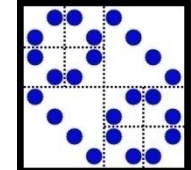
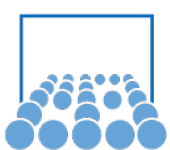


GPU

CUDA Programming (NVIDIA)

*Compute **U**nified **D**evice **A**rchitecture:*

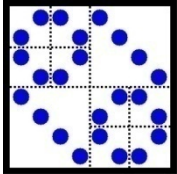
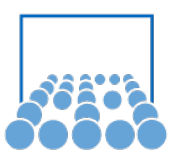
- C programming on GPUs
- Requires no knowledge of graphics APIs or GPU programming
- Access to naive instructions and memory
- Easy to get started
- Stable, free, documented, supported
- Windows and Linux



GPU

GPU is viewed as a compute device operating as a coprocessor to the main CPU (host)

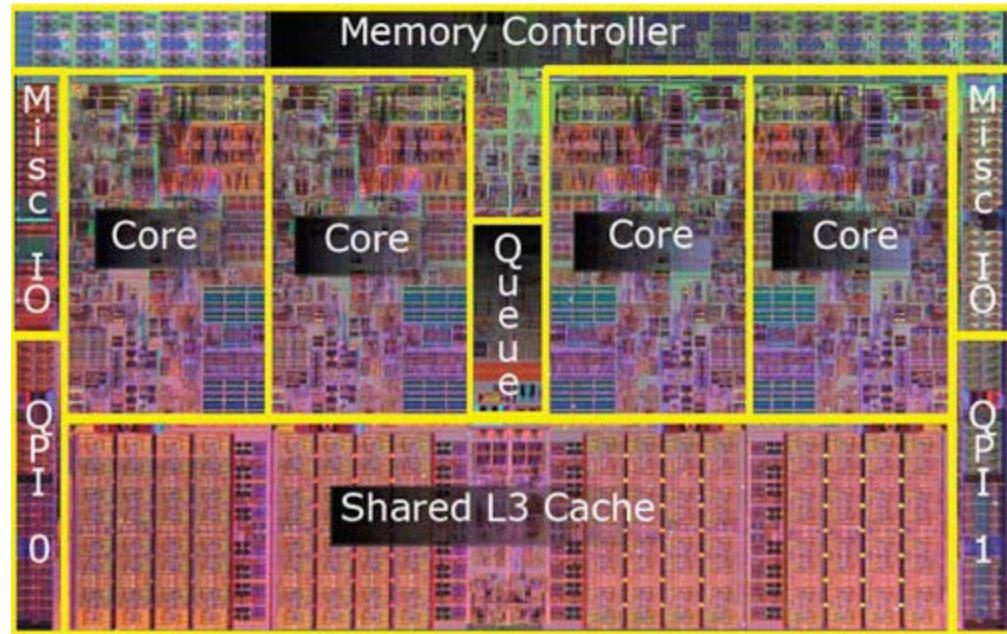
- For data parallel, cost intensive functions or functions that are executed many times, but on different data (loops)
- A function compiled for the device is called a kernel
- The kernel is executed on the device as many different threads
- Host (CPU) and device (GPU) manage their own memory
- Data can be copied between them

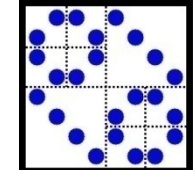


3rd Level of Parallelism



MPI
OpenMP





Grid of Thread Blocks

Computational Grid

Block (0, 1)	Block (1, 1)	Block (2, 1)
Block (0, 0)	Block (1, 0)	Block (2, 0)

Block (1,0)

Thread (0, 3)	Thread (1, 3)	Thread (2, 3)	Thread (3, 3)
Thread (0, 2)	Thread (1, 2)	Thread (2, 2)	Thread (3, 2)
Thread (0, 1)	Thread (1, 1)	Thread (2, 1)	Thread (3, 1)
Thread (0, 0)	Thread (1, 0)	Thread (2, 0)	Thread (3, 0)

The computational grid consists of a grid of thread blocks

Each thread executes the kernel

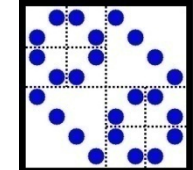
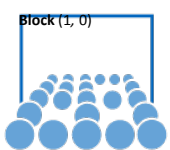
The application specifies the grid and the block dimensions

Grid layouts can be 1D, 2D, 3D

Maximal sizes are determined by GPU memory and kernel complexity

Each block has unique block ID

Each thread has unique thread ID (within its block)



Example: Matrix Addition

CPU Program

```
void add_matrix
( float* a, float* b, float* c, int N ) {
int index;
for ( int i = 0; i < N; ++i )
for ( int j = 0; j < N; ++j ) {
index = i + j*N;
c[index] = a[index] + b[index];
}
}

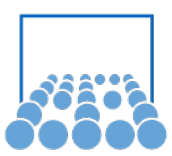
int main() {
add_matrix( a, b, c, N );
}
```

CUDA Program

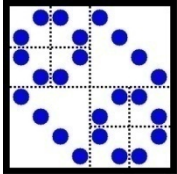
```
__global__ add_matrix
( float* a, float* b, float* c, int N ) {
int i = blockIdx.x * blockDim.x + threadIdx.x;
int j = blockIdx.y * blockDim.y + threadIdx.y;
int index = i + j*N;
if ( i < N && j < N )
c[index] = a[index] + b[index];
}

int main() {
dim3 dimBlock( blocksize, blocksize );
dim3 dimGrid( N/dimBlock.x, N/dimBlock.y );
add_matrix<<<dimGrid, dimBlock>>>( a, b, c, N );
}
```

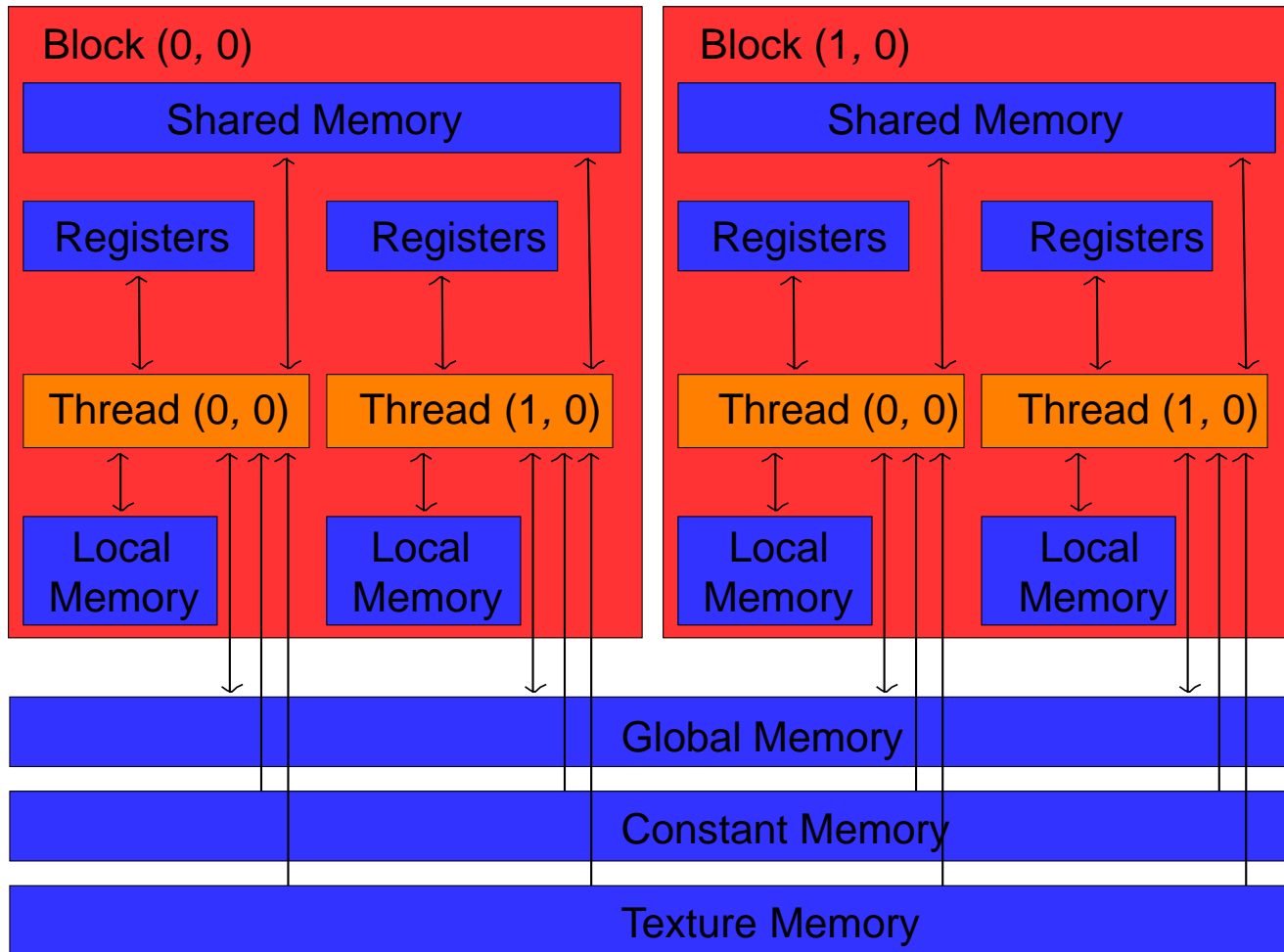
The nested for loops are replaced with an implicit grid

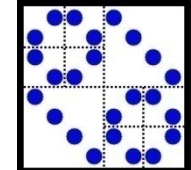
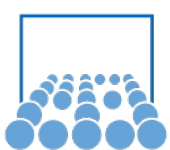


Memory Model



CUDA exposes all the different types of memory on the GPU





Memory Model II

Overview			
Registers	Per thread	Read-Write	
Local memory	Per thread	Read-Write	
Shared memory	Per block	Read-Write	For sharing data within a block
Global memory	Per grid	Read-Write	Not cached
Constant memory	Per grid	Read-only	Cached
Texture memory	Per grid	Read-only	Spatially cached

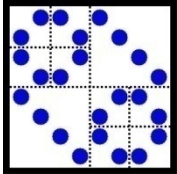
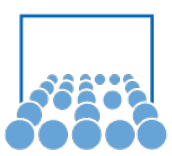
Explicit GPU memory allocation and deallocation via
`cudaMalloc()` and `cudaFree()`

Access to GPU memory via pointers

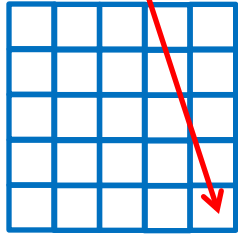
Copying between CPU and GPU memory:

A slow operation: minimize this!

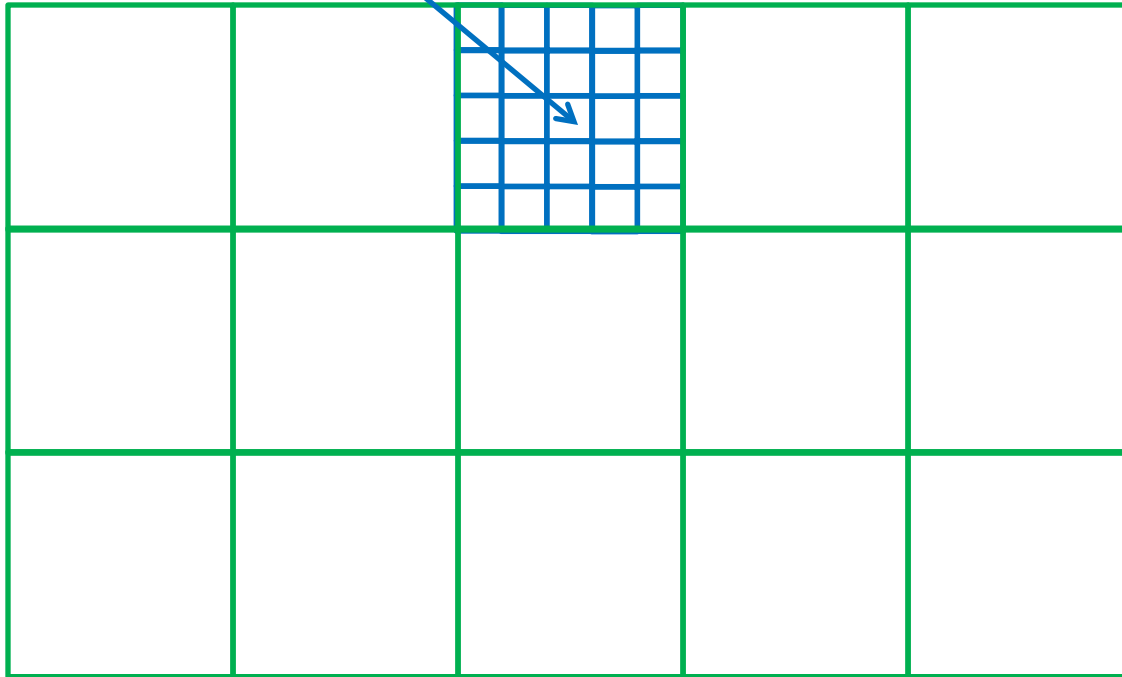
Multiple levels of parallelism:



Thread



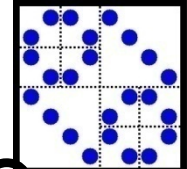
Thread block



Grid of thread blocks

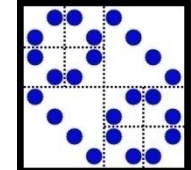
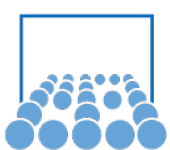
- Thread block:
 - Up to 512 threads per block
 - Communicate via shared memory
 - Threads guaranteed to be resident
 - `threadIdx, blockIdx`
 - `__syncthreads()`

- Grid of thread blocks:
 - `f<<<N, T>>>(a, b, c)`
 - Communication via global memory



CUDA Programming Language

- Minimal C extensions
- A runtime library
 - A host (CPU) component to control and access GPU(s)
 - A device component
 - A common component:
 - Built-in vector types, C standard library subset
- Function type qualifiers
 - Specify where to call and execute a function
 - `__device__` , `__global__` , `__host__`
- Variable type qualifiers
 - `__device__` , `__constant__` , `__shared__`
- Kernel execution directive
 - `foo<<<GridDim, BlockDim>>>(…)`
- Built-in variables for grid/block size and block/thread indices



CUDA Compiler

Source files must be compiled with the CUDA compiler `nvcc`

CUDA kernels are stored in files ending with `.cu`

NVCC uses the host compiler to compile CPU code

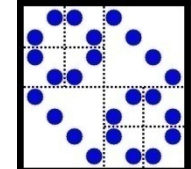
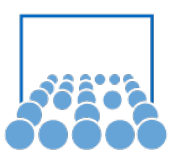
NVCC automatically handles `#include`'s and linking

Write kernels+CPU caller in `.cu` files

Compile with `nvcc`

Store signature of CPU called in header file

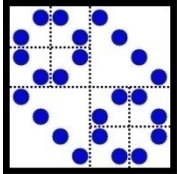
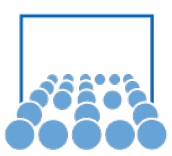
`#include` header file in C(++) sources



GPU Runtime Component

Only available on the GPU:

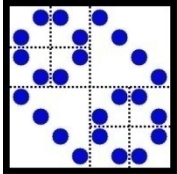
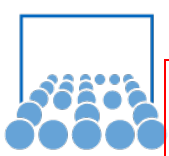
- Less accurate, faster math functions `__sin(x)`
- `syncthreads()` : wait until all threads in the block have reached this point
- Type conversion functions, with rounding mode
- Type casting functions
- Texture functions
- Atomic functions:
Guarantees that operation (like add) is performed on a variable without interference from other threads



CPU Runtime Component

Only available on the CPU:

- Device management:
 Get device properties, multi-GPU control, ...
- Memory management:
 `cudaMalloc()`, `cudaMemcpy()`, `cudaFree()`,...
- Texture management
- Asynchronous concurrent execution
- Built-in vector types: i.e. `float1`, `float2`, `int3`, `ushort4`, ...
- Mathematical functions: standard `math.h` on CPU,



```
const int N = 1024; const int blocksize = 16;
```

```
__global__
void add_matrix( float* a, float *b, float *c, int N )
{ int i = blockIdx.x * blockDim.x + threadIdx.x;
  int j = blockIdx.y * blockDim.y + threadIdx.y;
  int index = i + j*N;
  if ( i < N && j < N )
    c[index] = a[index] + b[index]; }
```

COMPUTE KERNEL

Store in source file
(i.e. MatrixAdd.cu)
Compile with
nvcc MatrixAdd.cu
Run

```
int main() {
float *a = new float[N*N]; float *b = new float[N*N]; float *c = new float[N*N];
for ( int i = 0; i < N*N; ++i ) {
a[i] = 1.0f; b[i] = 3.5f; }
```

CPU MEMORY ALLOCATION

```
float *ad, *bd, *cd;
const int size = N*N*sizeof(float); cudaMalloc( (void**)&ad, size );
cudaMalloc( (void**)&bd, size ); cudaMalloc( (void**)&cd, size );
```

GPU MEMORY ALLOCATION

```
cudaMemcpy( ad, a, size, cudaMemcpyHostToDevice );
cudaMemcpy( bd, b, size, cudaMemcpyHostToDevice );
```

COPY DATA TO GPU

```
dim3 dimBlock( blocksize, blocksize );
dim3 dimGrid( N/dimBlock.x, N/dimBlock.y );
add_matrix<<<dimGrid, dimBlock>>>( ad, bd, cd, N );
```

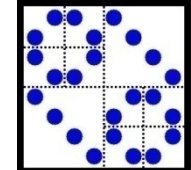
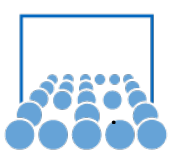
EXECUTE KERNEL

```
cudaMemcpy( c, cd, size, cudaMemcpyDeviceToHost );
```

COPY RESULTS CPU

```
cudaFree( ad ); cudaFree( bd ); cudaFree( cd );
delete[] a; delete[] b; delete[] c;
return EXIT_SUCCESS; }
```

CLEAN UP, RETURN



Execution Model

A GPU consists of N multiprocessors (MP)

Each MP has M scalar processors (SP)

Each MP processes batches of blocks

A block is processed by only one MP

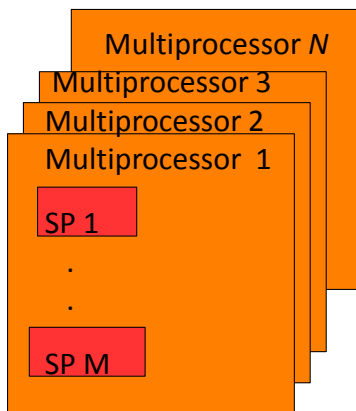
Each block is split into SIMD groups of threads called warps

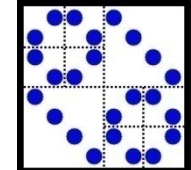
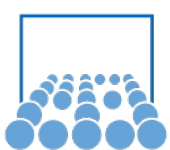
A warp is executed physically in parallel

A scheduler switches between warps

A warp contains threads of consecutive, increasing thread ID

The warp size is 32 threads today





Costs of Operations

4 clock cycles: Floating point: add, multiply, fused mult.-add
Integer add, bitwise op., compare, min, max

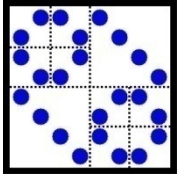
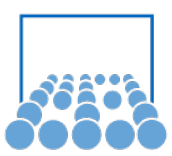
16 clock cycles: reciprocal, square root, __log, 32-bit int mult

32 clock cycles: __sin(x) , __cos(x) , __exp(x)

36 clock cycles: Floating point division

Particularly expensive: Integer division, modulo

Double precision will perform at half the speed



Right kind of memory

Constant memory:

Quite small, $\approx 20K$

As fast as register access if all threads in a warp
access the same location

Texture memory:

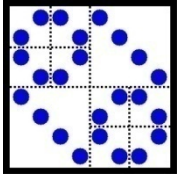
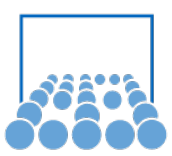
Spatially cached

Optimized for 2D locality

Neighbouring threads should read neighbouring addresses

No need to think about coalescing

Constraint: These memories can only be updated from CPU



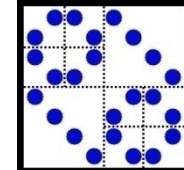
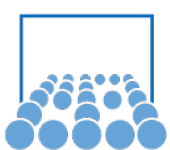
Accessing global memory

4 cycles to issue on memory fetch
but 400-600 cycles of latency!

Likely to be a performance bottleneck

Order of magnitude speedups possible:
Coalesce memory access

Use shared memory to reorder non-coalesced addressing



Coalescing (Assembling)

For best performance, global memory access should be coalesced

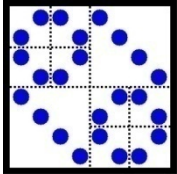
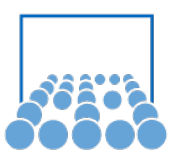
A memory access coordinated within a warp

A contiguous, aligned region of global memory

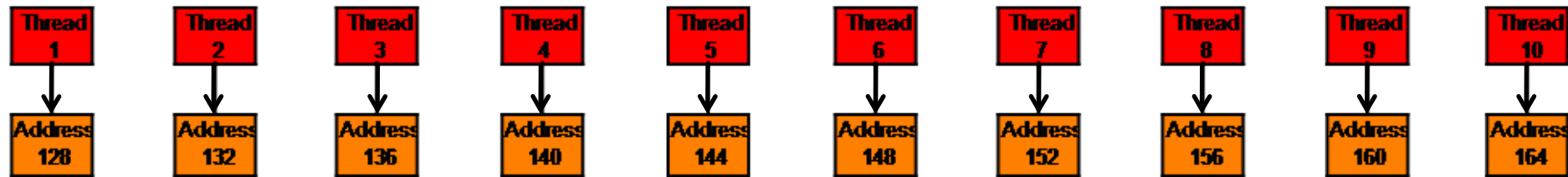
- 128 bytes – each thread reads a float or int
- 256 bytes – each thread reads a float2 or int2
- 512 bytes – each thread reads a float4 or int4
- float3s are not aligned

Warp base address (WBA) must be a multiple of $16 * \text{sizeof}(\text{type})$

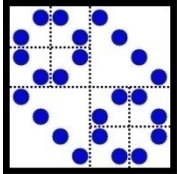
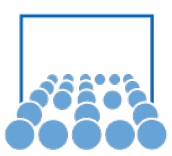
The k-th thread should access the element at $\text{WBA} + k$
These restrictions apply to both reading and writing



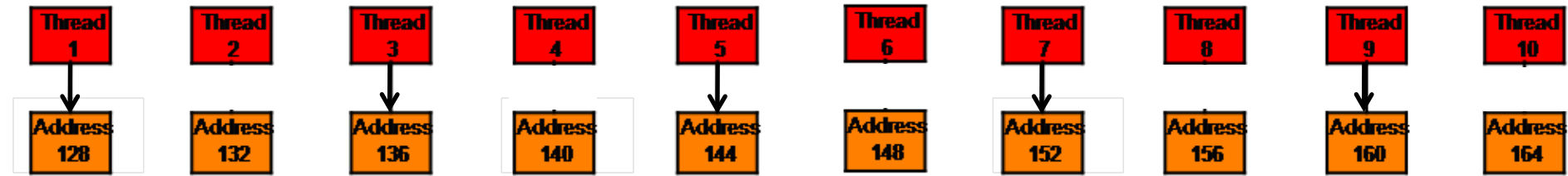
Coalesced memory access



Coalesced memory access:
Thread k accesses $WBA + k$

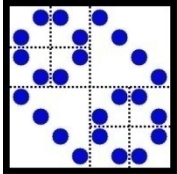
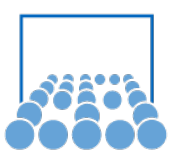


Coalesced memory access

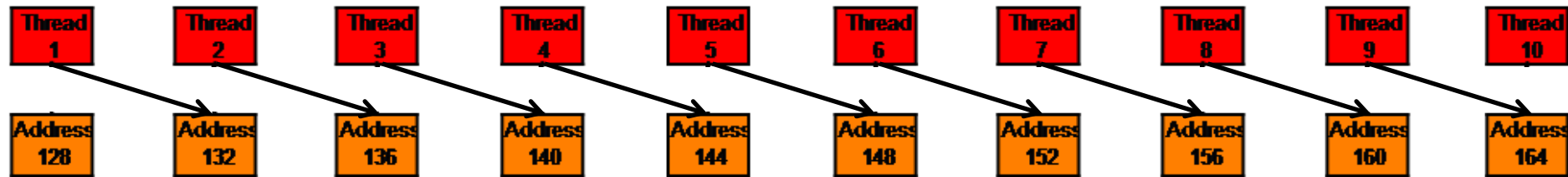


Coalesced memory access:
Thread k accesses $WBA + k$

Not all threads have to participate

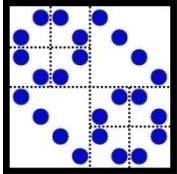
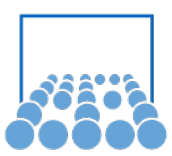


Coalesced memory access

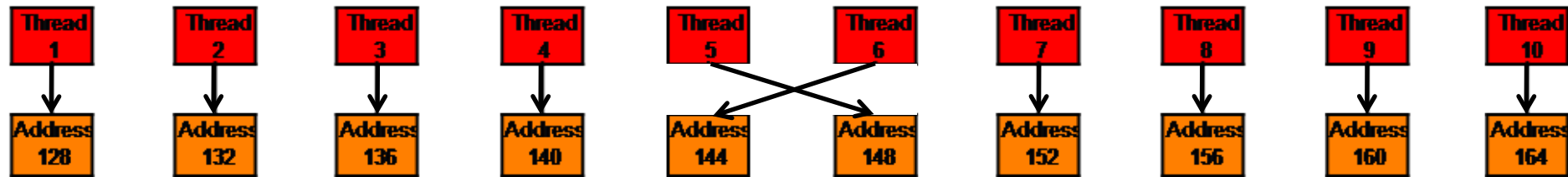


Non-coalesced memory access:

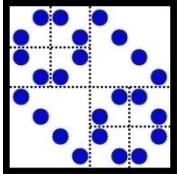
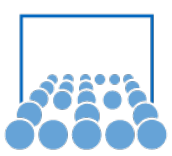
Misaligned starting address



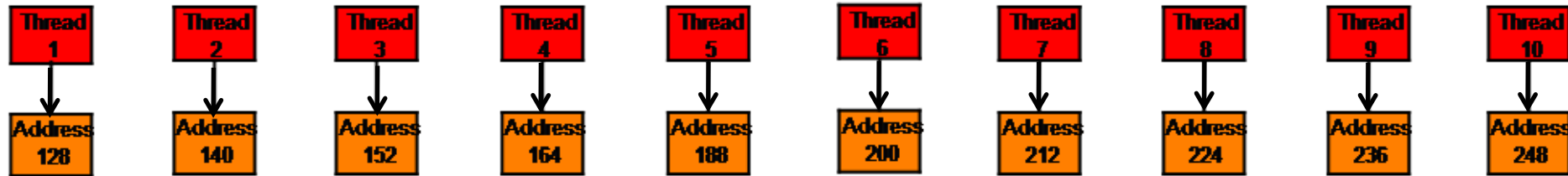
Coalesced memory access



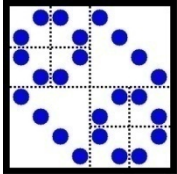
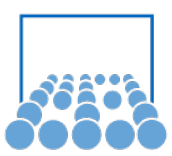
Non-coalesced memory access:
Non-sequential access



Coalesced memory access



Non-coalesced memory access:
Wrong size of type



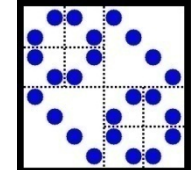
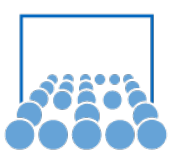
Software fro GPU

CUDA

OpenCL: Vendor-independent industrial standard

DirectCompute: GPU computing from Microsoft

...



OpenCL

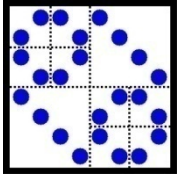
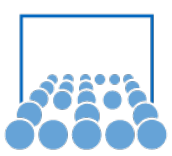
C-based language for GPU kernels + device kernels

Plus low-level device API (application programming interface)

- Same flavor as CUDA
- JIT (just in time) compilation of kernel programs
- Portable - but inevitable optimization required for every platform

Managed by Khronos group (non-profit organization)

- All major vendors participate
- This is the cross-vendor industry-standard



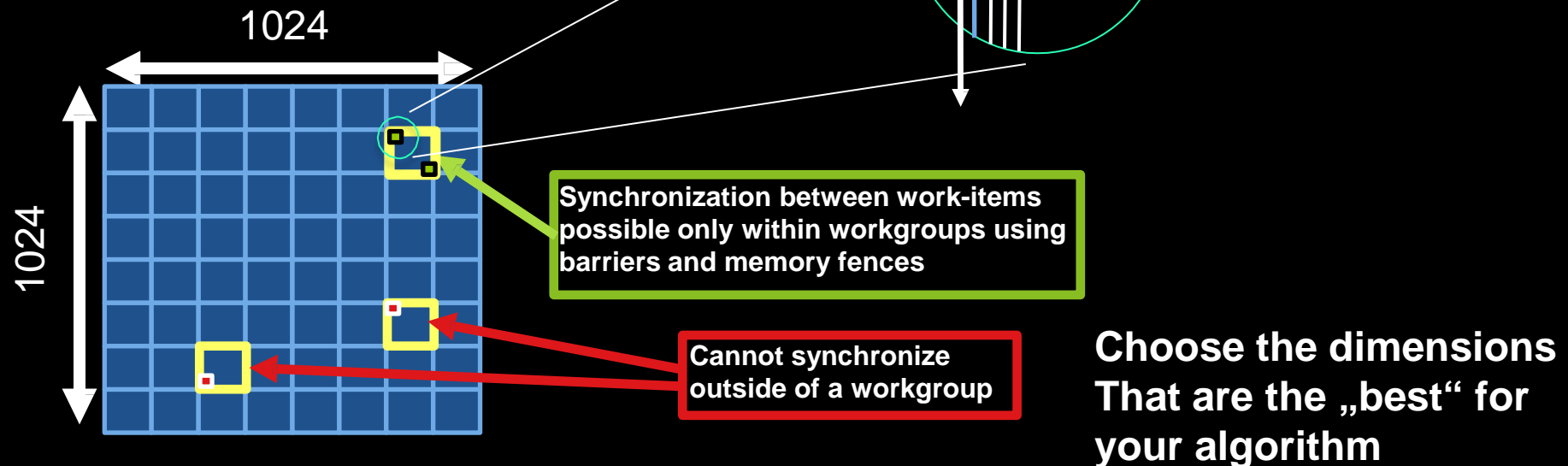
Work Items – N-D Range

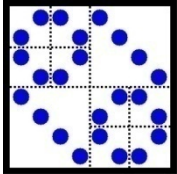
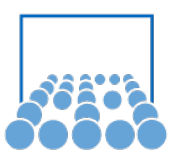
OpenCL execution model:

- define an N-d computation domain
- execute kernel for each point in the domain

Global dimensions: 1024 x 1024

Local dimensions: 128 x 128
work group – execute together





Parallelizing For Loop

Scalar:

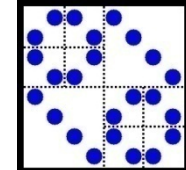
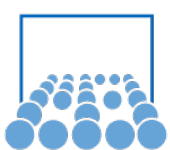
```
void
scalar_mul(const float *a,
           const float *b,
           float *c,
           int n)
{
    int i;
    for (i=0; i<n; i++)
        c[i] = a[i] * b[i];
}
```

Data Parallel:

Kernel executed n times,
once for each work item

```
kernel void
dp_mul(global const float *a,
        global const float *b,
        global float *c)
{
    int id = get_global_id(0);
    c[id] = a[id] * b[id];
} // execute over „n“ work-items
```

Get the index of the work item



OpenCL Objects

Setup:

- Devices (CPU, GPU, Cell,..)
- Contexts (collection of devices)
- Queues (submit work to the device)

Memory:

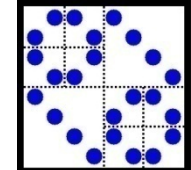
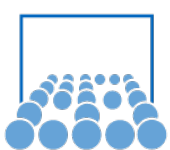
- Buffers (blocks of memory, kernels can access however they like)
- Images (2D or 3D formatted images)

Execution:

- Programs (collections of kernels)
- Kernels

Synchronisation:

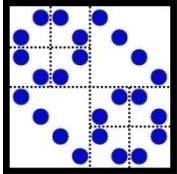
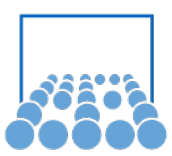
- Events



DirectCompute

GPGPU under Windows

- Microsoft API / Windows standard for all GPU vendors
- General-purpose GPU computing under Windows
- released with DirectX11 / Windows 7
- Supports all CUDA-enabled devices and ATI GPUs
- Low-level API for device management and launching of kernels
- Defines HLSL-based language for compute shaders (High Level Shading Language)



Application-level integration

Matlab:

Parallel Computing Toolbox,
Jacket / AccelerEyes

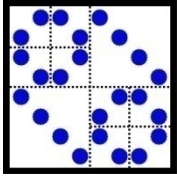
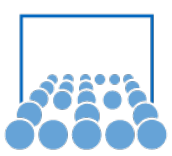
Mathematica

LabView

PetSc

Trilinos

OpenFoam, Ansys,..



Application-specific Libraries

CUsparse: NVIDIA library for sparse matrix vector operations

CUBLAS: NVIDIA library for dense linear algebra

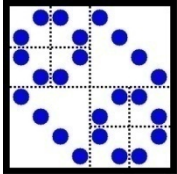
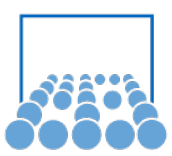
CUFFT: NVIDIA library for Fast Fourier Transforms

CUSP: NVIDIA library for sparse linear algebra

Thrust: A template library for CUDA applications (sort, reduce,..)

MAGMA: LAPACK for HPC on heterogeneous architectures

OpenCurrent: PDE, Gauss-Seidel, multigrid



Application-specific Libraries

ViennaCL: Scientific computing library, C++, OpenCL
BLAS, iterative methods, preconditioners

LAMA: Library for Accelerated math Applications
BLAS, various sparse matrix storage formats

FEAST: Finite Element Analysis and Solution Tools

PARDISO: Fast direct solver for sparse problems
Using BLAS functionality of GPUs