

# Parallel Programming & Cluster Computing

## **GPGPU: Number Crunching in Your Graphics Card**



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# Outline

- What is GPGPU?
- GPU Programming
- Digging Deeper: CUDA on NVIDIA
- CUDA Thread Hierarchy and Memory Hierarchy
- CUDA Example: Matrix-Matrix Multiply



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# What is GPGPU?



# Accelerators

No, not this ....



<http://gizmodo.com/5032891/nissans-eco-gas-pedal-fights-back-to-help-you-save-gas>



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# Accelerators

- In HPC, an accelerator is hardware component whose role is to speed up some aspect of the computing workload.
- In the olden days (1980s), supercomputers sometimes had *array processors*, which did vector operations on arrays, and PCs sometimes had *floating point accelerators*: little chips that did the floating point calculations in hardware rather than software.
- More recently, *Field Programmable Gate Arrays* (FPGAs) allow reprogramming deep into the hardware.





# Why Accelerators are Good

Accelerators are good because:

- they make your code run faster.



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# Why Accelerators are Bad

Accelerators are bad because:

- they're expensive;
- they're hard to program;
- your code on them isn't portable to other accelerators, so the labor you invest in programming them has a very short half-life.



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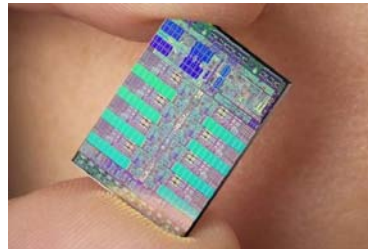


# The King of the Accelerators

The undisputed champion of accelerators is:  
the **graphics processing unit**.

[http://www.amd.com/us-en/assets/content\\_type/DigitalMedia/46928a\\_01\\_ATI-FirePro\\_V8700\\_angled\\_low\\_res.gif](http://www.amd.com/us-en/assets/content_type/DigitalMedia/46928a_01_ATI-FirePro_V8700_angled_low_res.gif)

[http://images.nvidia.com/products/quadro\\_fx\\_5800/Quadro\\_FX5800\\_low\\_3qtr.png](http://images.nvidia.com/products/quadro_fx_5800/Quadro_FX5800_low_3qtr.png)



<http://www.gamecyte.com/wp-content/uploads/2009/01/ibm-sony-toshiba-cell.jpg>



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# Why GPU?

- Graphics Processing Units (GPUs) were originally designed to accelerate graphics tasks like image rendering.
- They became very very popular with videogamers, because they've produced better and better images, and lightning fast.
- And, prices have been extremely good, ranging from three figures at the low end to four figures at the high end.



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# GPUs are Popular

- Chips are expensive to design (hundreds of millions of \$\$\$), expensive to build the factory for (billions of \$\$\$), but cheap to produce.
- In 2006 – 2007, GPUs sold at a rate of about 80 million cards per year, generating about \$20 billion per year in revenue.

[http://www.xbitlabs.com/news/video/display/20080404234228\\_Shipments\\_of\\_Discrete\\_Graphi\\_cs\\_Cards\\_on\\_the\\_Rise\\_but\\_Prices\\_Down\\_Jon\\_Peddie\\_Research.html](http://www.xbitlabs.com/news/video/display/20080404234228_Shipments_of_Discrete_Graphi_cs_Cards_on_the_Rise_but_Prices_Down_Jon_Peddie_Research.html)

- This means that the GPU companies have been able to recoup the huge fixed costs.





# GPU Do Arithmetic

- GPUs mostly do stuff like rendering images.
- This is done through mostly floating point arithmetic – the same stuff people use supercomputing for!



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# GPU Programming





# Hard to Program?

- In the olden days – that is, until just the last few years – programming GPUs meant either:
  - using a graphics standard like OpenGL (which is mostly meant for rendering), or
  - getting fairly deep into the graphics rendering pipeline.
- To use a GPU to do general purpose number crunching, you had to make your number crunching pretend to be graphics.
- This was hard. So most people didn't bother.





# Easy to Program?

More recently, GPU manufacturers have worked hard to make GPUs easier to use for general purpose computing.

This is known as *General Purpose Graphics Processing Units.*



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# How to Program a GPU

- Proprietary programming language or extensions
  - NVIDIA: CUDA (C/C++)
  - AMD/ATI: StreamSDK/Brook+ (C/C++)
- OpenCL (Open Computing Language): an industry standard for doing number crunching on GPUs.
- Portland Group Fortran and C compilers with accelerator directives.
- Others are popping up or in development now ....





# NVIDIA CUDA

- NVIDIA proprietary
- Formerly known as “Compute Unified Device Architecture”
- Extensions to C to allow better control of GPU capabilities
- Modest extensions but major rewriting of the code
- Portland Group Inc (PGI) now has a Fortran implementation of CUDA available in their Fortran compiler.







# CUDA Example Part 1

```
// example1.cpp : Defines the entry point for the console applicati
on.
//

#include "stdafx.h"

#include <stdio.h>
#include <cuda.h>

// Kernel that executes on the CUDA device
__global__ void square_array(float *a, int N)
{
    int idx = blockIdx.x * blockDim.x + threadIdx.x;
    if (idx<N) a[idx] = a[idx] * a[idx];
}
```

<http://llpanorama.wordpress.com/2008/05/21/my-first-cuda-program/>





# CUDA Example Part 2

```
// main routine that executes on the host
int main(void)
{
    float *a_h, *a_d; // Pointer to host & device arrays
    const int N = 10; // Number of elements in arrays
    size_t size = N * sizeof(float);
    a_h = (float *)malloc(size); // Allocate array on host
    cudaMalloc((void **) &a_d, size); // Allocate array on device
    // Initialize host array and copy it to CUDA device
    for (int i=0; i<N; i++) a_h[i] = (float)i;
    cudaMemcpy(a_d, a_h, size, cudaMemcpyHostToDevice);
    // Do calculation on device:
    int block_size = 4;
    int n_blocks = N/block_size + (N%block_size == 0 ? 0:1);
    square_array <<< n_blocks, block_size >>> (a_d, N);
    // Retrieve result from device and store it in host array
    cudaMemcpy(a_h, a_d, sizeof(float)*N, cudaMemcpyDeviceToHost);
    // Print results
    for (int i=0; i<N; i++) printf("%d %f\n", i, a_h[i]);
    // Cleanup
    free(a_h); cudaFree(a_d);
}
```





# AMD/ATI Brook+

- AMD/ATI proprietary
- Formerly known as “Close to Metal” (CTM)
- Extensions to C to allow better control of GPU capabilities
- No Fortran version available



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# Brook+ Example Part 1

```
float4 matmult_kernel (int y, int x, int k,  
                        float4 M0[], float4 M1[])  
{  
    float4 total = 0;  
    for (int c = 0; c < k / 4; c++)  
    {  
        total += M0[y][c] * M1[x][c];  
    }  
    return total;  
}
```

[http://developer.amd.com/gpu\\_assets/Stream\\_Computing\\_Overview.pdf](http://developer.amd.com/gpu_assets/Stream_Computing_Overview.pdf)





## Brook+ Example Part 2

```
void matmult (float4 A[], float4 B'[], float4 C[])
{
    for (int i = 0; i < n; i++)
    {
        for (j = 0; j < m / 4; j+)
        {
            launch_thread{
                C[i][j] =
                    matmult_kernel(j, i, k, A, B');}
        }
    }
    sync_threads{}
}
```





# OpenCL

- Open Computing Language
- Open standard developed by the Khronos Group, which is a consortium of many companies (including NVIDIA, AMD and Intel, but also lots of others)
- Initial version of OpenCL standard released in Dec 2008.
- Many companies will create their own implementations.
- Apple was first to market, with an OpenCL implementation included in Mac OS X v10.6 (“Snow Leopard”) in 2009.





# OpenCL Example Part 1

```
// create a compute context with GPU device
context = clCreateContextFromType(0, CL_DEVICE_TYPE_GPU, NULL, NULL,
    NULL);
// create a work-queue
queue = clCreateWorkQueue(context, NULL, NULL, 0);
// allocate the buffer memory objects
memobjs[0] =
    clCreateBuffer(context,
        CL_MEM_READ_ONLY | CL_MEM_COPY_HOST_PTR,
        sizeof(float)*2*num_entries, srcA);
memobjs[1] =
    clCreateBuffer(context, CL_MEM_READ_WRITE,
        sizeof(float)*2*num_entries, NULL);
// create the compute program
program =
    clCreateProgramFromSource(context, 1, &fft1D_1024_kernel_src, NULL);
// build the compute program executable
clBuildProgramExecutable(program, false, NULL, NULL);
// create the compute kernel
kernel = clCreateKernel(program, "fft1D_1024");
```





# OpenCL Example Part 2

```
// create N-D range object with work-item dimensions
global_work_size[0] = n;
local_work_size[0] = 64;
range = clCreateNDRangeContainer(context, 0, 1, global_work_size,
    local_work_size);
// set the args values
clSetKernelArg(kernel, 0, (void *)&memobjs[0], sizeof(cl_mem), NULL);
clSetKernelArg(kernel, 1, (void *)&memobjs[1], sizeof(cl_mem), NULL);
clSetKernelArg(kernel, 2, NULL,
    sizeof(float)*(local_work_size[0]+1)*16, NULL);
clSetKernelArg(kernel, 3, NULL,
    sizeof(float)*(local_work_size[0]+1)*16, NULL);
// execute kernel
clExecuteKernel(queue, kernel, NULL, range, NULL, 0, NULL);
```







# OpenCL Example Part 3

```
// This kernel computes FFT of length 1024. The 1024 length FFT
// is decomposed into calls to a radix 16 function, another
// radix 16 function and then a radix 4 function
kernel void fft1D_1024 (
    global float2 *in, __global float2 *out,
    local float *sMemx, __local float *sMemy)
{
    int tid = get_local_id(0);
    int blockIdx = get_group_id(0) * 1024 + tid;
    float2 data[16];
    // starting index of data to/from global memory
    in = in + blockIdx;
    out = out + blockIdx;
    globalLoads(data, in, 64); // coalesced global reads
```





# OpenCL Example Part 4

```
fftRadix16Pass(data);          // in-place radix-16 pass
twiddleFactorMul(data, tid, 1024, 0);
// local shuffle using local memory
localShuffle(data, sMemx, sMemy, tid,
              (((tid & 15) * 65) + (tid >> 4)));
fftRadix16Pass(data);          // in-place radix-16 pass
twiddleFactorMul(data, tid, 64, 4); // twiddle factor multiplication
localShuffle(data, sMemx, sMemy, tid,
              (((tid >> 4) * 64) + (tid & 15)));
// four radix-4 function calls
fftRadix4Pass(data);
fftRadix4Pass(data + 4);
fftRadix4Pass(data + 8);
fftRadix4Pass(data + 12);
// coalesced global writes
globalStores(data, out, 64);
}
```





# Portland Group Accelerator Directives

- Proprietary directives in Fortran and C
- Similar to OpenMP in structure
- If the compiler doesn't understand these directives, it ignores them, so the same code can work with an accelerator or without, and with the PGI compilers or other compilers.
- In principle, this will be able to work on a variety of accelerators, but the first instance is NVIDIA; PGI recently announced a deal with AMD/ATI.
- The directives tell the compiler what parts of the code happen in the accelerator; the rest happens in the regular hardware.





# PGI Accelerator Example

```
!$acc region
  do k = 1,n1
    do i = 1,n3
      c(i,k) = 0.0
      do j = 1,n2
        c(i,k) = c(i,k) +
&          a(i,j) * b(j,k)
      enddo
    enddo
  enddo
!$acc end region
```

<http://www.pgroup.com/resources/accel.htm>



# Digging Deeper: CUDA on NVIDIA





# NVIDIA Tesla

- NVIDIA now offers a GPU platform named Tesla.
- It consists essentially of their highest end graphics card, minus the video out connector.



[http://images.nvidia.com/products/geforce\\_gtx\\_480/geforce\\_gtx\\_480\\_3qtr\\_low.png](http://images.nvidia.com/products/geforce_gtx_480/geforce_gtx_480_3qtr_low.png)



[http://images.nvidia.com/products/tesla\\_C2050\\_C2070/Tesla\\_C2050\\_C2070\\_3qtr\\_low\\_new.png](http://images.nvidia.com/products/tesla_C2050_C2070/Tesla_C2050_C2070_3qtr_low_new.png)



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# NVIDIA Tesla C2050 Card Specs

- 448 GPU cores
- 1.15 GHz
- Single precision floating point performance:  
1030.4 GFLOPs (2 single precision flops per clock per core)
- Double precision floating point performance:  
515.2 GFLOPs (1 double precision flop per clock per core)
- Internal RAM: 3 GB DDR5
- Internal RAM speed: 144 GB/sec (compared 21-25 GB/sec for regular RAM)
- Has to be plugged into a PCIe slot (at most 8 GB/sec)





# NVIDIA Tesla S2050 Server Specs

- 4 C2050 cards inside a 1U server (looks like a Sooner node)
- 1.15 GHz
- Single Precision (SP) floating point performance: 4121.6 GFLOPs
- Double Precision (DP) floating point performance: 2060.8 GFLOPs
- Internal RAM: 12 GB total (3 GB per GPU card)
- Internal RAM speed: 576 GB/sec aggregate
- Has to be plugged into two PCIe slots (at most 16 GB/sec)







# Compare x86 vs S2050

Let's compare the best dual socket x86 server today vs S2050.

	<b>Dual socket, AMD 2.3 GHz 12-core</b>	<b>NVIDIA Tesla S2050</b>
<b>Peak DP FLOPs</b>	220.8 GFLOPs DP	2060.8 GFLOPs DP (9.3x)
<b>Peak SP FLOPs</b>	441.6 GFLOPs SP	4121.6 GFLOPs SP (9.3x)
<b>Peak RAM BW</b>	25 GB/sec	576 GB/sec (23x)
<b>Peak PCIe BW</b>	N/A	16 GB/sec
<b>Needs x86 server to attach to?</b>	No	Yes
<b>Power/Heat</b>	~450 W	~900 W + ~400 W (~2.9x)
<b>Code portable?</b>	Yes	No (CUDA) Yes (PGI, OpenCL)





# Compare x86 vs S2050

Here are some interesting measures:

	Dual socket, AMD 2.3 GHz 12-core	NVIDIA Tesla S2050
<b>DP GFLOPs/Watt</b>	~0.5 GFLOPs/Watt	~1.6 GFLOPs/Watt (~3x)
<b>SP GFLOPs/Watt</b>	~1 GFLOPs/Watt	~3.2 GFLOPs/Watt (~3x)
<b>DP GFLOPs/sq ft</b>	~590 GFLOPs/sq ft	~2750 GFLOPs/sq ft (4.7x)
<b>SP GFLOPs/sq ft</b>	~1180 GFLOPs/sq ft	~5500 GFLOPs/sq ft (4.7x)
<b>Racks per PFLOP DP</b>	142 racks/PFLOP DP	32 racks/PFLOP DP (23%)
<b>Racks per PFLOP SP</b>	71 racks/PFLOP SP	16 racks/PFLOP SP (23%)

OU's Sooner is 34.5 TFLOPs DP, which is just over **1 rack** of S2050.





# What Are the Downsides?

- You have to rewrite your code into CUDA or OpenCL or PGI accelerator directives.
  - CUDA: Proprietary, but maybe portable soon
  - OpenCL: portable but cumbersome
  - PGI accelerator directives: not clear whether you can have most of the code live inside the GPUs.
- BUT: Many groups are coming out with GPGPU code development tools that may help a lot, such as:
  - Fortran-to-CUDA-C converter (NCAR)
  - CUDA C automatic optimizer (memory, threading etc)
  - OpenMP-to-CUDA converter
  - CUDA-to-x86 converter (CUDA code on non-CUDA system)





# Programming for Performance

The biggest single performance bottleneck on GPU cards today is the PCIe slot:

- PCIe 2.0 x16: 8 GB/sec
- 1600 MHz Front Side Bus: 25 GB/sec
- GDDR4 GPU card RAM: 144 GB/sec per card

Your goal:

- At startup, move the data from x86 server RAM into GPU RAM.
- Do almost all the work inside the GPU.
- Use the x86 server only for I/O and message passing, to minimize the amount of data moved through the PCIe slot.





# Does CUDA Help?

Example Applications	URL	Speedup
Seismic Database	<a href="http://www.headwave.com">http://www.headwave.com</a>	66x – 100x
Mobile Phone Antenna Simulation	<a href="http://www.accelware.com">http://www.accelware.com</a>	45x
Molecular Dynamics	<a href="http://www.ks.uiuc.edu/Research/vmd">http://www.ks.uiuc.edu/Research/vmd</a>	21x – 100x
Neuron Simulation	<a href="http://www.evolvedmachines.com">http://www.evolvedmachines.com</a>	100x
MRI Processing	<a href="http://bic-test.beckman.uiuc.edu">http://bic-test.beckman.uiuc.edu</a>	245x – 415x
Atmospheric Cloud Simulation	<a href="http://www.cs.clemson.edu/~jesteel/clouds.html">http://www.cs.clemson.edu/~jesteel/clouds.html</a>	50x

[http://www.nvidia.com/object/IO\\_43499.html](http://www.nvidia.com/object/IO_43499.html)



# CUDA

# Thread Hierarchy and Memory Hierarchy

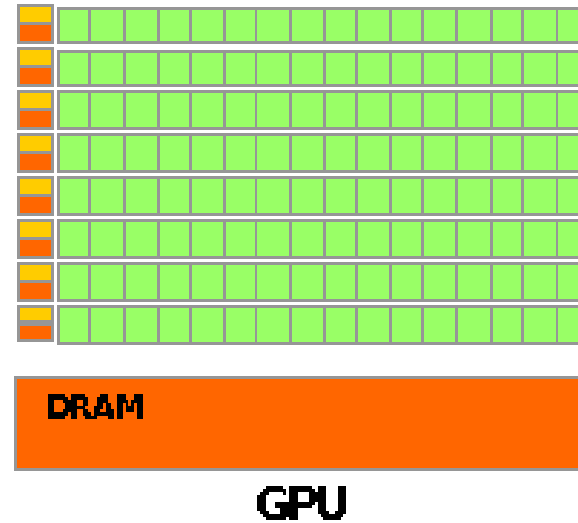
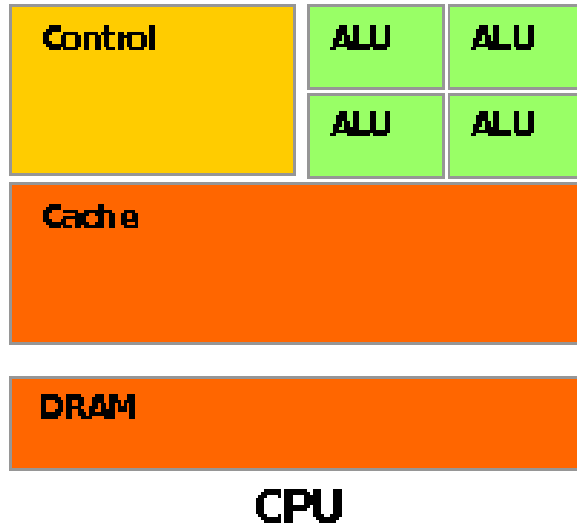


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Some of these slides provided by Paul Gray, University of Northern Iowa



# CPU vs GPU Layout



Source: NVIDIA CUDA Programming Guide





# Buzzword: Kernel

In CUDA, a *kernel* is code (typically a function) that can be run inside the GPU.

Typically, the kernel code operates in lock-step on the stream processors inside the GPU.







# Buzzword: Thread

In CUDA, a *thread* is an execution of a kernel with a given index.

Each thread uses its index to access a specific subset of the elements of a target array, such that the collection of all threads cooperatively processes the entire data set.

So these are very much like threads in the OpenMP or pthreads sense – they even have shared variables and private variables.





# Buzzword: Block

In CUDA, a ***block*** is a group of threads.

- Just like OpenMP threads, these could execute concurrently or independently, and in no particular order.
- Threads can be coordinated somewhat, using the `_syncthreads()` function as a barrier, making all threads stop at a certain point in the kernel before moving on en mass. (This is like what happens at the end of an OpenMP loop.)





# Buzzword: Grid

In CUDA, a ***grid*** is a group of (thread) blocks, with no synchronization at all among the blocks.



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# NVIDIA GPU Hierarchy

- Grids map to GPUs
- Blocks map to the MultiProcessors (MP)
  - Blocks are never split across MPs, but an MP can have multiple blocks
- Threads map to Stream Processors (SP)
- Warps are groups of (32) threads that execute simultaneously

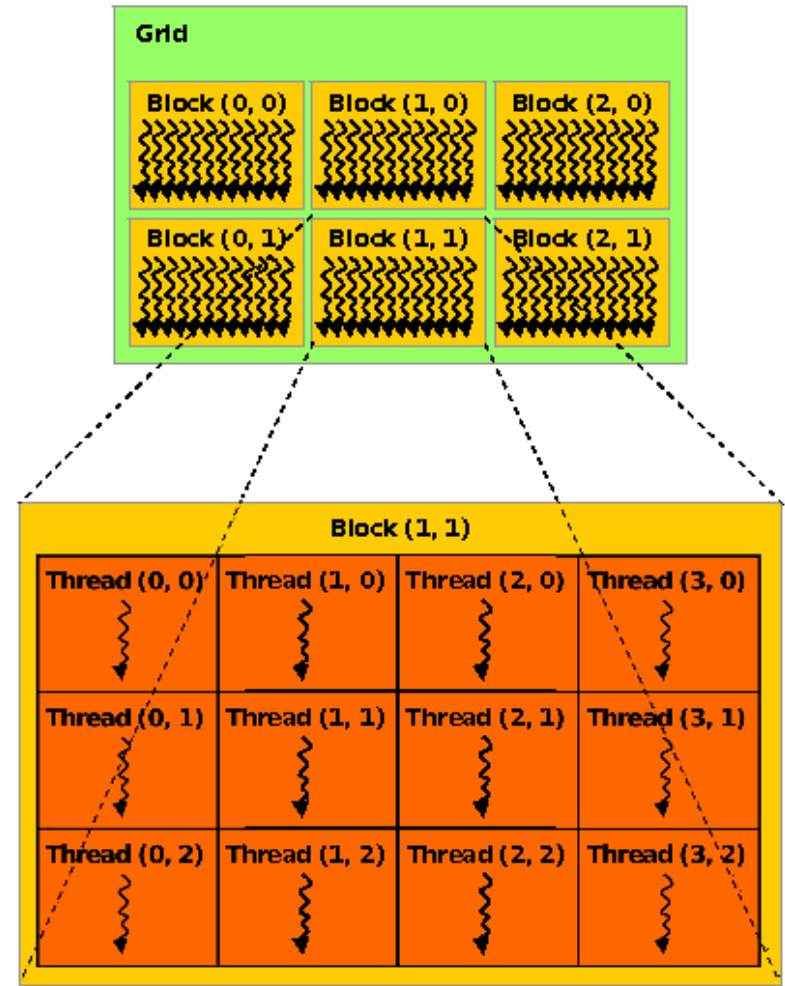


Image Source:  
NVIDIA CUDA Programming Guide



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# CUDA Built-in Variables

- **blockIdx.x**, **blockIdx.y**, **blockIdx.z** are built-in variables that returns the block ID in the x-axis, y-axis and z-axis of the block that is executing the given block of code.
- **threadIdx.x**, **threadIdx.y**, **threadIdx.z** are built-in variables that return the thread ID in the x-axis, y-axis and z-axis of the thread that is being executed by this stream processor in this particular block.

So, you can express your collection of blocks, and your collection of threads within a block, as a 1D array, a 2D array or a 3D array.

These can be helpful when thinking of your data as 2D or 3D.





## \_\_global\_\_ Keyword

In CUDA, if a function is declared with the `__global__` keyword, that means that it's intended to be executed inside the GPU.

In CUDA, the term for the GPU is *device*, and the term for the x86 server is *host*.

So, a kernel runs on a device, while the main function and so on run on the host.

Note that a host can play host to multiple devices; for example, an S2050 server contains 4 C2050 GPU cards, and if a single host has two PCIe slots, then both of the PCIe plugs of the S2050 can be plugged into that same host.





# Copying Data from Host to Device

If data need to move from the host (where presumably the data are initially input or generated), then a copy has to exist in both places.

Typically, what's copied are arrays, though of course you can also copy a scalar (the address of which is treated as an array of length 1).

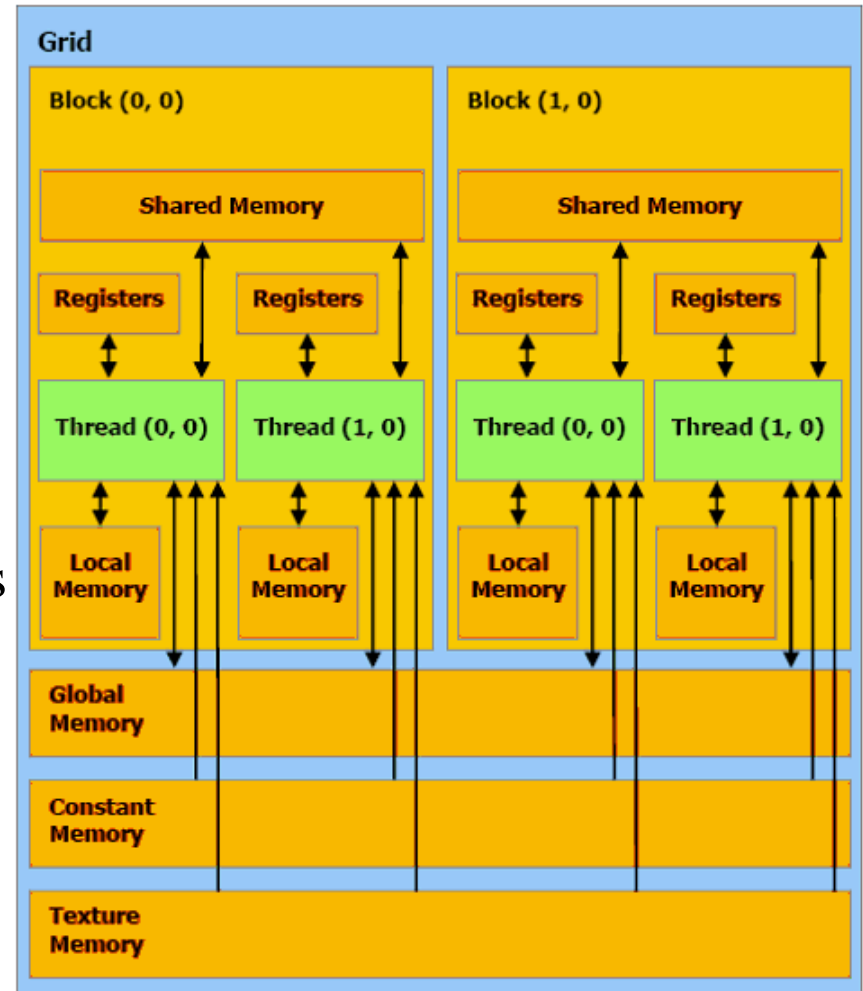




# CUDA Memory Hierarchy #1

CUDA has a hierarchy of several kinds of memory:

- Host memory (x86 server)
- Device memory (GPU)
  - **Global**: visible to all threads in all blocks – largest, slowest
  - **Shared**: visible to all threads in a particular block – medium size, medium speed
  - **Local**: visible only to a particular thread – smallest, fastest



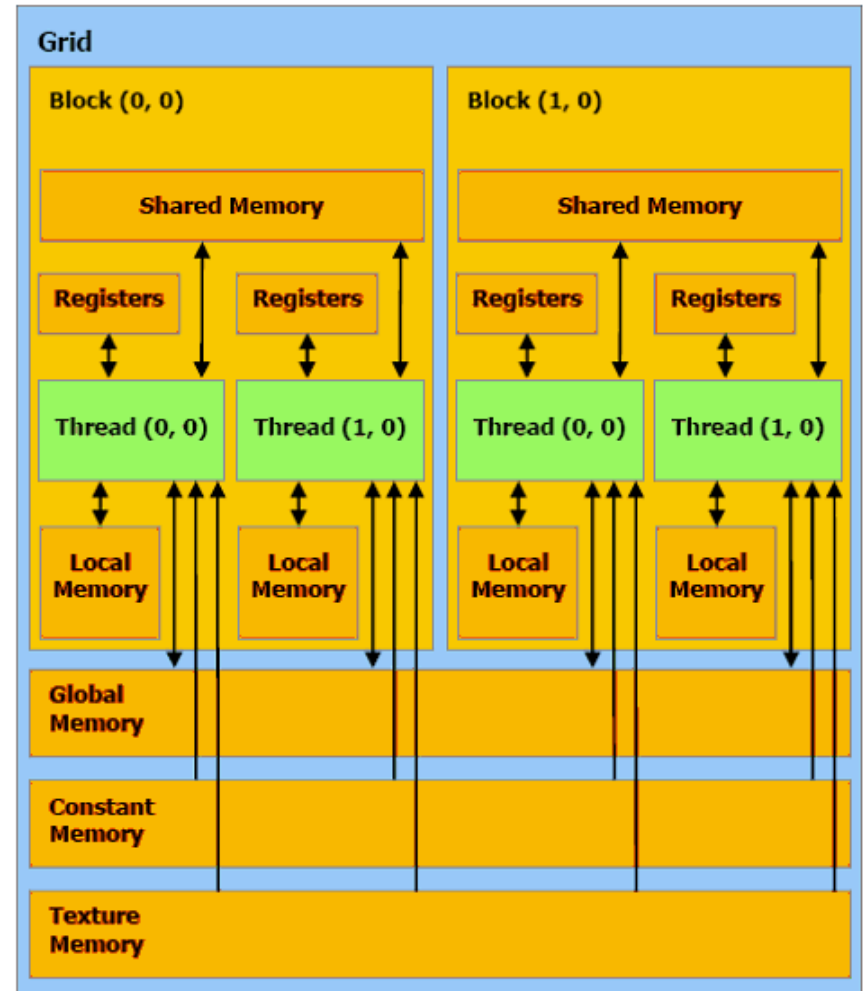




# CUDA Memory Hierarchy #2

CUDA has a hierarchy of several kinds of memory:

- Host memory (x86 server)
- Device memory (GPU)
  - **Constant**: visible to all threads in all blocks; read only
  - **Texture**: visible to all threads in all blocks; read only



# CUDA Example: Matrix-Matrix Multiply



[http://developer.download.nvidia.com/compute/cuda/sdk/  
website/Linear\\_Algebra.html#matrixMul](http://developer.download.nvidia.com/compute/cuda/sdk/website/Linear_Algebra.html#matrixMul)



# Matrix-Matrix Multiply Main Part 1

```
float* host_A;
float* host_B;
float* host_C;
float* device_A;
float* device_B;
float* device_C;

host_A = (float*) malloc(mem_size_A);
host_B = (float*) malloc(mem_size_B);
host_C = (float*) malloc(mem_size_C);

cudaMalloc((void**) &device_A, mem_size_A);
cudaMalloc((void**) &device_B, mem_size_B);
cudaMalloc((void**) &device_C, mem_size_C);

// Set up the initial values of A and B here.

// Henry says: I've oversimplified this a bit from
// the original example code.
```





# Matrix-Matrix Multiply Main Part 2

```
// copy host memory to device
cudaMemcpy(device_A, host_A, mem_size_A,
           cudaMemcpyHostToDevice);
cudaMemcpy(device_B, host_B, mem_size_B,
           cudaMemcpyHostToDevice);

// setup execution parameters
dim3 threads(BLOCK_SIZE, BLOCK_SIZE);
dim3 grid(WC / threads.x, HC / threads.y);

// execute the kernel
matrixMul<<< grid, threads >>>(device_C,
                                device_A, device_B, WA, WB);

// copy result from device to host
cudaMemcpy(host_C, device_C, mem_size_C,
           cudaMemcpyDeviceToHost);
```





# Matrix Matrix Multiply Kernel Part 1

```
__global__ void matrixMul( float* C, float* A, float* B, int wA, int wB)
{
    // Block index
    int bx = blockIdx.x;
    int by = blockIdx.y;

    // Thread index
    int tx = threadIdx.x;
    int ty = threadIdx.y;

    // Index of the first sub-matrix of A processed by the block
    int aBegin = wA * BLOCK_SIZE * by;

    // Index of the last sub-matrix of A processed by the block
    int aEnd   = aBegin + wA - 1;

    // Step size used to iterate through the sub-matrices of A
    int aStep  = BLOCK_SIZE;

    // Index of the first sub-matrix of B processed by the block
    int bBegin = BLOCK_SIZE * bx;

    // Step size used to iterate through the sub-matrices of B
    int bStep  = BLOCK_SIZE * wB;

    // Csub is used to store the element of the block sub-matrix
    // that is computed by the thread
    float Csub = 0;
```





# Matrix Matrix Multiply Kernel Part 2

```
// Loop over all the sub-matrices of A and B
// required to compute the block sub-matrix
for (int a = aBegin, b = bBegin;
     a <= aEnd;
     a += aStep, b += bStep) {

    // Declaration of the shared memory array As used to
    // store the sub-matrix of A
    __shared__ float As[BLOCK_SIZE][BLOCK_SIZE];

    // Declaration of the shared memory array Bs used to
    // store the sub-matrix of B
    __shared__ float Bs[BLOCK_SIZE][BLOCK_SIZE];

    // Load the matrices from device memory
    // to shared memory; each thread loads
    // one element of each matrix
    AS(ty, tx) = A[a + wA * ty + tx];
    BS(ty, tx) = B[b + wB * ty + tx];

    // Synchronize to make sure the matrices are loaded
    __syncthreads();
}
```





# Matrix Matrix Multiply Kernel Part 3

```
// Multiply the two matrices together;
// each thread computes one element
// of the block sub-matrix
for (int k = 0; k < BLOCK_SIZE; ++k)
    Csub += AS(ty, k) * BS(k, tx);

// Synchronize to make sure that the preceding
// computation is done before loading two new
// sub-matrices of A and B in the next iteration
__syncthreads();
}
```

```
// Write the block sub-matrix to device memory;
// each thread writes one element
int c = wB * BLOCK_SIZE * by + BLOCK_SIZE * bx;
C[c + wB * ty + tx] = Csub;
}
```





# Would We Really Do It This Way?

We wouldn't really do matrix-matrix multiply this way.

NVIDIA has developed a CUDA implementation of the BLAS libraries, which include a highly tuned matrix-matrix multiply routine.

(We'll learn about BLAS next time.)

There's also a CUDA FFT library, if your code needs Fast Fourier Transforms.







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**Thanks for your  
attention!**



**Questions?**

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