

GPU Computing with CUDA

Lecture 2 - CUDA Memories

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Outline of lecture

- ▶ Recap of Lecture 1
- ▶ Warp scheduling
- ▶ CUDA Memory hierarchy
- ▶ Introduction to the Finite Difference Method

Recap

▶ GPU

- A real alternative in scientific computing
- Massively parallel commodity hardware
 - ▶ Market driven development
 - ▶ Cheap!

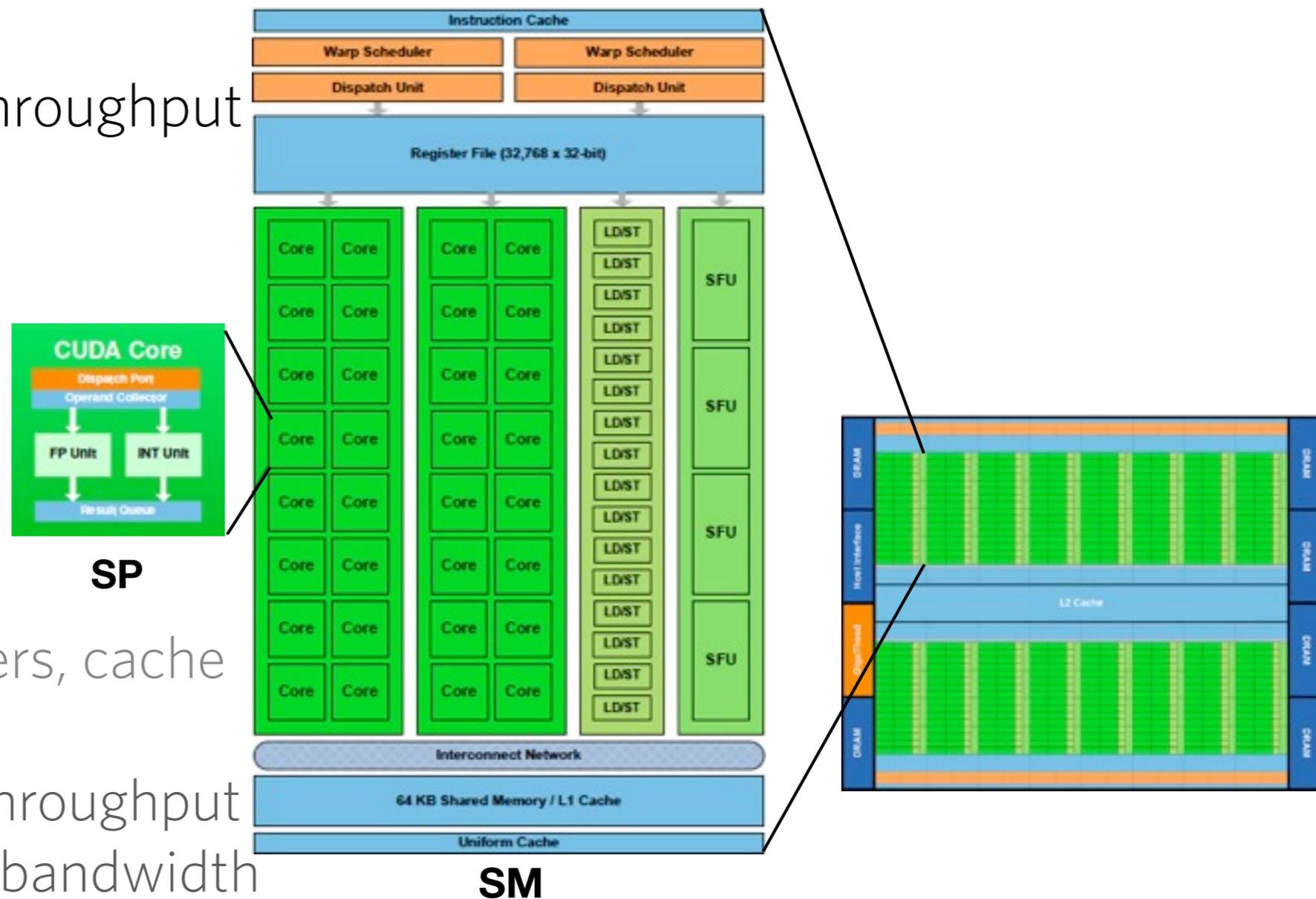
▶ CUDA

- NVIDIA technology capable of programming massively parallel processors with general purpose
- C/C++ with extensions

Recap

▶ GPU chip

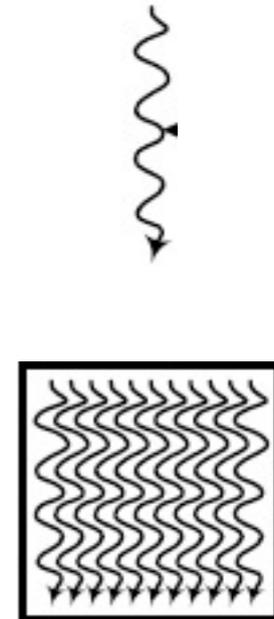
- Designed for high throughput tasks
- Many simple cores
- Fermi:
 - ▶ 16 SMs
 - 32 SPs (cores)
 - Shared, registers, cache
 - SFU, CU
 - ▶ 1TFLOP peak throughput
 - ▶ 144GB/s peak bandwidth



Recap

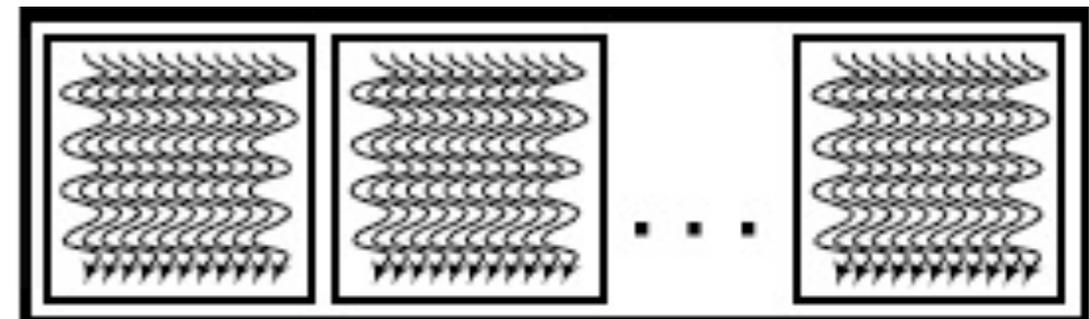
▶ Programming model

- Based on **threads**
- Thread hierarchy: grouped in **thread blocks**
- Threads in a thread block can communicate



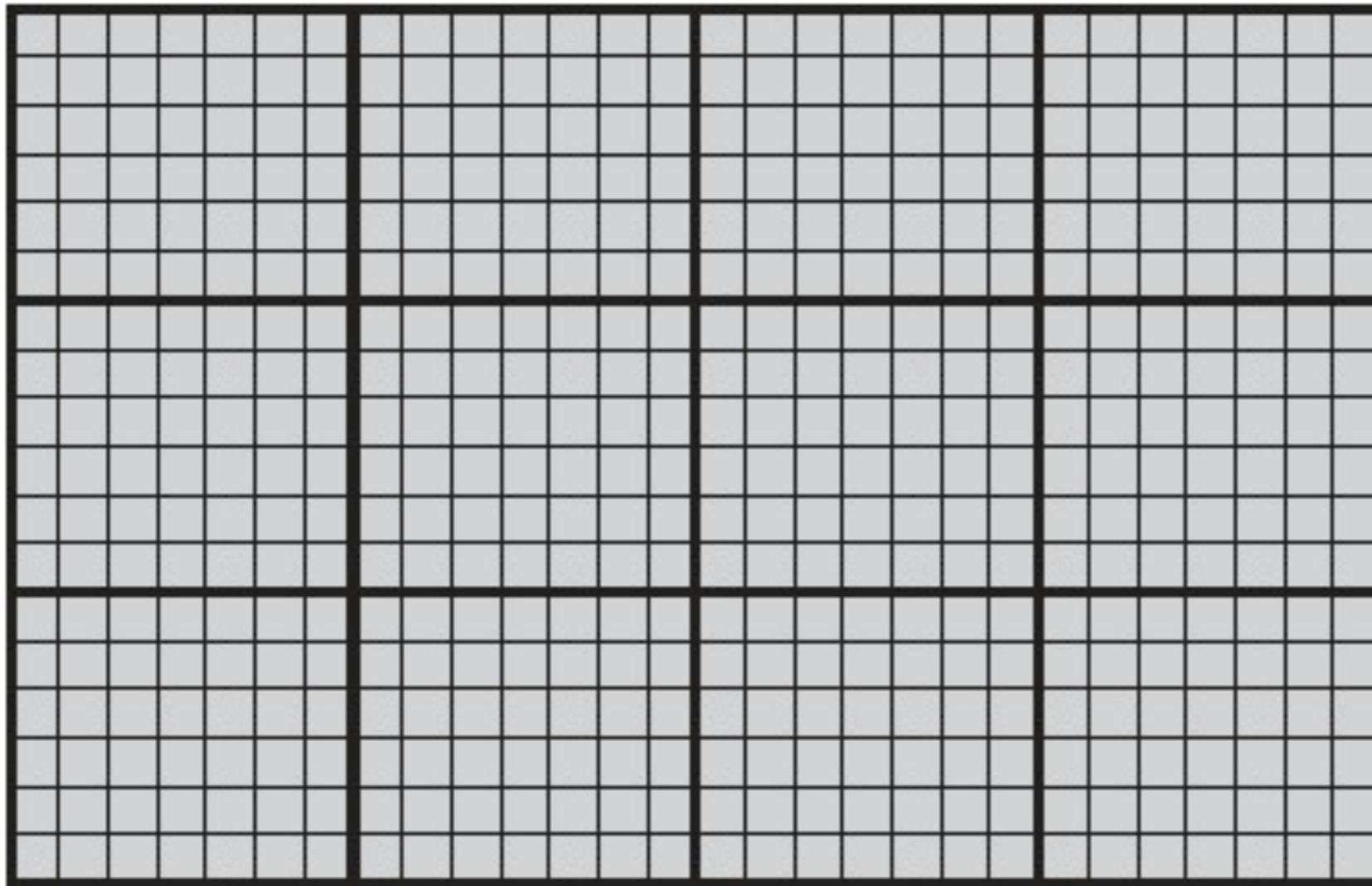
▶ Challenge: parallel thinking

- Data parallelism
- Load balancing
- Conflicts
- Latency hiding



CUDA - Programming model

- ▶ Connecting the hardware and the software



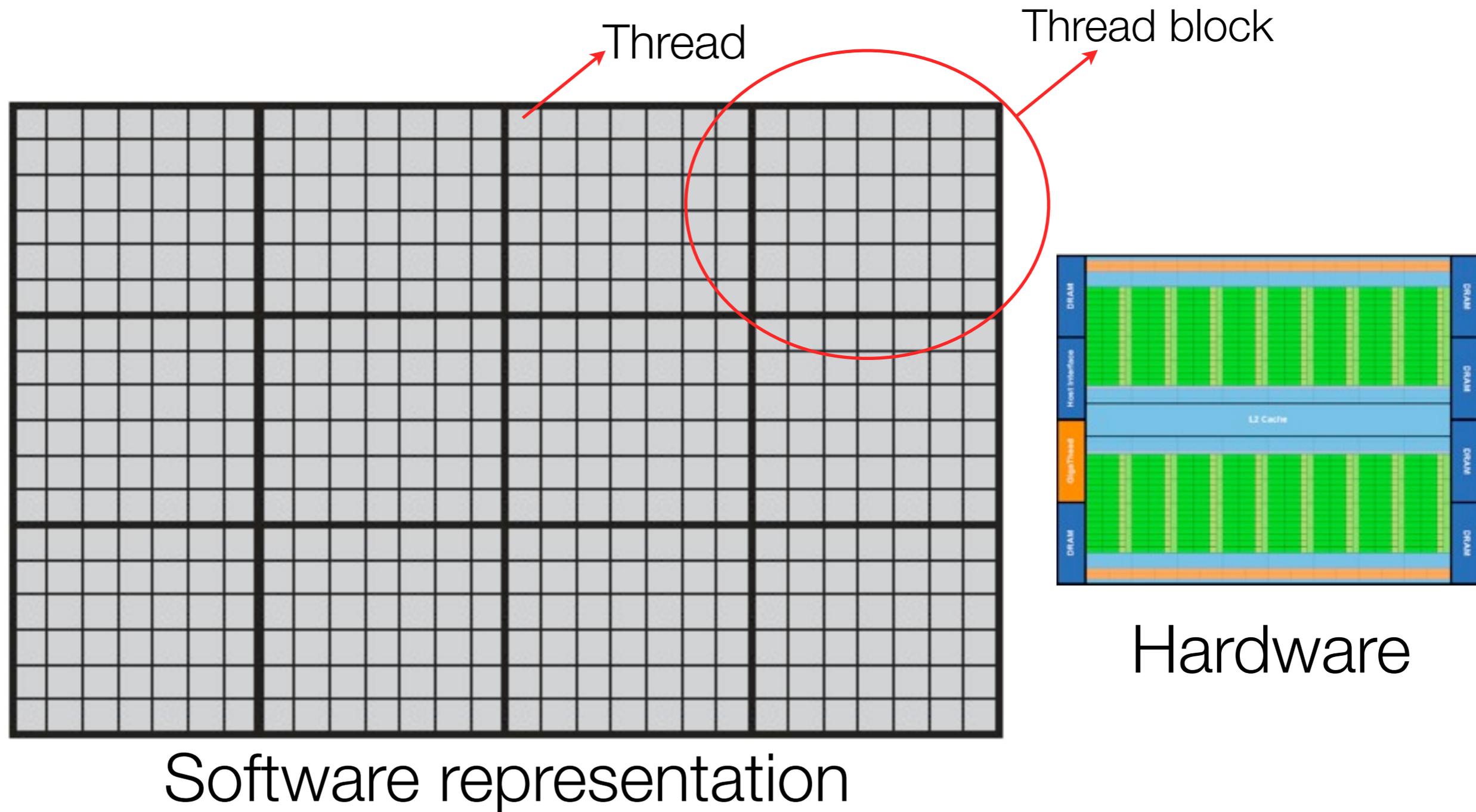
Software representation



Hardware

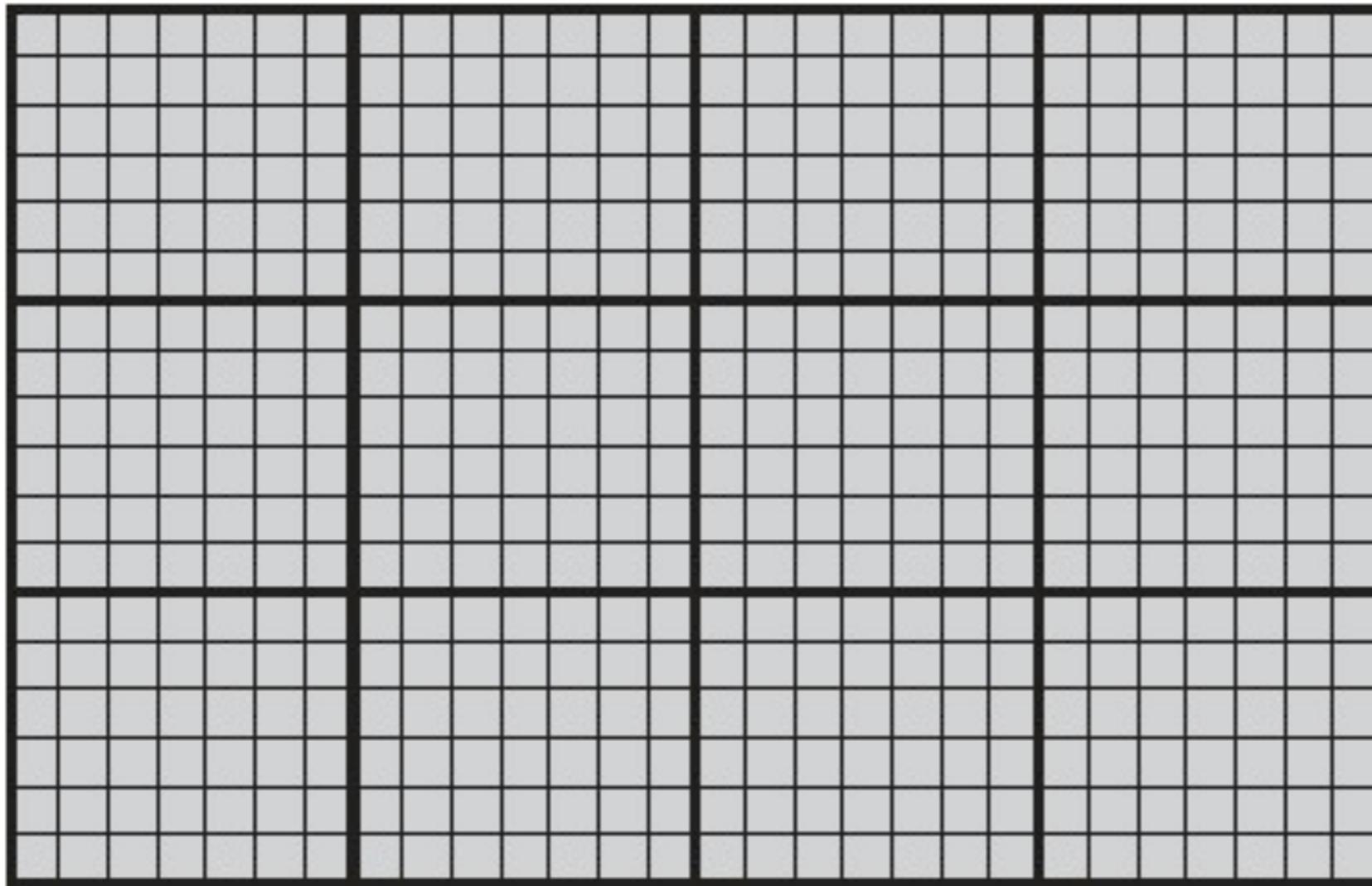
CUDA - Programming model

- ▶ Connecting the hardware and the software



CUDA - Programming model

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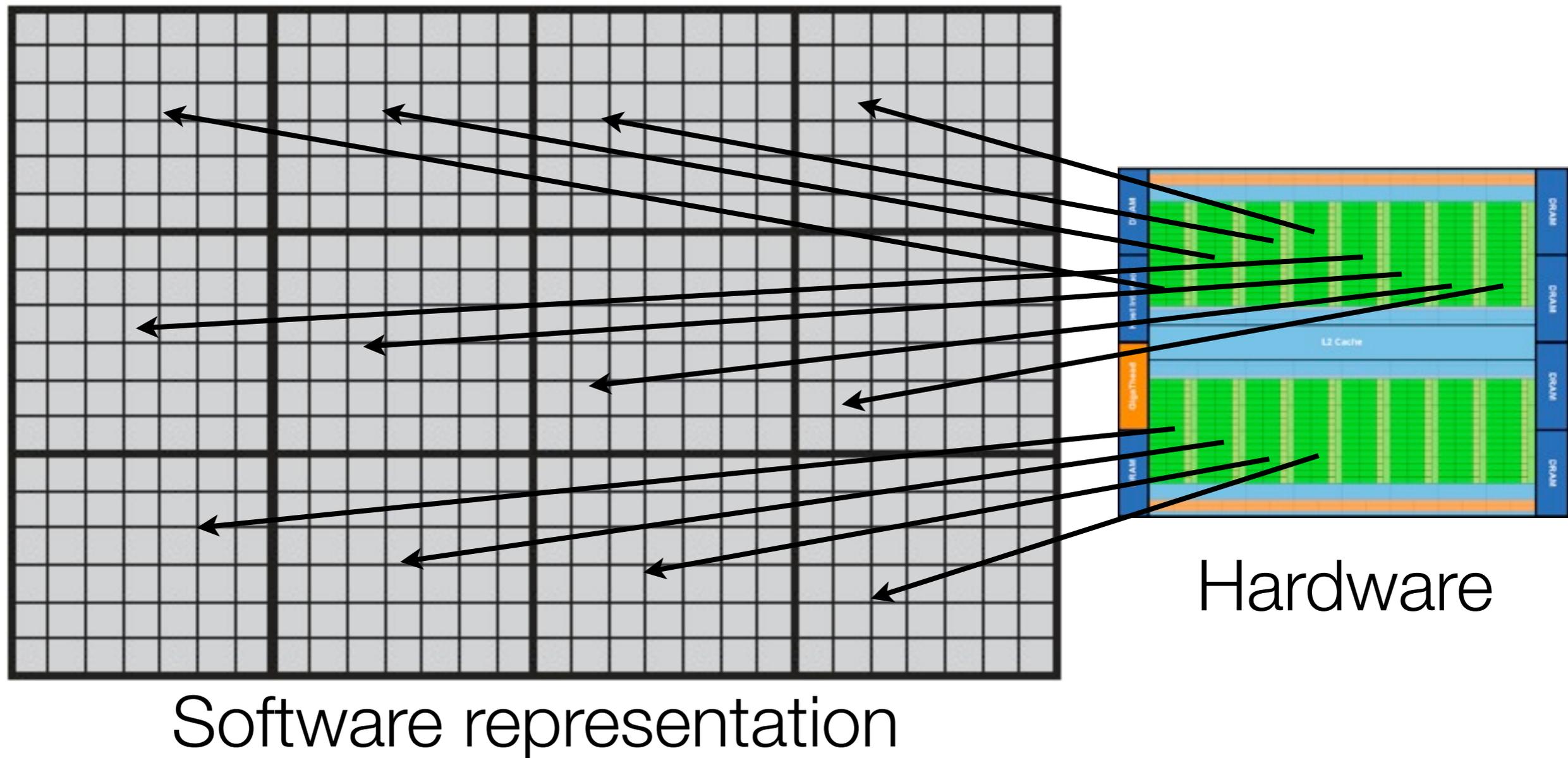
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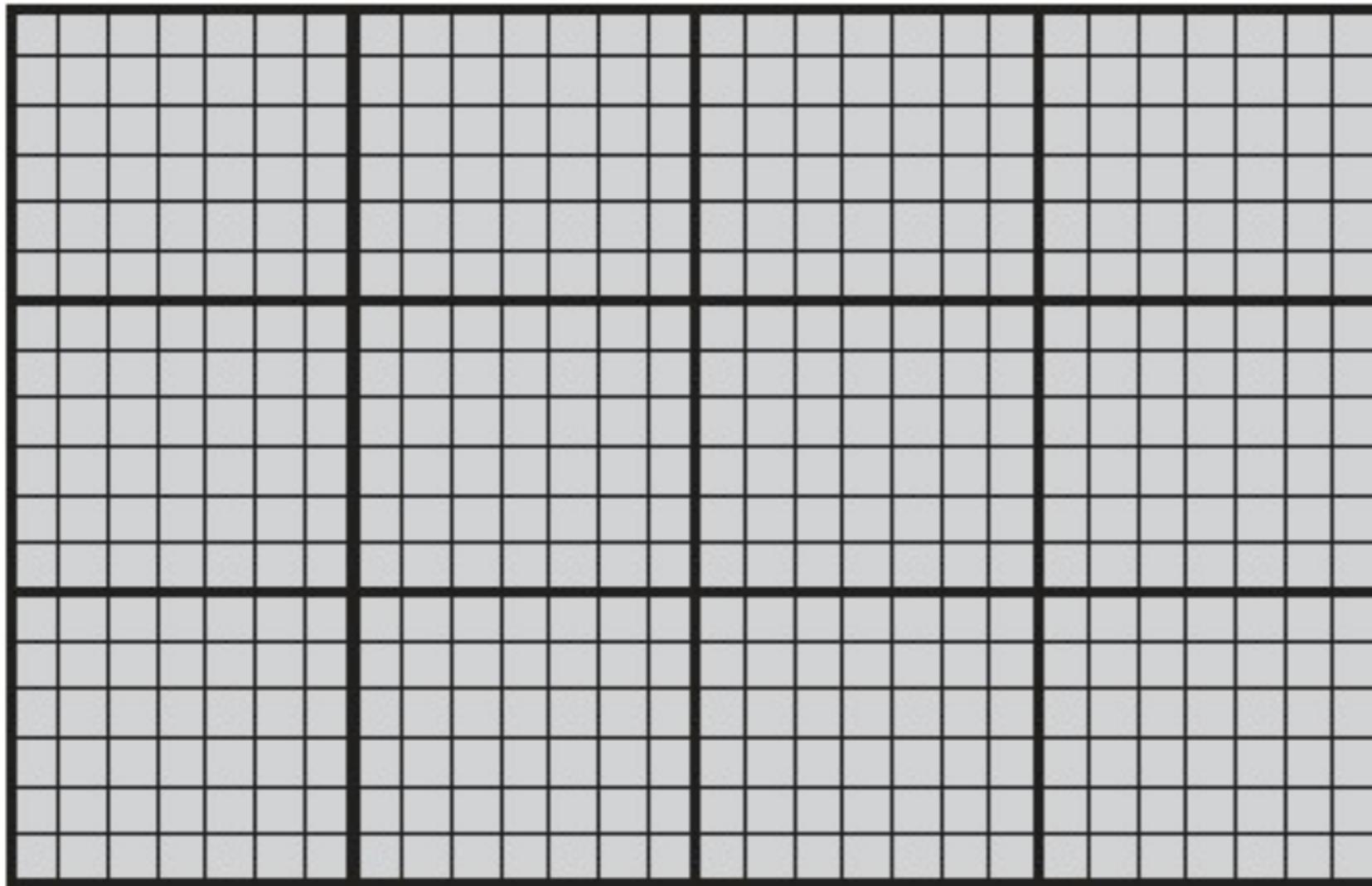
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CUDA - Programming model

- ▶ Connecting the hardware and the software



Software representation



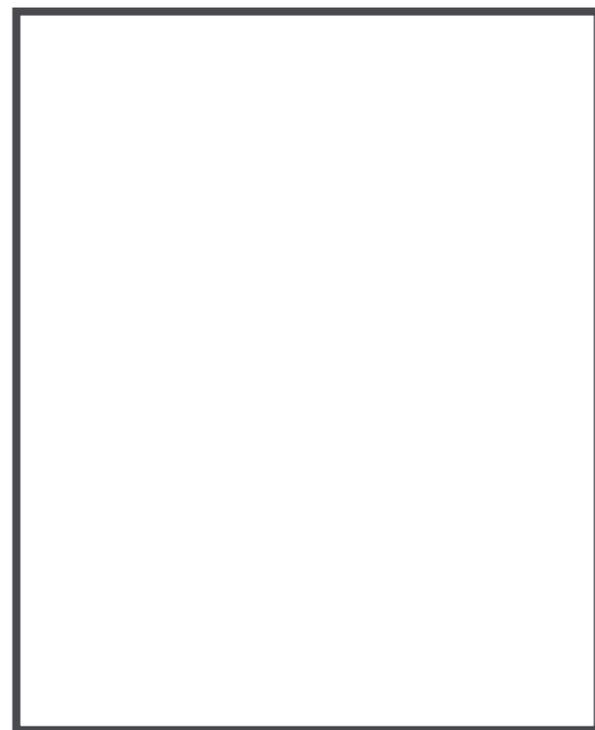
Hardware

How a Streaming Multiprocessor (SM) works

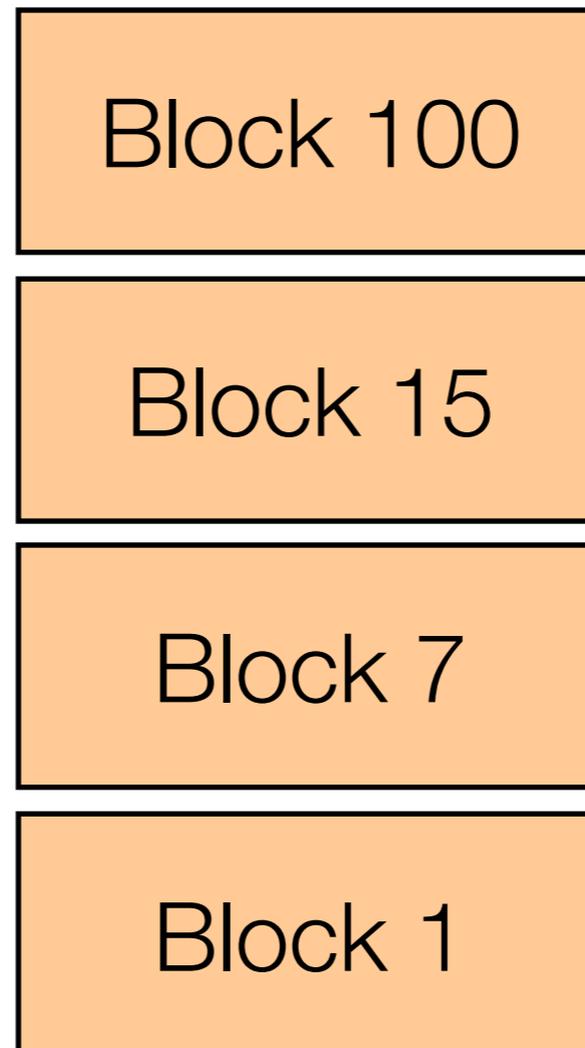
- ▶ CUDA Threads are grouped in thread blocks
 - All threads of the same thread block are executed in the same SM at the same time
 - ▶ SMs have shared memory, then threads within a thread block can communicate
 - The entirety of the threads of a thread block must be executed before there is space to schedule another thread block

How a Streaming Multiprocessor (SM) works

- ▶ Hardware schedules thread blocks onto available SMs
 - No guarantee of order of execution
 - If a SM has more resources the hardware will schedule more blocks

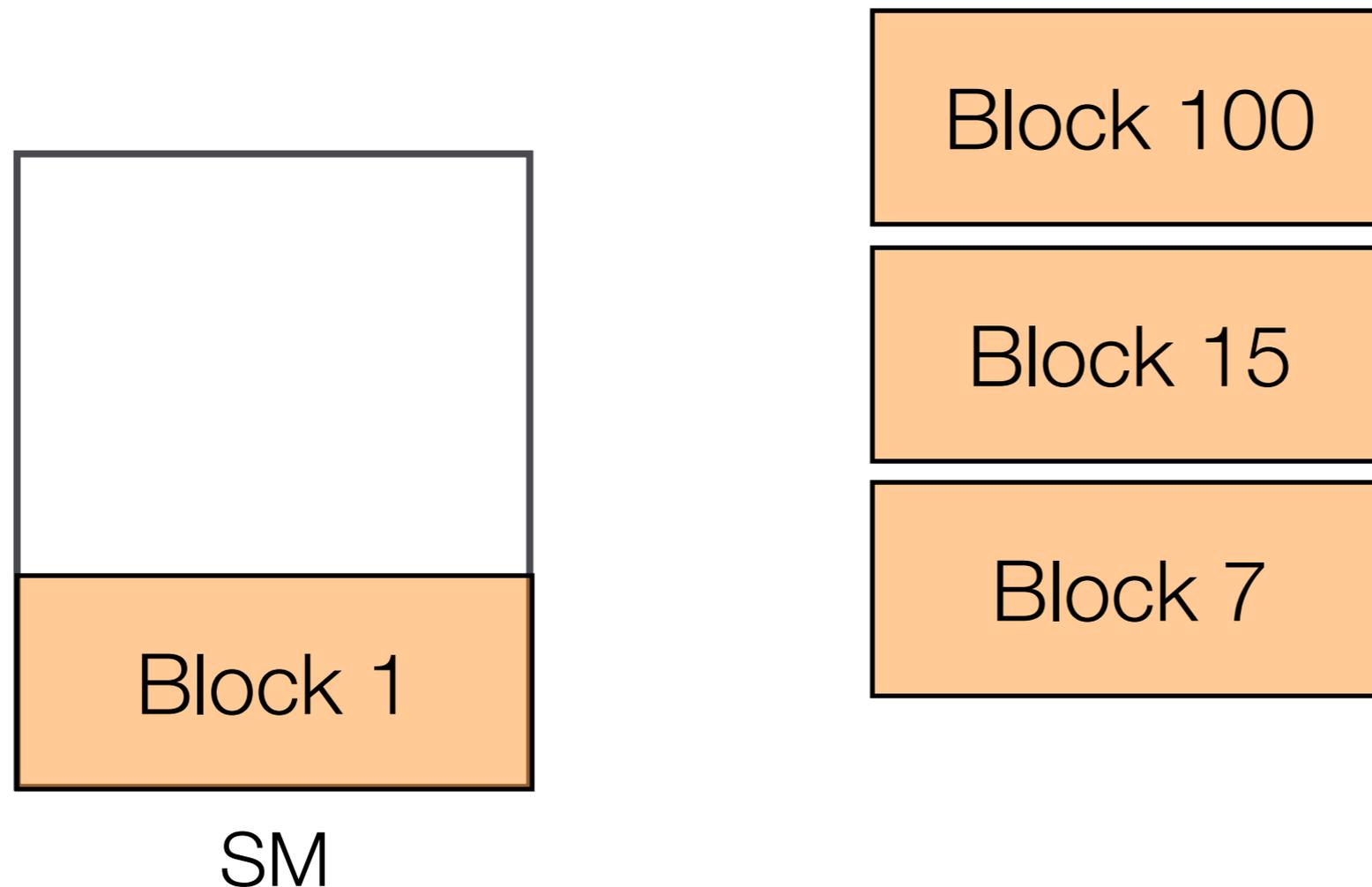


SM



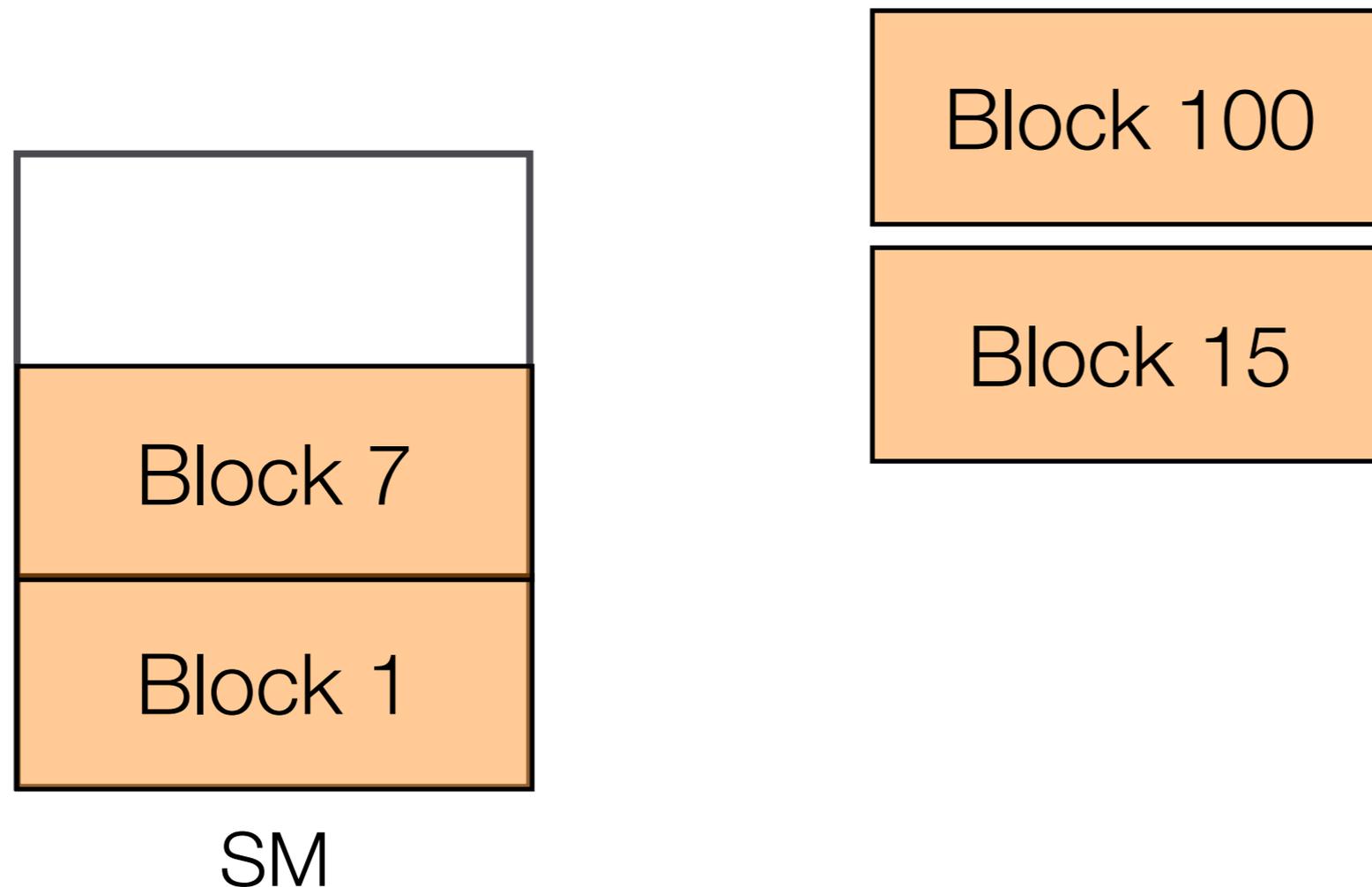
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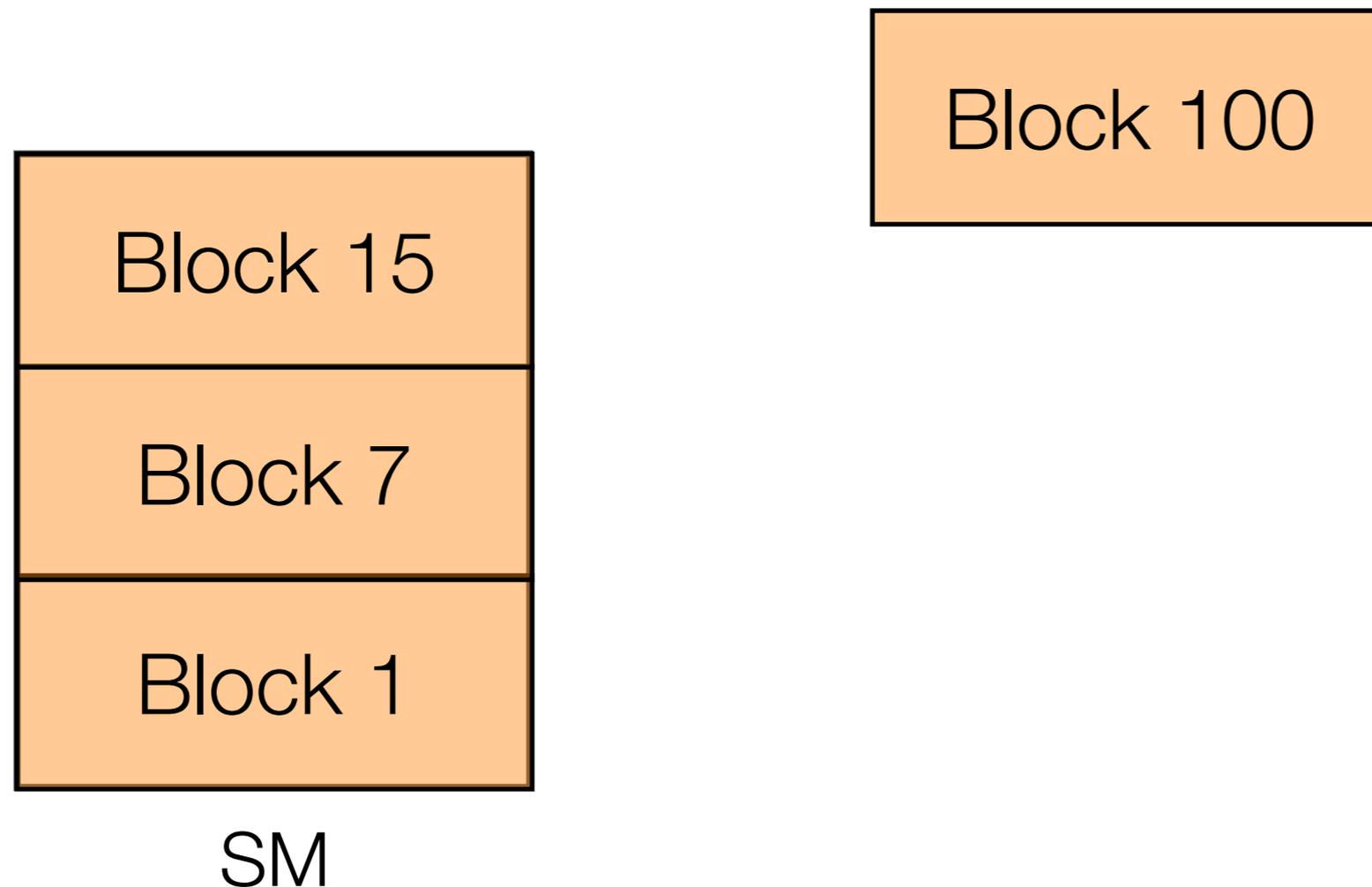
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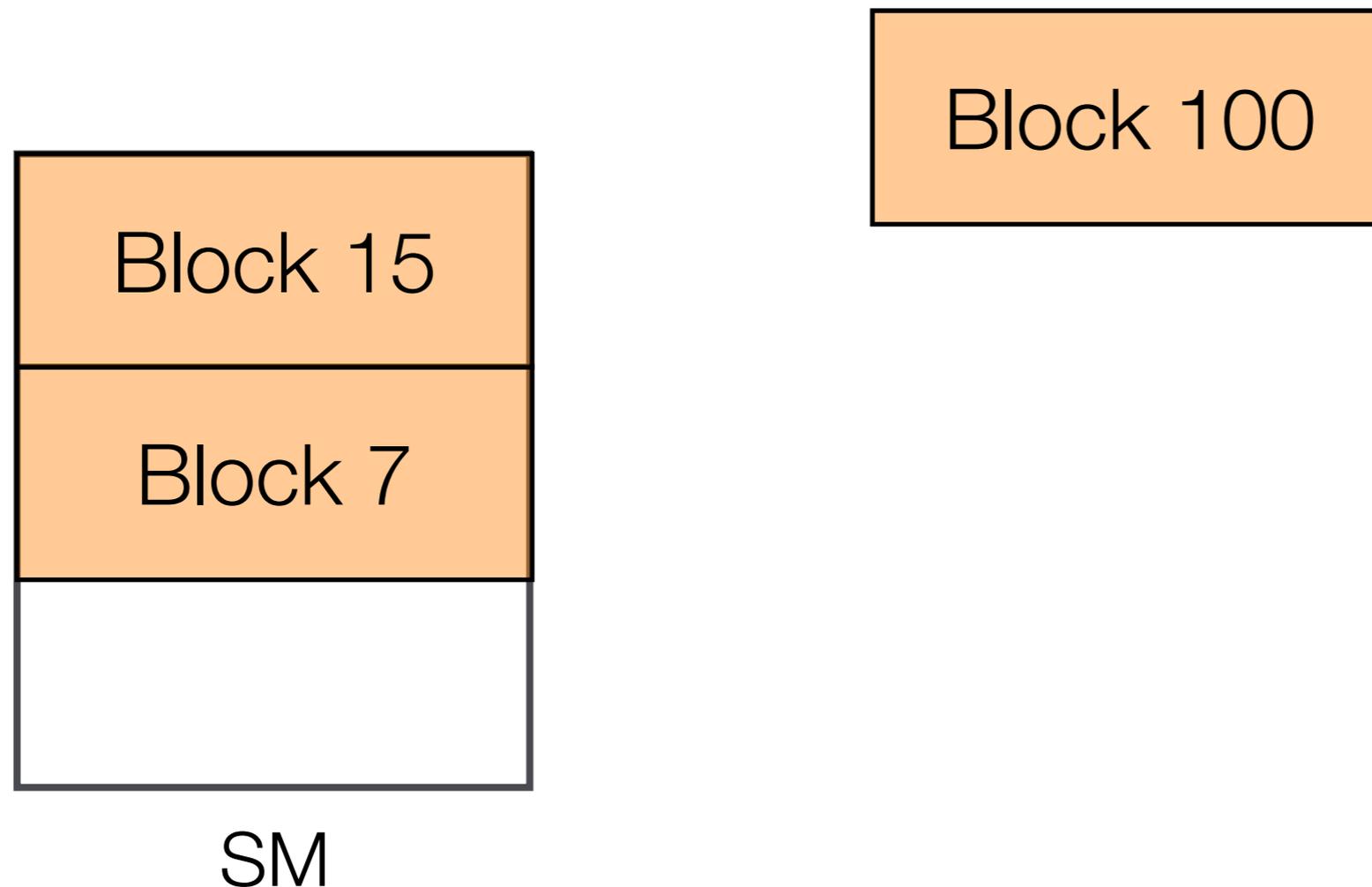
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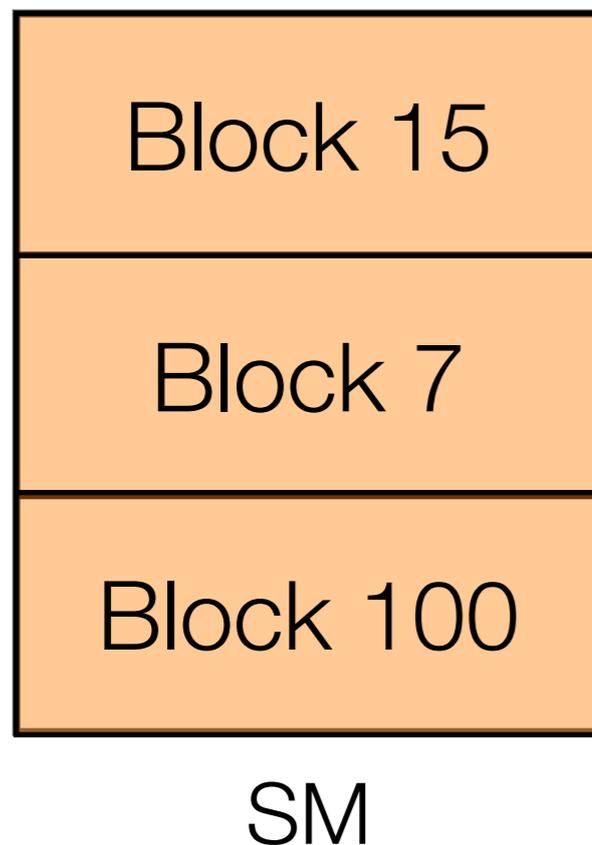
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Warps

- ▶ Inside the SM, threads are launched in groups of 32 called **warps**
 - Warps share the control part (warp scheduler)
 - At any time, only one warp is executed per SM
 - Threads in a warp will be executing the same instruction
 - Half warps for compute capability 1.X
- Fermi:
 - ▶ Maximum number of active threads $1024 * 8 * 32 = 262144$

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- Fermi:

Max threads
per block

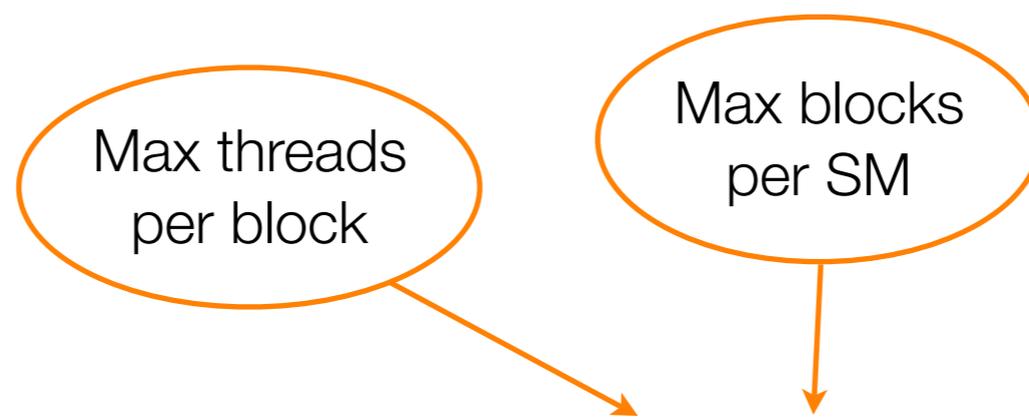


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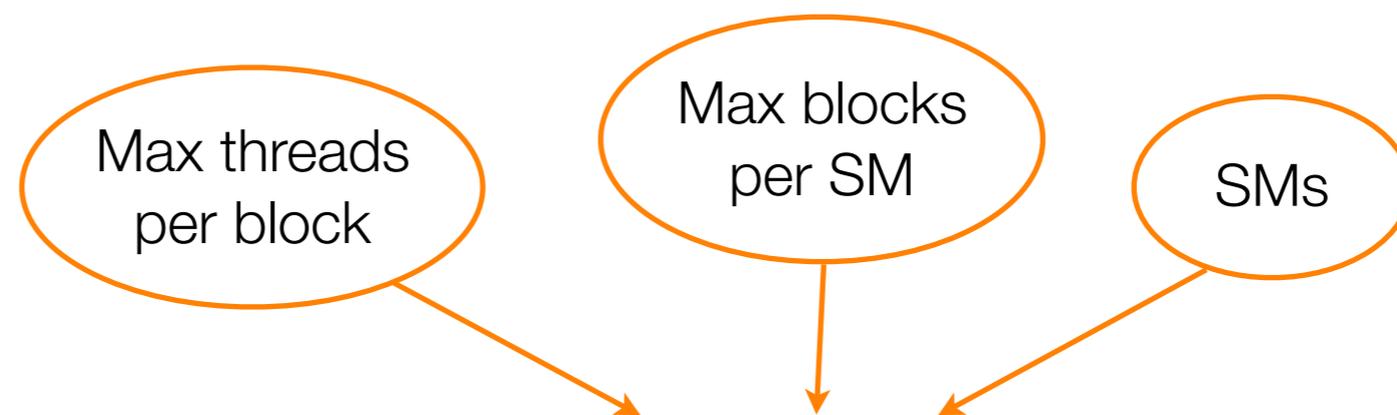


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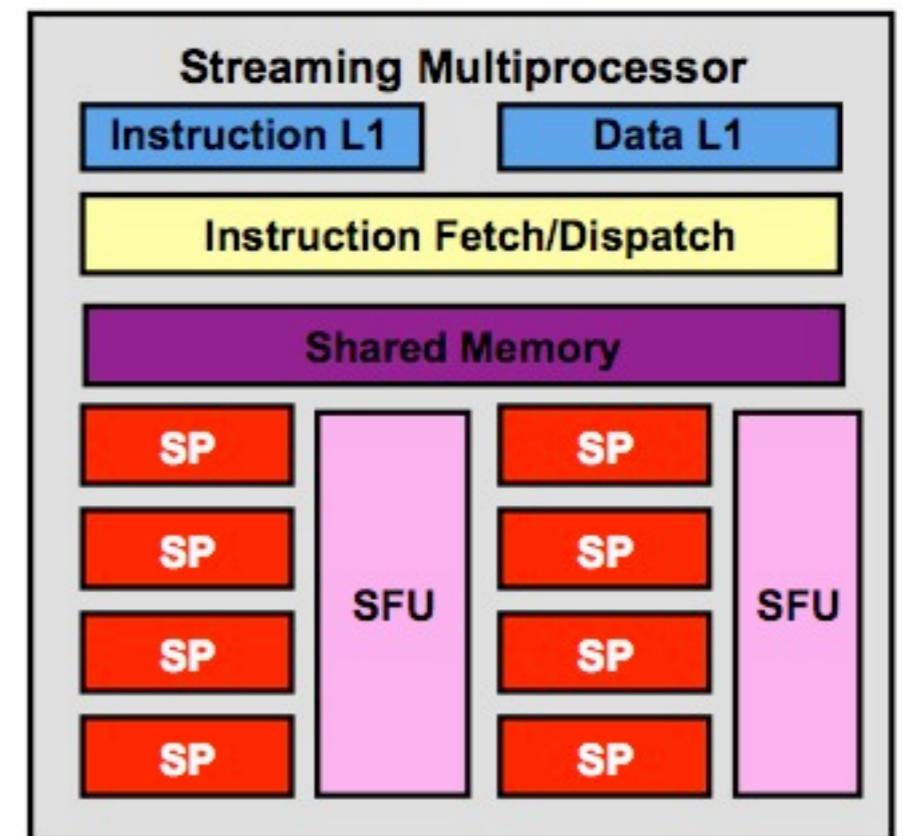
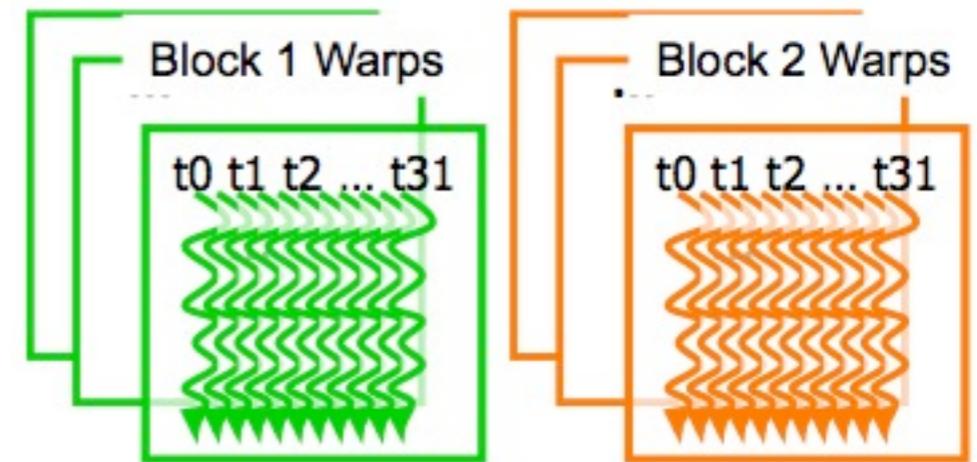
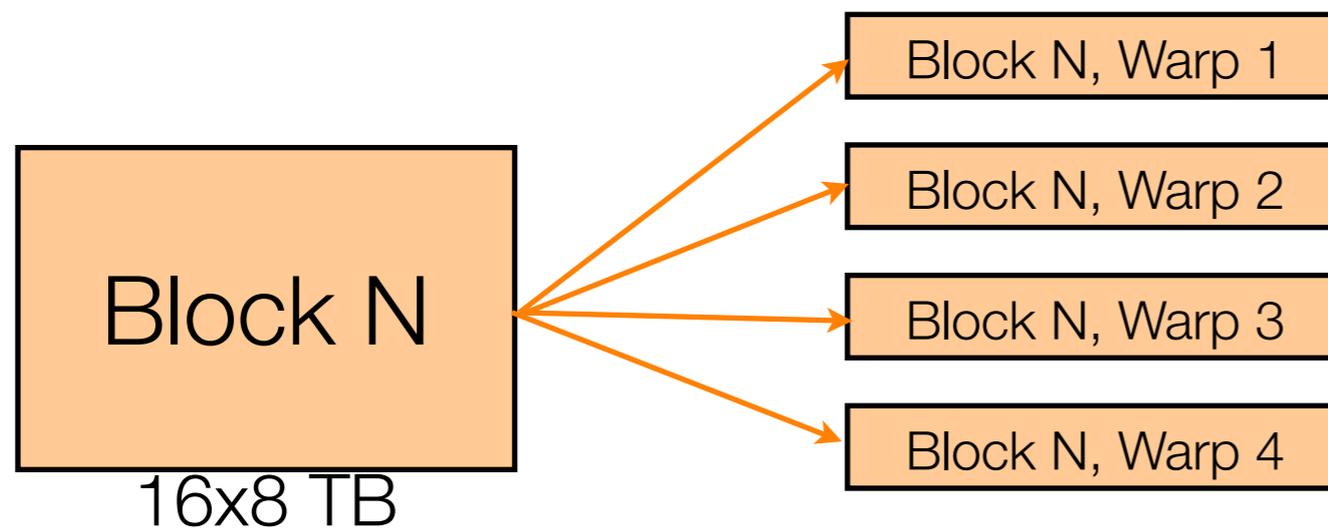
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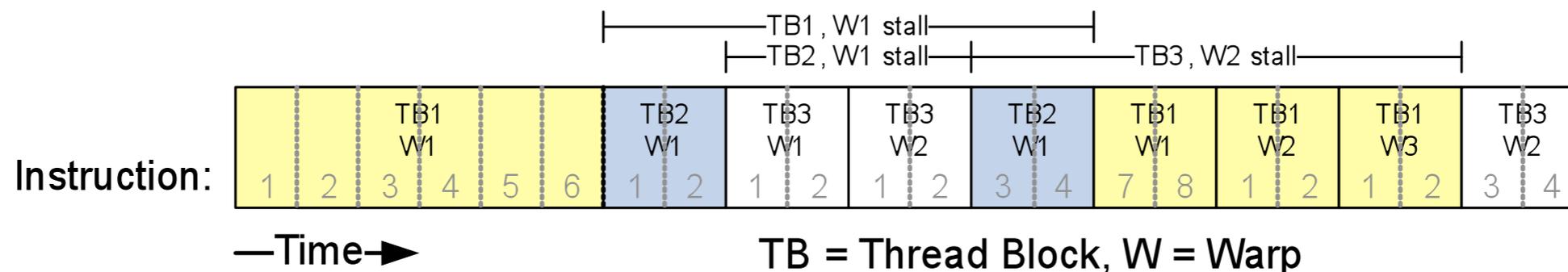
Warp designation

- ▶ Hardware separates threads of a block into warps
 - All threads in a warp correspond to the same thread block
 - Threads are placed in a warp sequentially
 - Threads are scheduled in warps



Warp scheduling

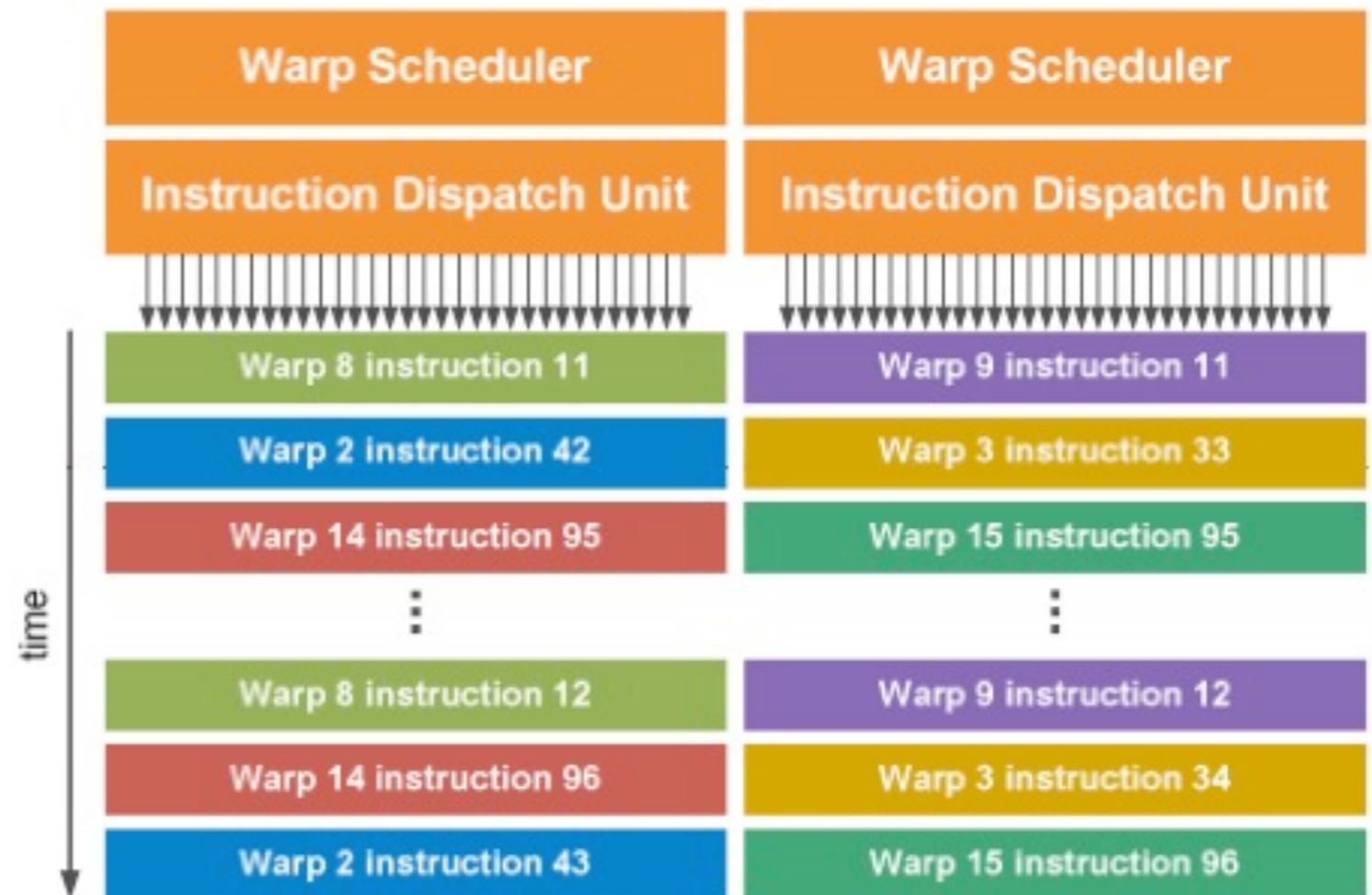
- ▶ The SM implements a zero-overhead warp scheduling
 - Warps whose next instruction has its operands ready for consumption are eligible for execution
 - Eligible warps are selected for execution on a prioritized scheduling policy
 - All threads in a warp execute the same instruction



Warp scheduling

► Fermi

- Double warp scheduler
- Each SM has two warp schedulers and two instruction units



Memory hierarchy

- ▶ CUDA works in both the CPU and GPU
 - One has to keep track of which memory is operating on (host - device)
 - Within the GPU there are also different memory spaces

- Each thread can:
 - Read/write per-thread **registers**
 - Read/write per-thread **local memory**
 - Read/write per-block **shared memory**
 - Read/write per-grid **global memory**
 - Read/only per-grid **constant memory**

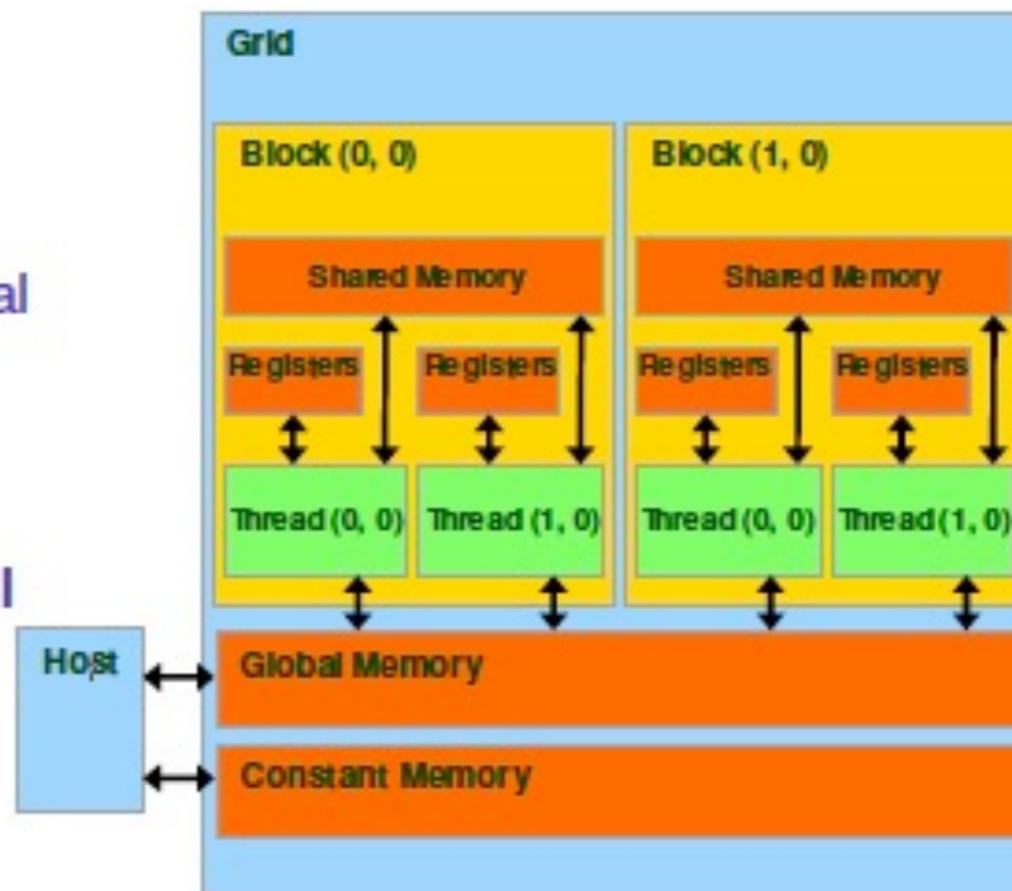
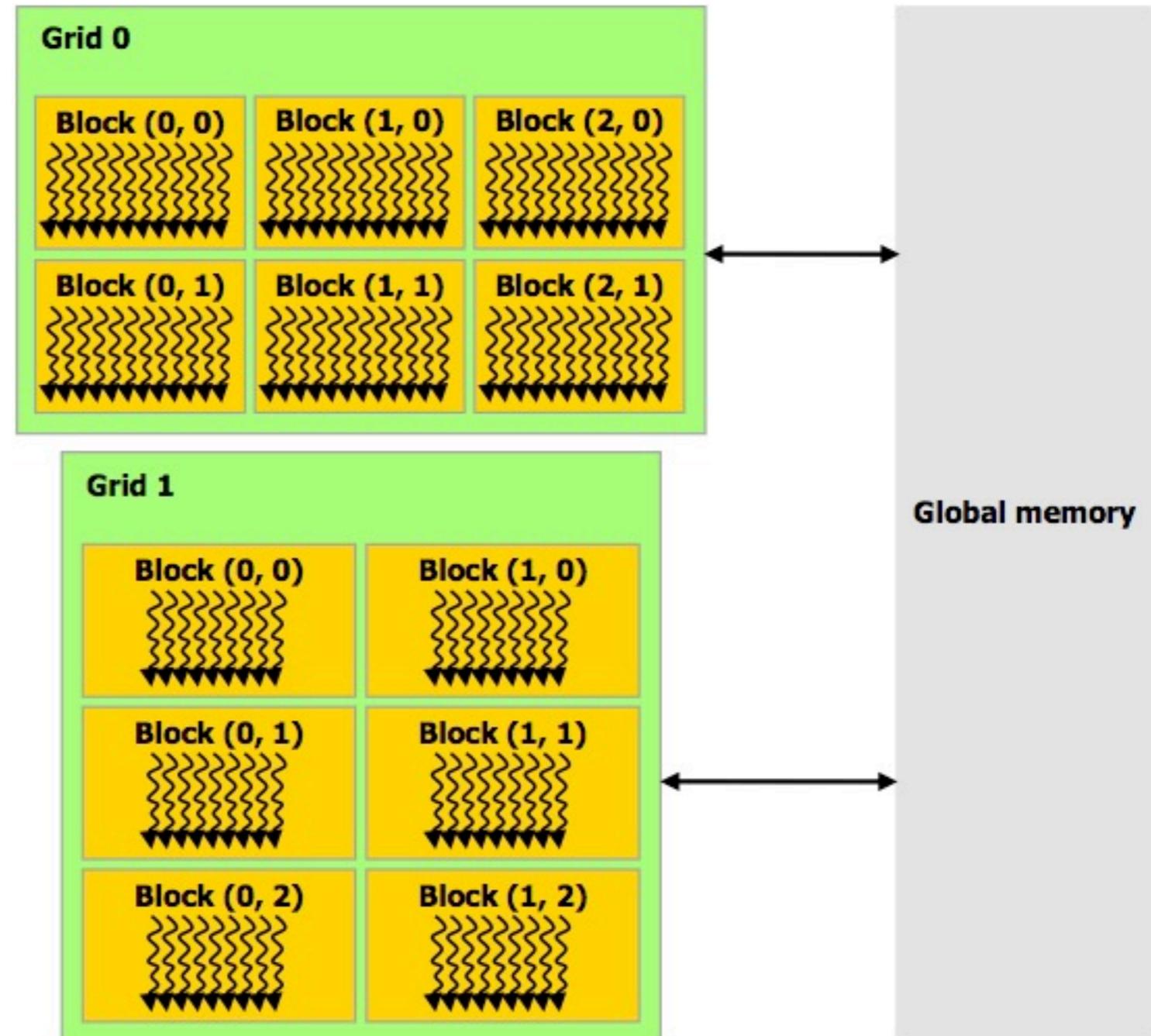


Figure 4.2 GeForce 8800GTX Implementation of CUDA Memories

Memory hierarchy

- ▶ Global memory
 - Main GPU memory
 - Communicates with host
 - Can be seen by all threads
 - Order of GB
 - Off chip, slow (~400 cycles)

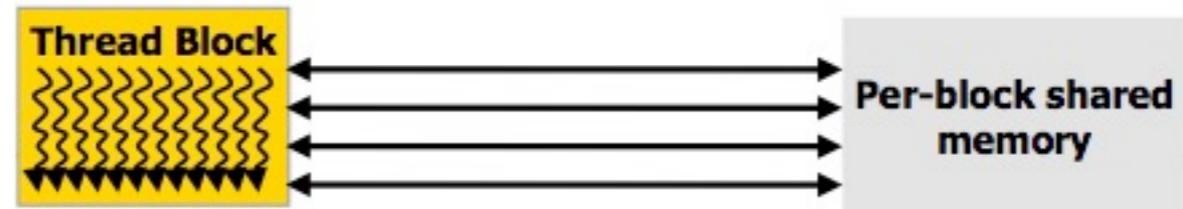


```
__device__ float variable;
```

Memory hierarchy

▶ Shared memory

- Per SM
- Seen by threads of the same thread block
- Order of kB
- On chip, fast (~4 cycles)



```
__shared__ float variable;
```

▶ Registers

- Private to each thread
- On chip, fast

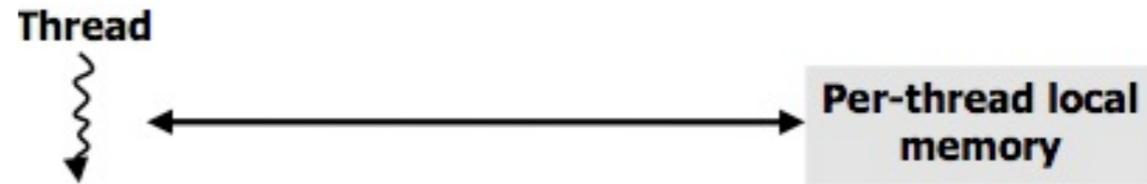


```
float variable;
```

Memory hierarchy

▶ Local memory

- Private to each thread
- Off chip, slow
- Register overflows



```
float variable [10];
```

▶ Constant memory

- Read only
- Off chip, but fast (cached)
- Seen by all threads
- 64kB with 8kB cache

```
__constant__ float variable;
```

Memory hierarchy

▶ Texture memory

- Seen by all threads
- Read only
- Off chip, but fast (cached) if cache hit
- Cache optimized for 2D locality
- Binds to global

```
texture<type, dim> tex_var;  
cudaChannelFormatDesc();  
cudaBindTexture2D(...);  
tex2D(tex_var, x_index, y_index);
```

Memory hierarchy

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Initialize



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Options



Memory hierarchy

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Initialize

Options

Bind

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Initialize

Options

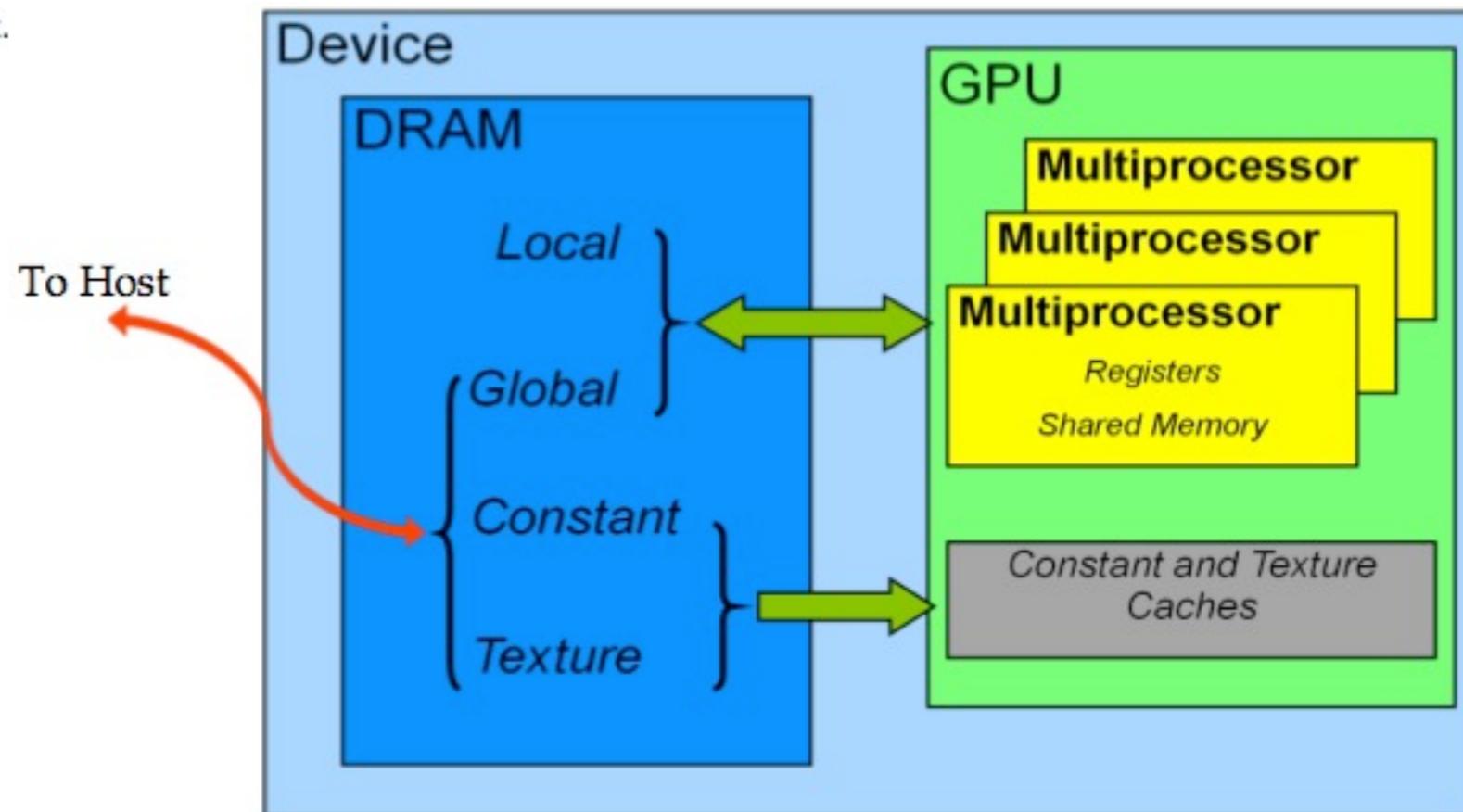
Bind

Fetch

Memory hierarchy

| Memory | Location on/off chip | Cached | Access | Scope | Lifetime |
|----------|----------------------|--------|--------|----------------------|-----------------|
| Register | On | n/a | R/W | 1 thread | Thread |
| Local | Off | † | R/W | 1 thread | Thread |
| Shared | On | n/a | R/W | All threads in block | Block |
| Global | Off | † | R/W | All threads + host | Host allocation |
| Constant | Off | Yes | R | All threads + host | Host allocation |
| Texture | Off | Yes | R | All threads + host | Host allocation |

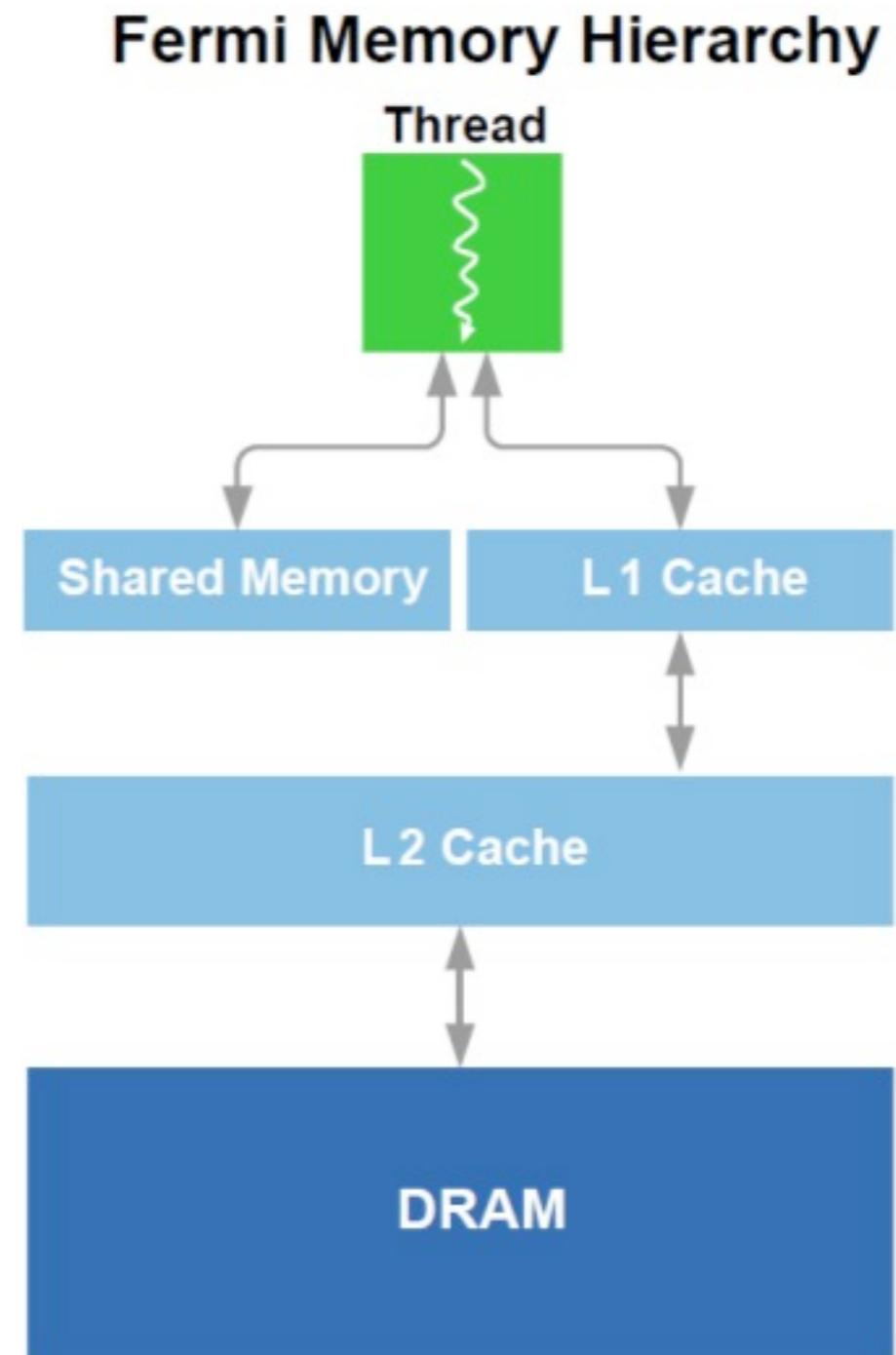
†Cached only on devices of compute capability 2.x.



Memory hierarchy

▶ Case of Fermi

- Added an L1 cache to each SM
- Shared + cache = 64kB:
 - ▶ Shared = 48kB, cache = 16kB
 - ▶ Shared = 16kB, cache = 48kB



Memory hierarchy

- ▶ Use you memory strategically
 - Read only: **__constant__** (fast)
 - Read/write and communicate within a block: **__shared__** (fast and communication)
 - Read/write inside thread: registers (fast)
 - Data locality: texture

Resource limits

- ▶ Number of thread blocks per SM at the same time is limited by
 - Shared memory usage
 - Registers
 - No more than 8 thread blocks per SM
 - Number of threads

Resource limits - Examples

Number of blocks

How big should my blocks be? 8x8, 16x16 or 64x64?

▶ 8x8

8*8 = 64 threads per block, 1024/64 = 16 blocks. An SM can have up to 8 blocks: only 512 threads will be active at the same time

Max threads
per block



▶ 16x16

16*16 = 256 threads per block, 1024/256 = 4 blocks. An SM can take all blocks, then all 1024 threads will be active and achieve full capacity unless other resource overrule

▶ 64x64

64*64 = 4096 threads per block: doesn't fit in a SM

Resource limits - Examples

Registers

We have a kernel that uses 10 registers. With 16x16 block, how many blocks can run in G80 (max 8192 registers per SM, 768 threads per SM)?

$10 * 16 * 16 = 2560$. SM can hold $8192 / 2560 = 3$ blocks, meaning we will use $3 * 16 * 16 = 768$ threads, which is within limits.

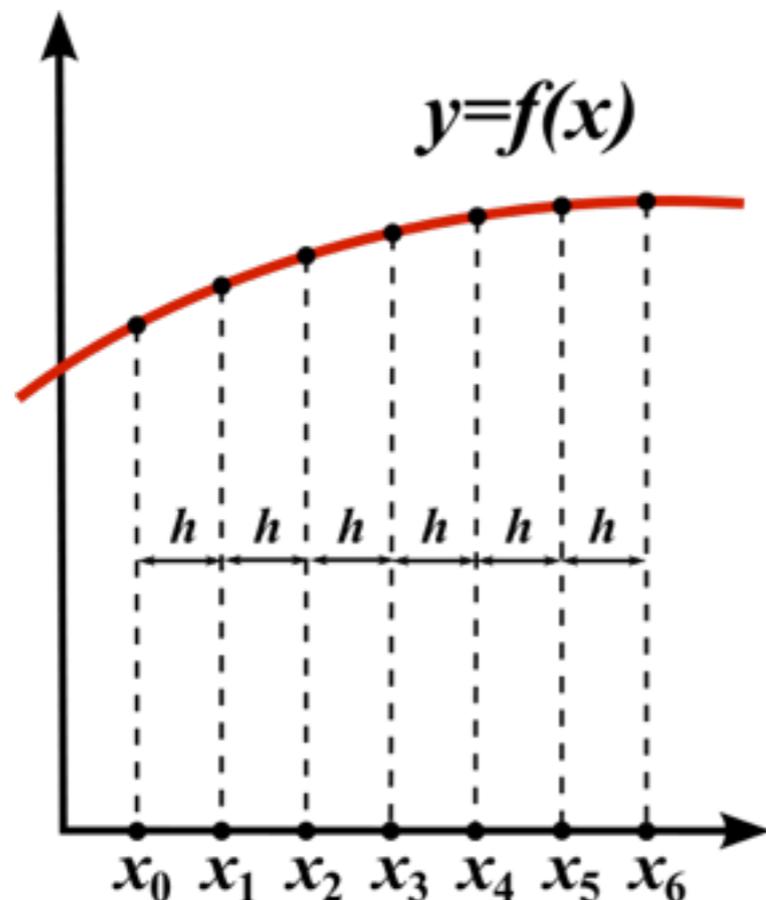
If we add one more register, the number of registers grows to $11 * 16 * 16 = 2816$. SM can hold $8192 / 2816 = 2$ blocks, meaning we will use $2 * 16 * 16 = 512$ threads. Now as less threads are running per SM is more difficult to have enough warps to have the GPU always busy!

Introduction to Finite Difference Method

- ▶ Throughout the course many examples will be taken from Finite Difference Method
- ▶ Also, next lab involves implementing a stencil code
- ▶ Just want to make sure we're all in the same page!

Introduction to Finite Difference Method

- ▶ Finite Difference Method (FDM) is a numerical method for solution of differential equations
- ▶ Domain is discretized with a mesh
- ▶ Derivative at a node are approximated by the linear combination of the values of points nearby (including itself)



$$\frac{\partial y}{\partial x_i} \approx \frac{y_i - y_{i-1}}{x_i - x_{i-1}}$$

Example using first order
one sided difference

FDM - Accuracy and order

- ▶ From Taylor's polynomial, we get

$$f(x_0 + h) = f(x_0) + f'(x_0) \cdot h + R_1(x)$$

$$f'(x_0) = \frac{f(a + h) - f(a)}{h} - R(x)$$

- ▶ Sources of error

- Roundoff (Machine)
- Truncation ($R(x)$): gives the order of the method

FDM - Stability

- ▶ Stability of the algorithm may be conditional

$$\frac{\partial u}{\partial t} = \frac{\partial^2 u}{\partial x^2}$$

- ▶ Explicit method

$$\frac{u_j^{n+1} - u_j^n}{k} = \frac{u_{j+1}^n - 2u_j^n + u_{j-1}^n}{h^2} \quad \text{Conditionally stable} \quad \frac{k}{h^2} < 0.5$$

- ▶ Implicit

$$\frac{u_j^{n+1} - u_j^n}{k} = \frac{u_{j+1}^{n+1} - 2u_j^{n+1} + u_{j-1}^{n+1}}{h^2} \quad \text{Unconditionally stable}$$

- ▶ Crank Nicolson

$$\frac{u_j^{n+1} - u_j^n}{k} = \frac{1}{2} \left(\frac{u_{j+1}^n - 2u_j^n + u_{j-1}^n}{h^2} + \frac{u_{j+1}^{n+1} - 2u_j^{n+1} + u_{j-1}^{n+1}}{h^2} \right)$$

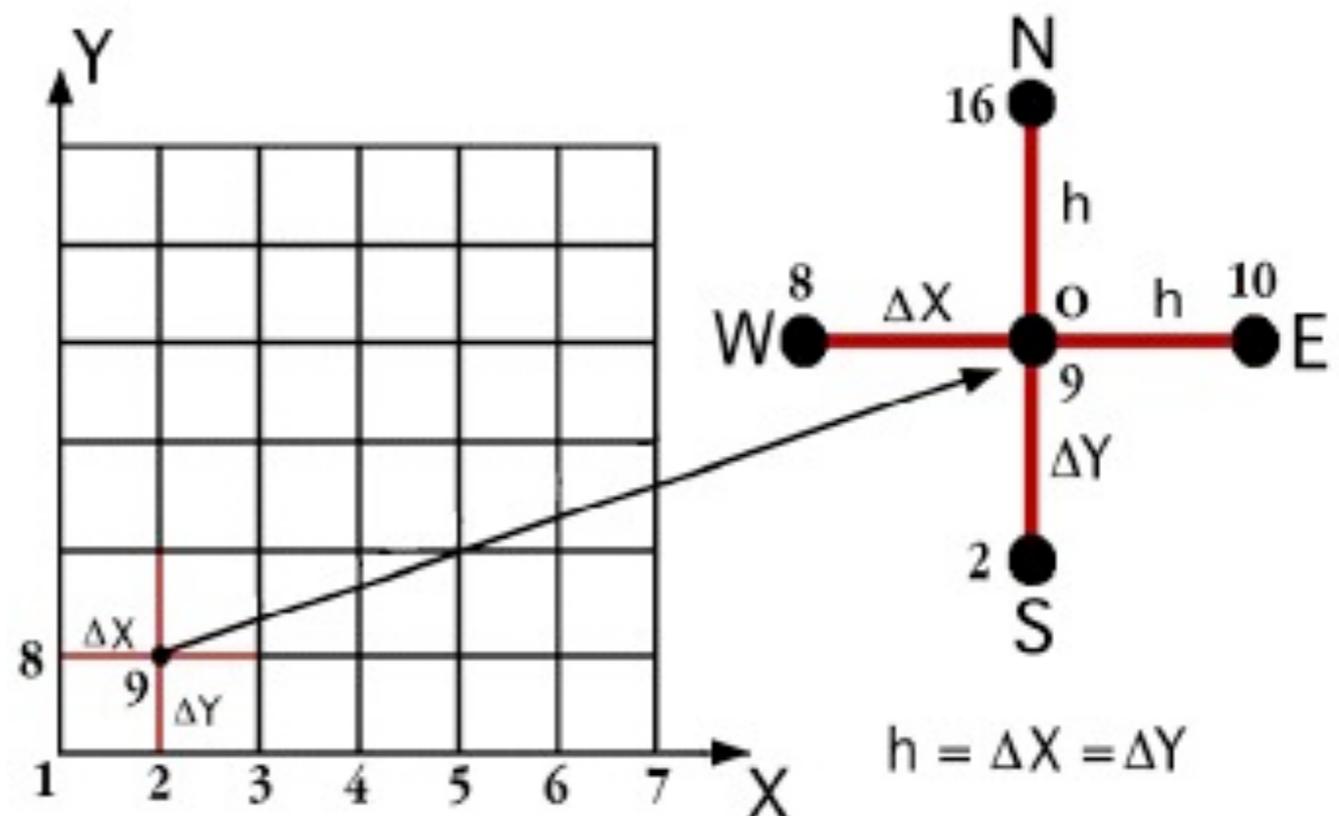
Unconditionally stable

FDM - 2D example

- ▶ We're going to be working with this example in the labs
- ▶ Explicit diffusion

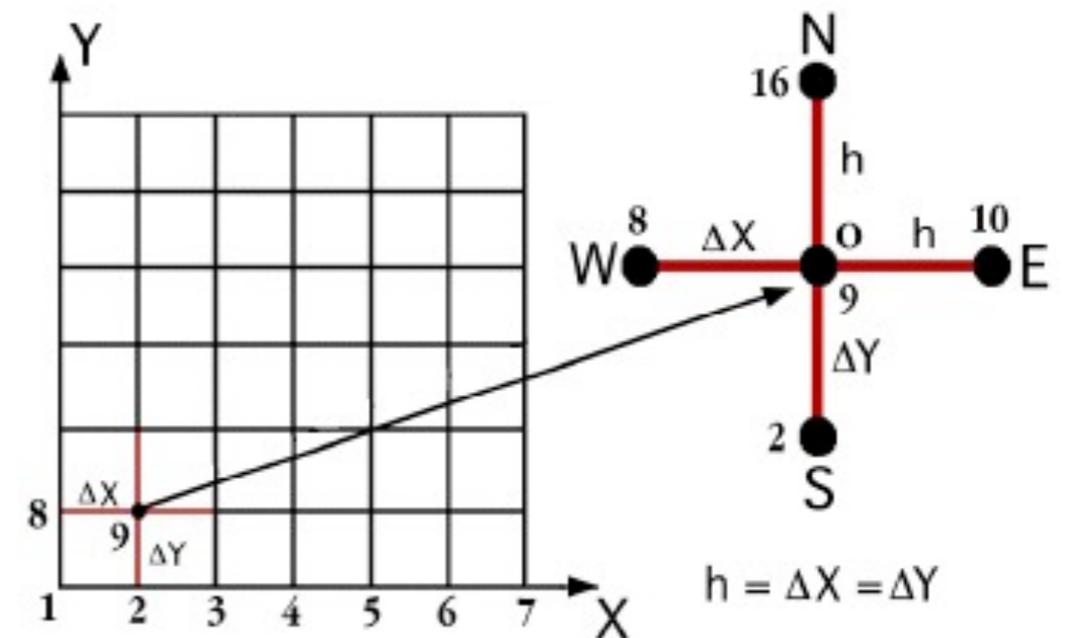
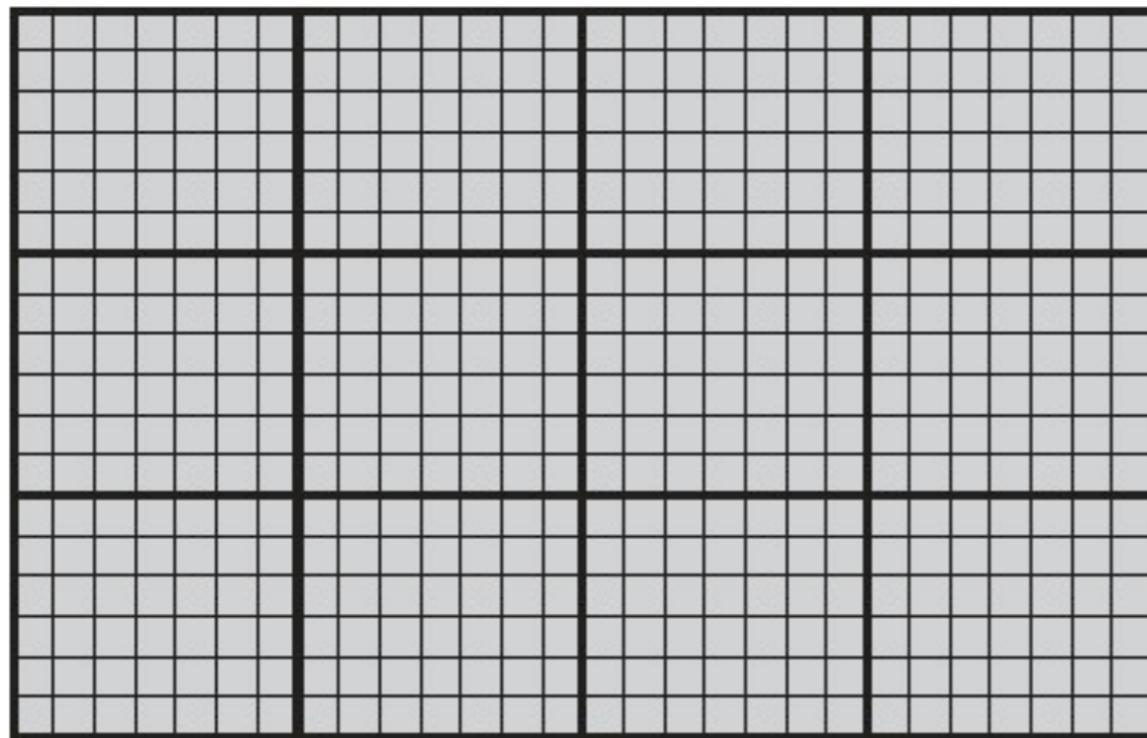
$$\frac{\partial u}{\partial t} = \alpha \frac{\partial^2 u}{\partial x^2}$$

$$u_{i,j}^{n+1} = u_{i,j}^n + \frac{\alpha k}{h^2} (u_{i,j+1}^n + u_{i,j-1}^n + u_{i+1,j}^n + u_{i-1,j}^n - 4u_{i,j}^n)$$



FDM on the GPU

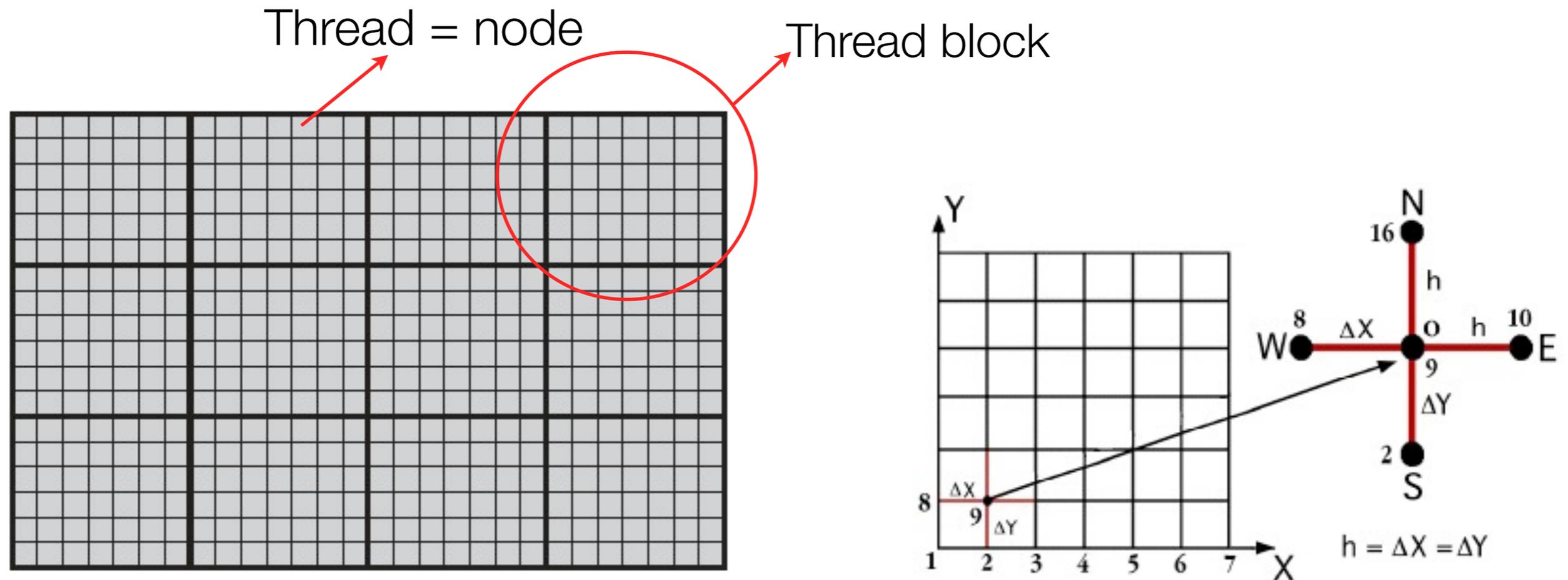
- ▶ Mapping the physical problem to the software and hardware
- ▶ Each thread will operate on one node: node indexing groups them naturally!



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FDM on the GPU

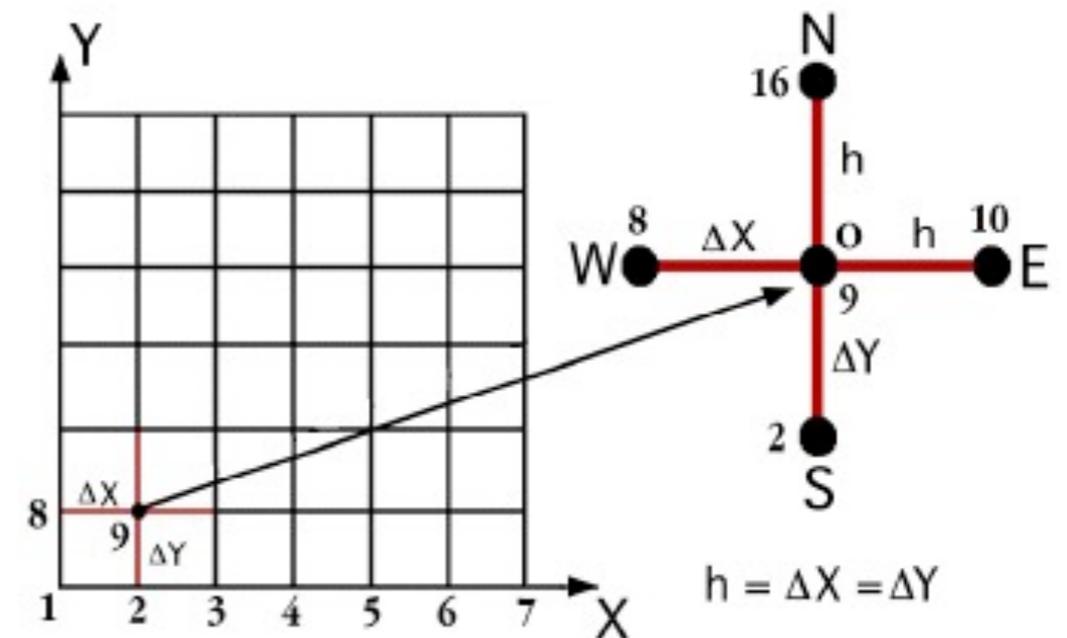
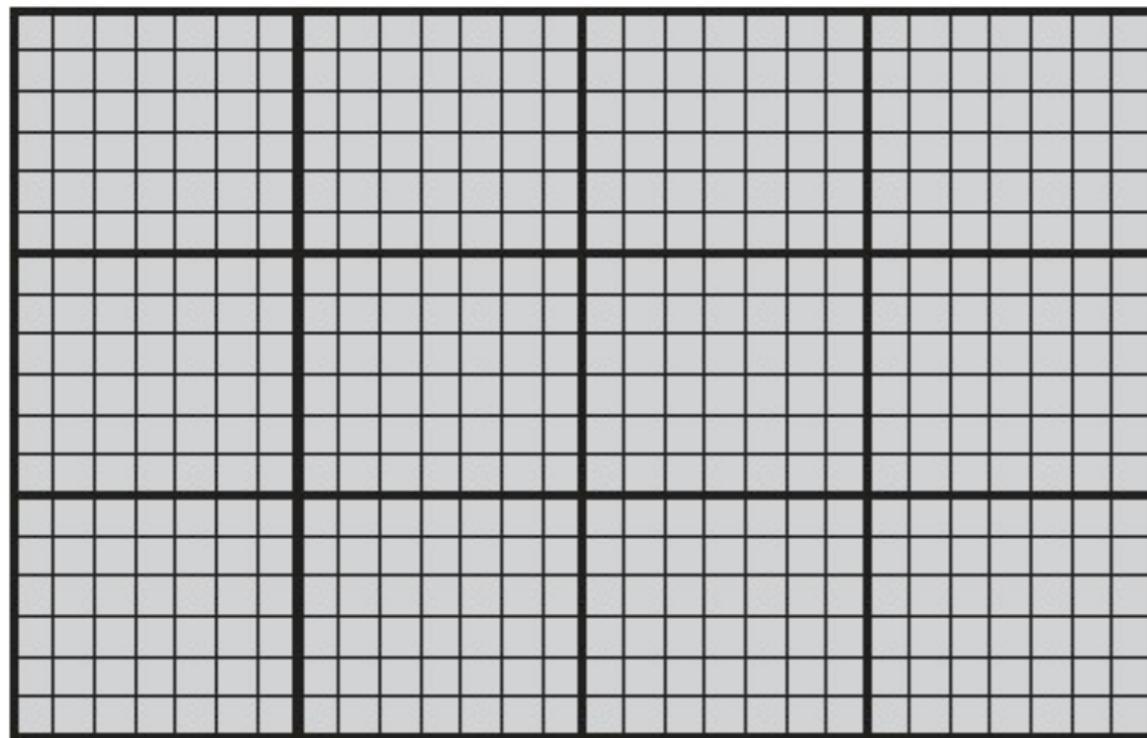
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FDM on the GPU

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