

# 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;  
}
```

```
#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.

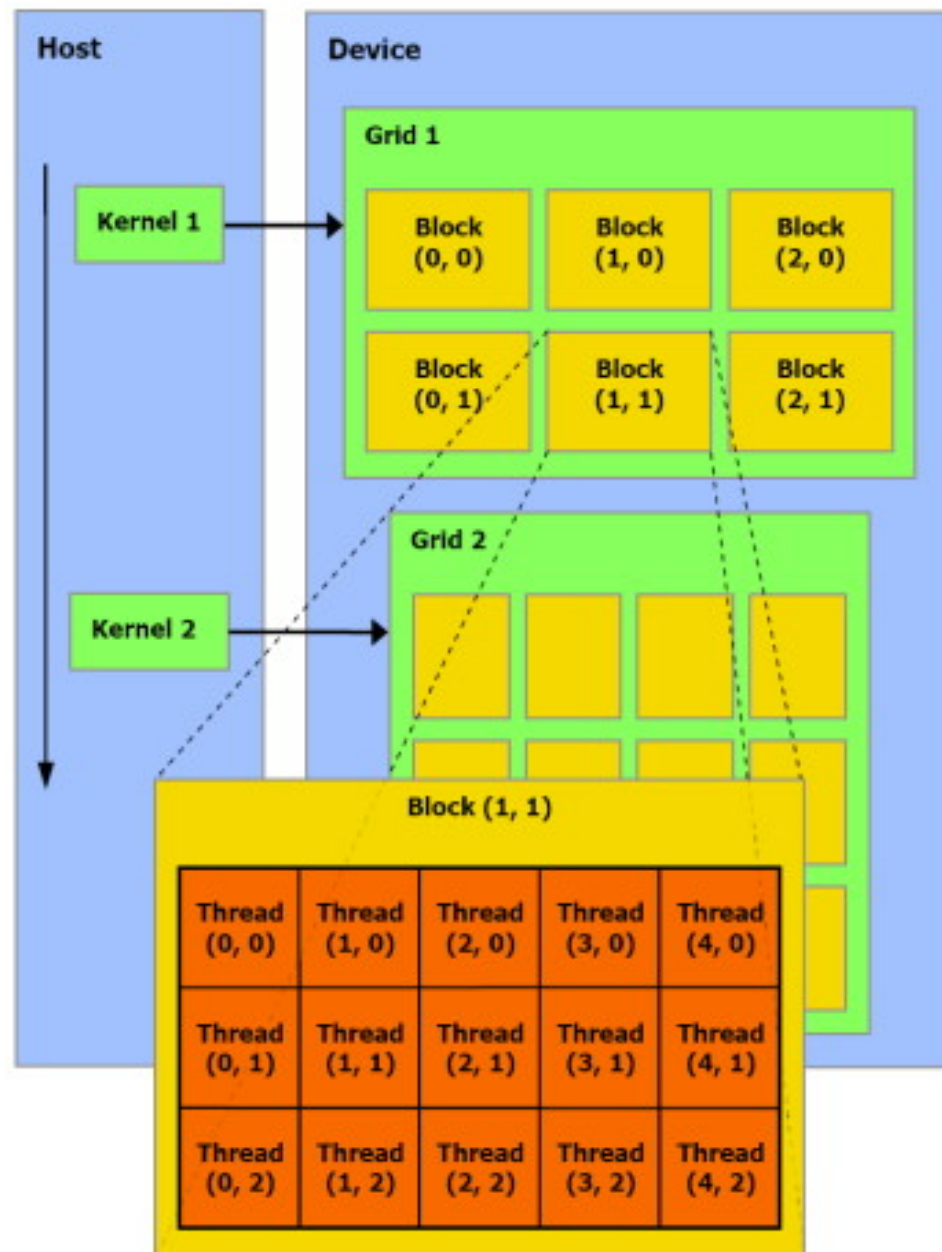




# ORGANIZATION OF THREADS

**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 changed isExecuted() to:

```
__global__ void isExecuted(int *dev_a, int blockid,  
                           int threadid){  
  
    if(blockIdx.x == blockid && threadIdx.x == threadid)  
        *dev_a = 1;  
    else  
        *dev_a = 0;  
  
}
```

```

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

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

```

Suppose:

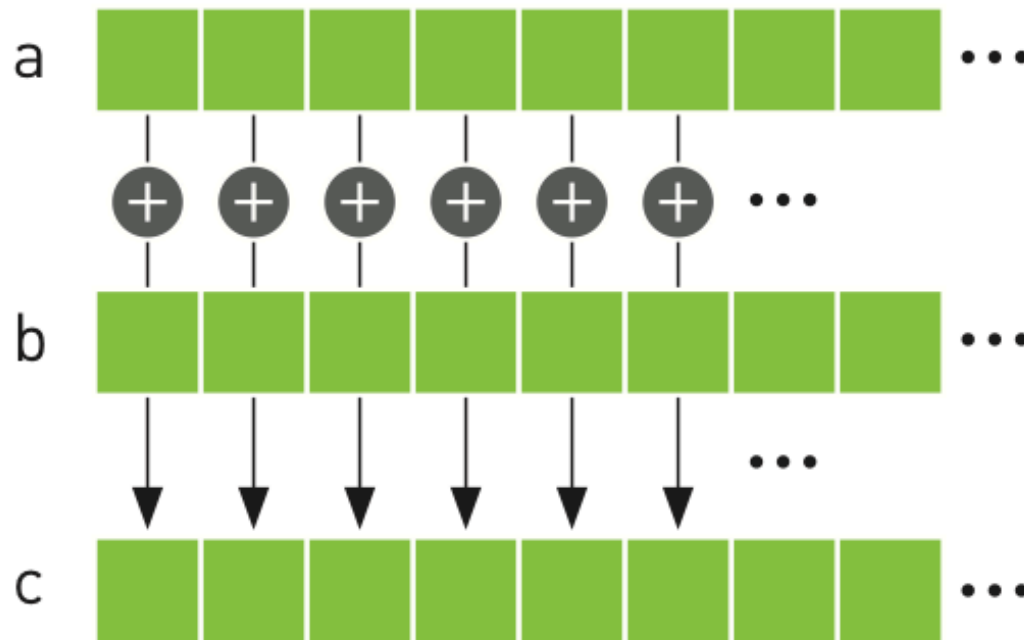
- I run `isExecuted(dev_a, 0, 0)`
- Block 0 thread 0 finishes before Block 0 thread 1 finishes.

Then, here's what would happen:

1. Block 0 thread 0 would set `*dev_a` to 1 because our arguments are `blockid = 0` and `threadid = 0`.
2. Block 0 thread 1 would set `*dev_a` back to 0 because `threadid` does not equal 1.

So `isExecuted(dev_a, 1, 1)` might return 0 even it block 1 thread 1 is among the threads that executed.

## EXAMPLE: VECTOR SUMMATION (Sanders, et. al.)



*Figure 4.1* Summing two vectors

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

```

#define N    10

int main( void ) {
    int a[N], b[N], c[N];
    int *dev_a, *dev_b, *dev_c;

    // allocate the memory on the GPU
    HANDLE_ERROR( cudaMalloc( (void**)&dev_a, N * sizeof(int) ) );
    HANDLE_ERROR( cudaMalloc( (void**)&dev_b, N * sizeof(int) ) );
    HANDLE_ERROR( cudaMalloc( (void**)&dev_c, N * sizeof(int) ) );

    // fill the arrays 'a' and 'b' on the CPU
    for (int i=0; i<N; i++) {
        a[i] = -i;
        b[i] = i * i;
    }
}

```

```

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

add<<<N,1>>>( dev_a, dev_b, dev_c );

// copy the array 'c' back from the GPU to the CPU
HANDLE_ERROR( cudaMemcpy( c, dev_c, N * sizeof(int),
                          cudaMemcpyDeviceToHost ) );

// display the results
for (int i=0; i<N; i++) {
    printf( "%d + %d = %d\n", a[i], b[i], c[i] );
}

```

```
// free the memory allocated on the GPU  
cudaFree( dev_a );  
cudaFree( dev_b );  
cudaFree( dev_c );  
  
return 0;  
}
```



```
[landau@impact1 vectorsums]$ nvcc vectorsums.cu -o vectorsums.out
[landau@impact1 vectorsums]$ ./vectorsums.out
0 + 0 = 0
-1 + 1 = 0
-2 + 4 = 2
-3 + 9 = 6
-4 + 16 = 12
-5 + 25 = 20
-6 + 36 = 30
-7 + 49 = 42
-8 + 64 = 56
-9 + 81 = 72
[landau@impact1 vectorsums]$
```

You can download the code, along with other simple CUDA C examples, at [https://github.com/jarad/gpuIntroduction/tree/master/CUDA\\_C\\_sandbox](https://github.com/jarad/gpuIntroduction/tree/master/CUDA_C_sandbox).

# REFERENCES

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