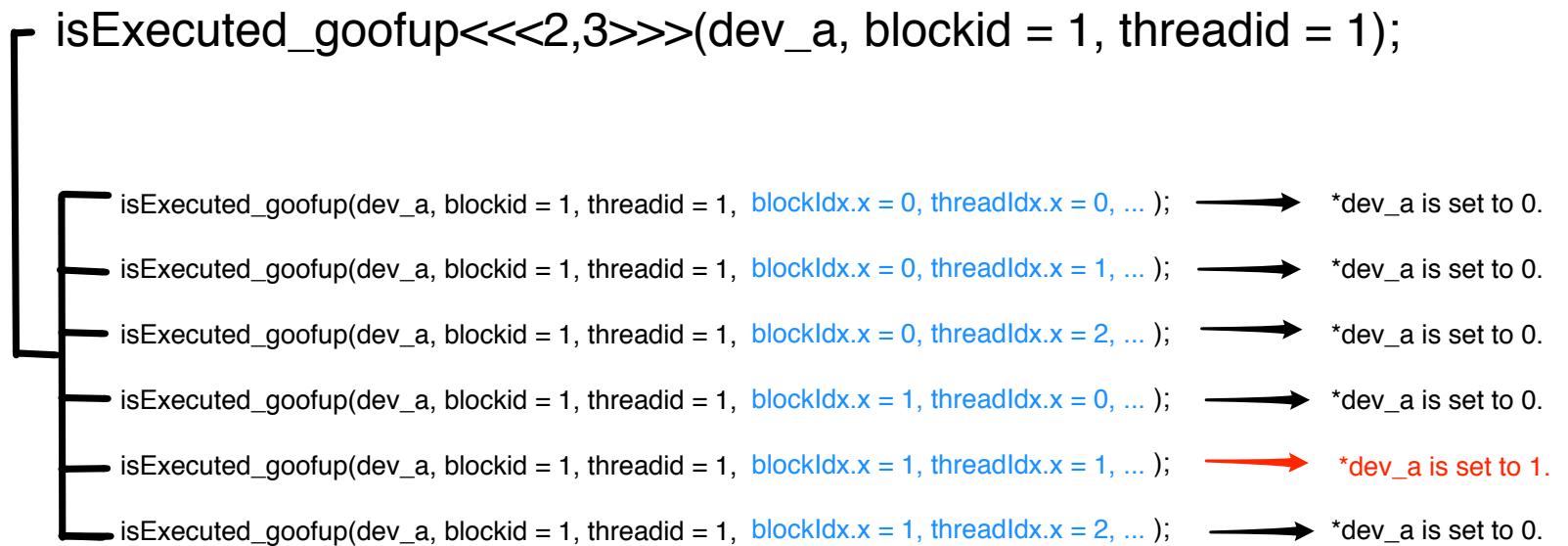
```
a[0] = 0
```

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!



## 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.