

# Computer Architecture

## Lecture 26: GPU Programming

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Fall 2022

06 January 2023

# Agenda for Today

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- GPU as an accelerator
  - Program structure
    - Bulk synchronous programming model
  - Memory hierarchy and memory management
  - Performance considerations
    - Memory access
    - SIMD utilization
    - Atomic operations
    - Data transfers
- Collaborative computing

# Recommended Readings

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- CUDA Programming Guide
  - <https://docs.nvidia.com/cuda/cuda-c-programming-guide/index.html>
- Hwu and Kirk, “Programming Massively Parallel Processors,” Third Edition, 2017

# An Example GPU

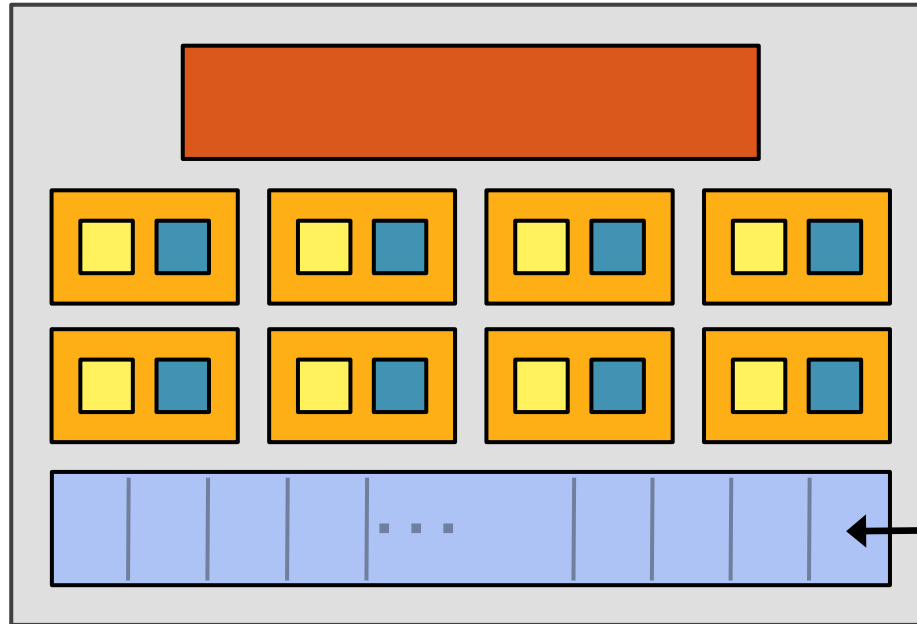
# Recall: NVIDIA GeForce GTX 285

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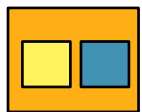
- NVIDIA-speak:
  - 240 stream processors
  - “SIMT execution”
- Generic speak:
  - 30 cores
  - 8 SIMD functional units per core



# NVIDIA GeForce GTX 285 “core”



64 KB of storage  
for thread contexts  
(registers)



= SIMD functional unit, control  
shared across 8 units



= multiply-add



= multiply



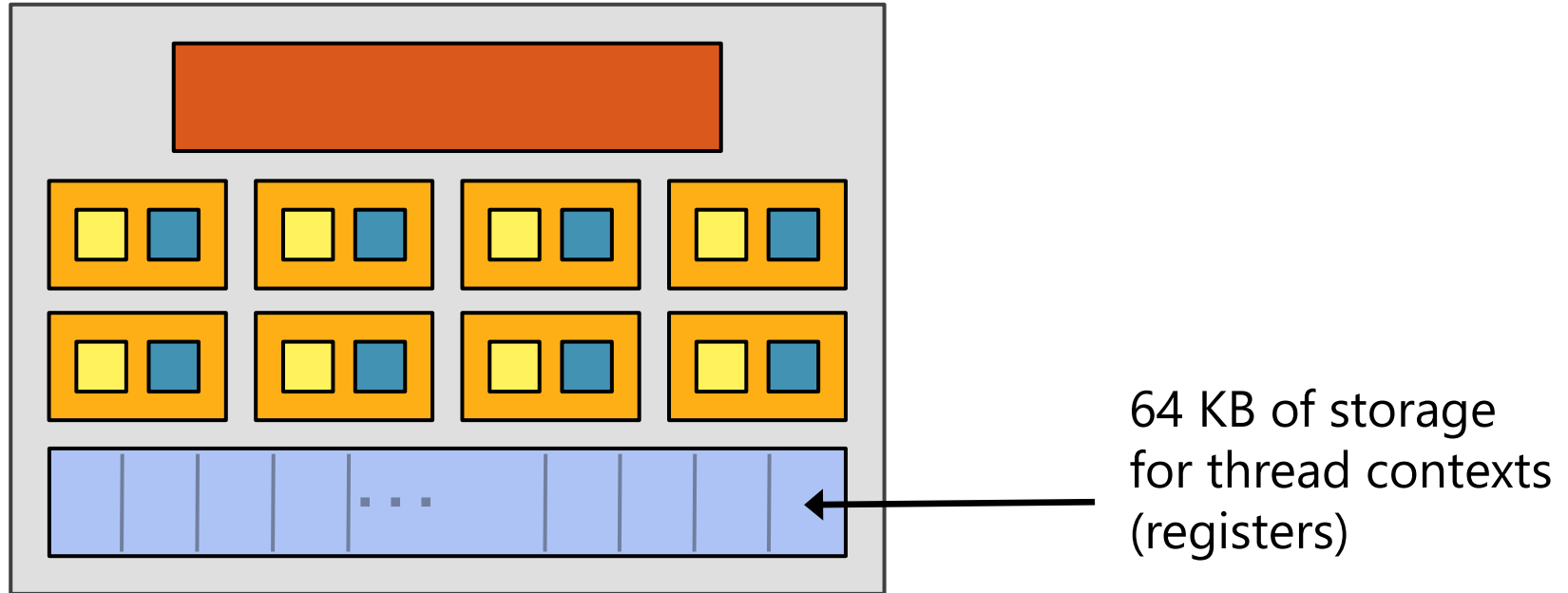
= instruction stream decode



= execution context storage

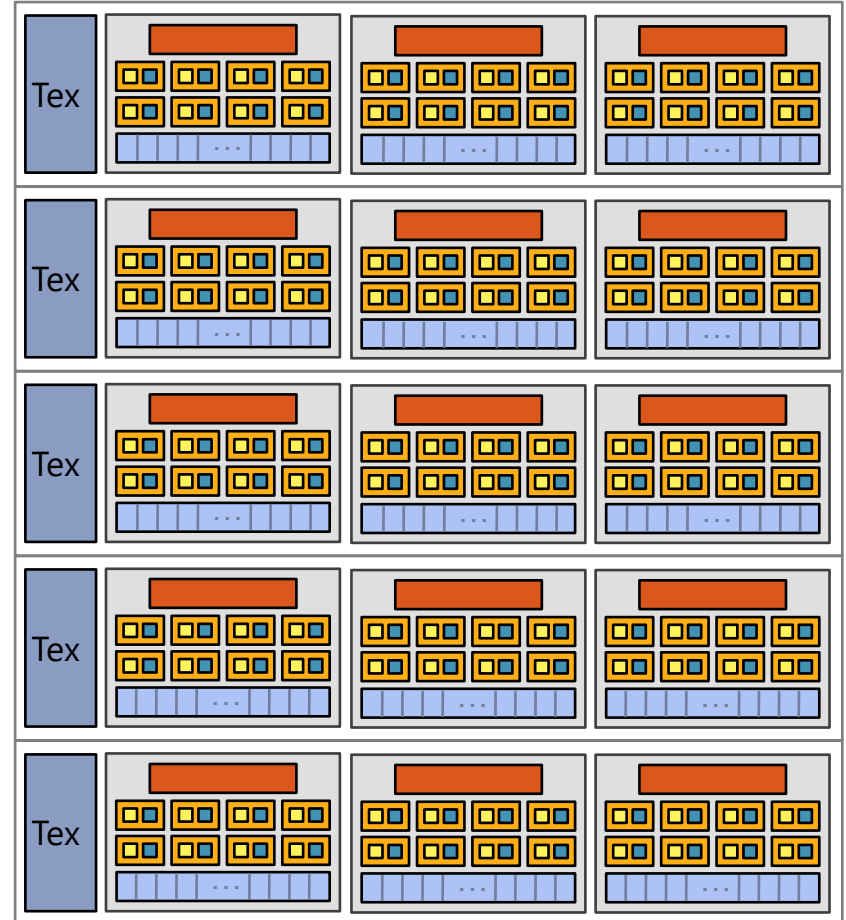
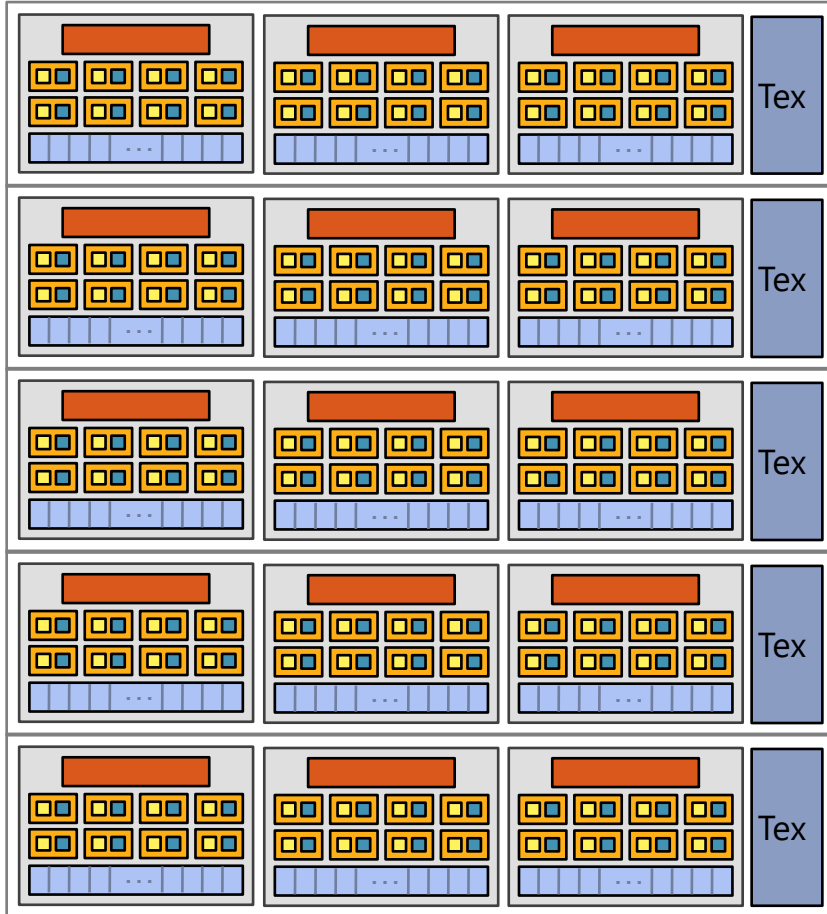
# NVIDIA GeForce GTX 285 “core”

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- Groups of 32 **threads** share instruction stream (each group is a Warp)
- Up to 32 warps are simultaneously interleaved
- Up to 1024 thread contexts can be stored

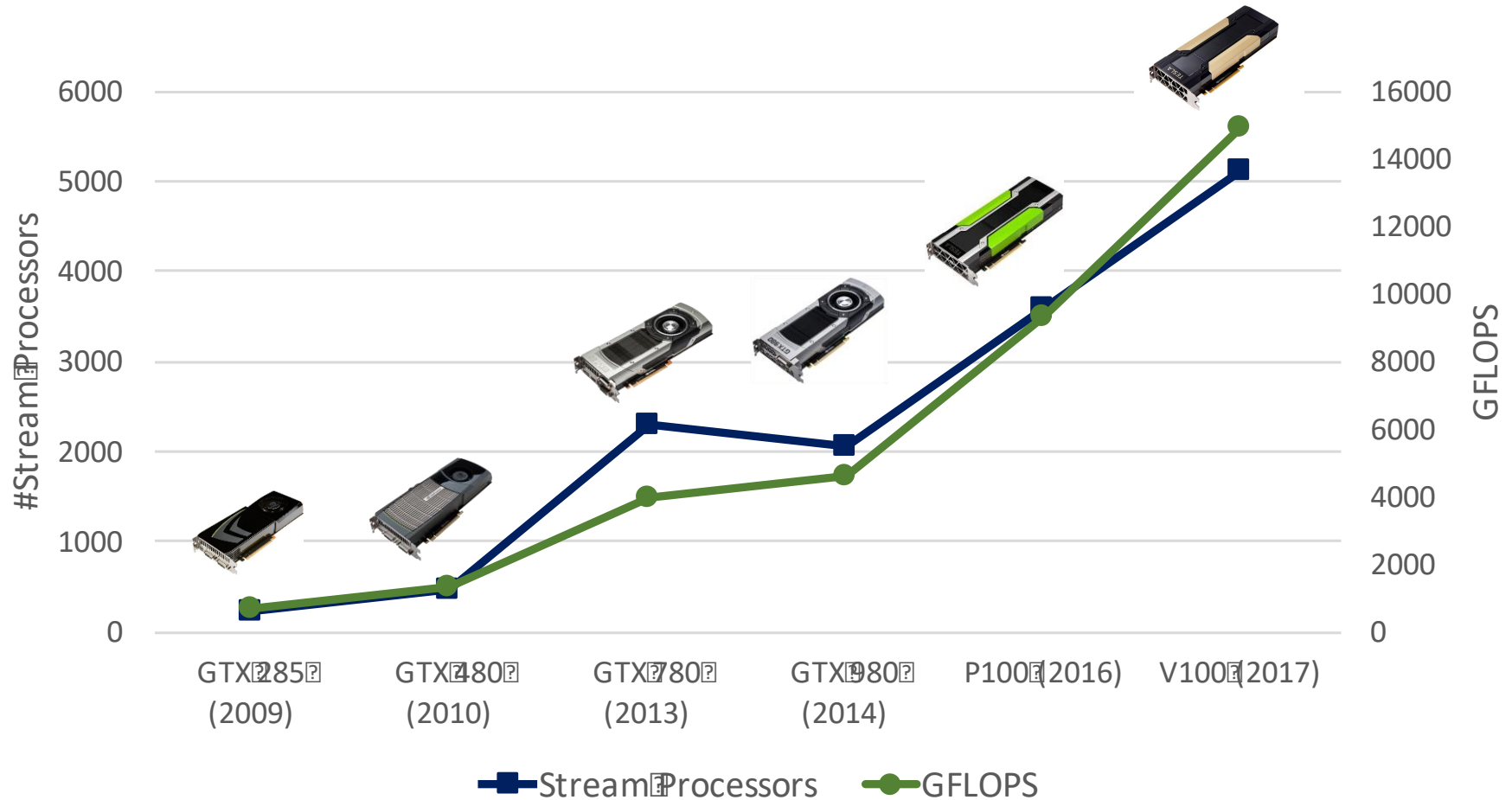
# NVIDIA GeForce GTX 285



30 cores on the GTX 285: 30,720 threads



# Recall: Evolution of NVIDIA GPUs



# Recall: NVIDIA V100

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- NVIDIA-speak:
  - 5120 stream processors
  - “SIMT execution”
  
- Generic speak:
  - 80 cores
  - 64 SIMD functional units per core
  
  - Specialized Functional Units for Machine Learning (tensor “cores” in NVIDIA-speak)



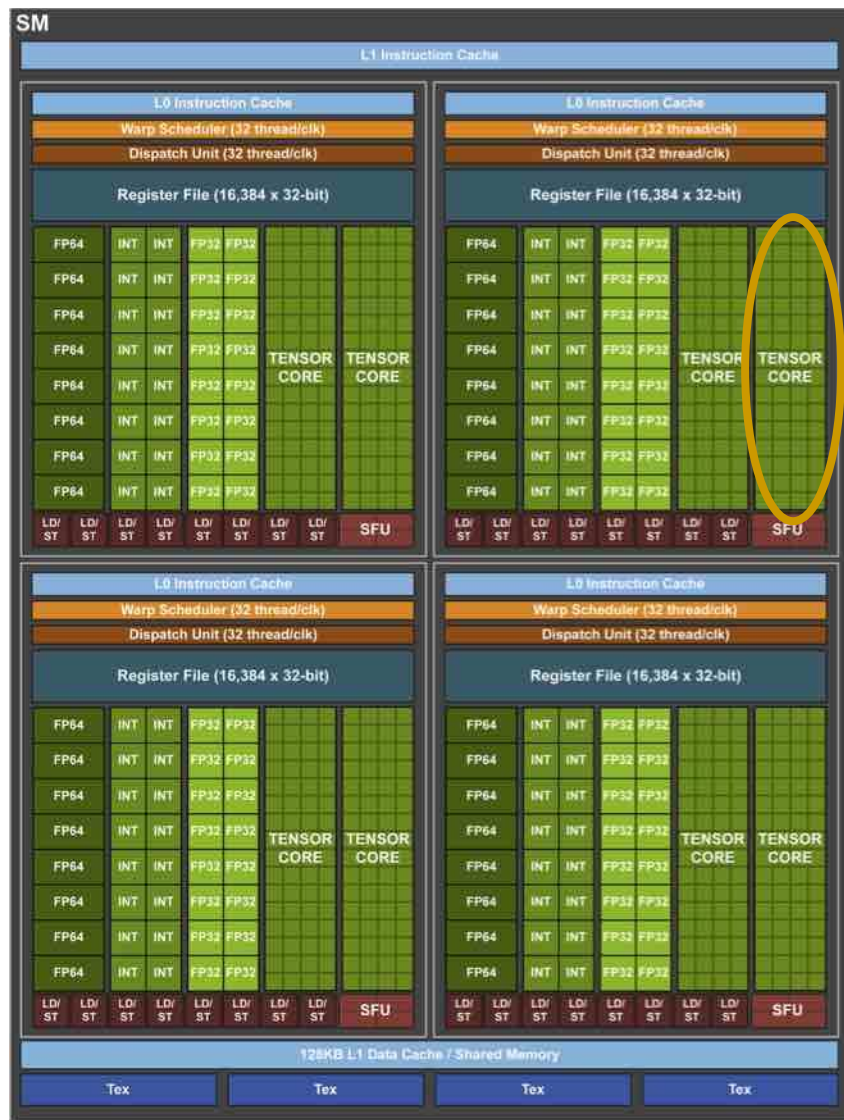
# Recall: NVIDIA V100 Block Diagram



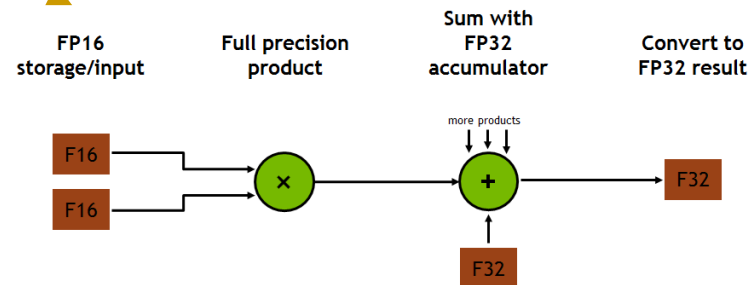
<https://devblogs.nvidia.com/inside-volta/>

80 cores on the V100

# Recall: NVIDIA V100 Core



15.7 TFLOPS Single Precision  
 7.8 TFLOPS Double Precision  
 125 TFLOPS for Deep Learning (Tensor "cores")



$$D = \begin{pmatrix} A_{0,0} & A_{0,1} & A_{0,2} & A_{0,3} \\ A_{1,0} & A_{1,1} & A_{1,2} & A_{1,3} \\ A_{2,0} & A_{2,1} & A_{2,2} & A_{2,3} \\ A_{3,0} & A_{3,1} & A_{3,2} & A_{3,3} \end{pmatrix} \begin{pmatrix} B_{0,0} & B_{0,1} & B_{0,2} & B_{0,3} \\ B_{1,0} & B_{1,1} & B_{1,2} & B_{1,3} \\ B_{2,0} & B_{2,1} & B_{2,2} & B_{2,3} \\ B_{3,0} & B_{3,1} & B_{3,2} & B_{3,3} \end{pmatrix} + \begin{pmatrix} C_{0,0} & C_{0,1} & C_{0,2} & C_{0,3} \\ C_{1,0} & C_{1,1} & C_{1,2} & C_{1,3} \\ C_{2,0} & C_{2,1} & C_{2,2} & C_{2,3} \\ C_{3,0} & C_{3,1} & C_{3,2} & C_{3,3} \end{pmatrix}$$

FP16 or FP32      FP16      FP16      FP16 or FP32

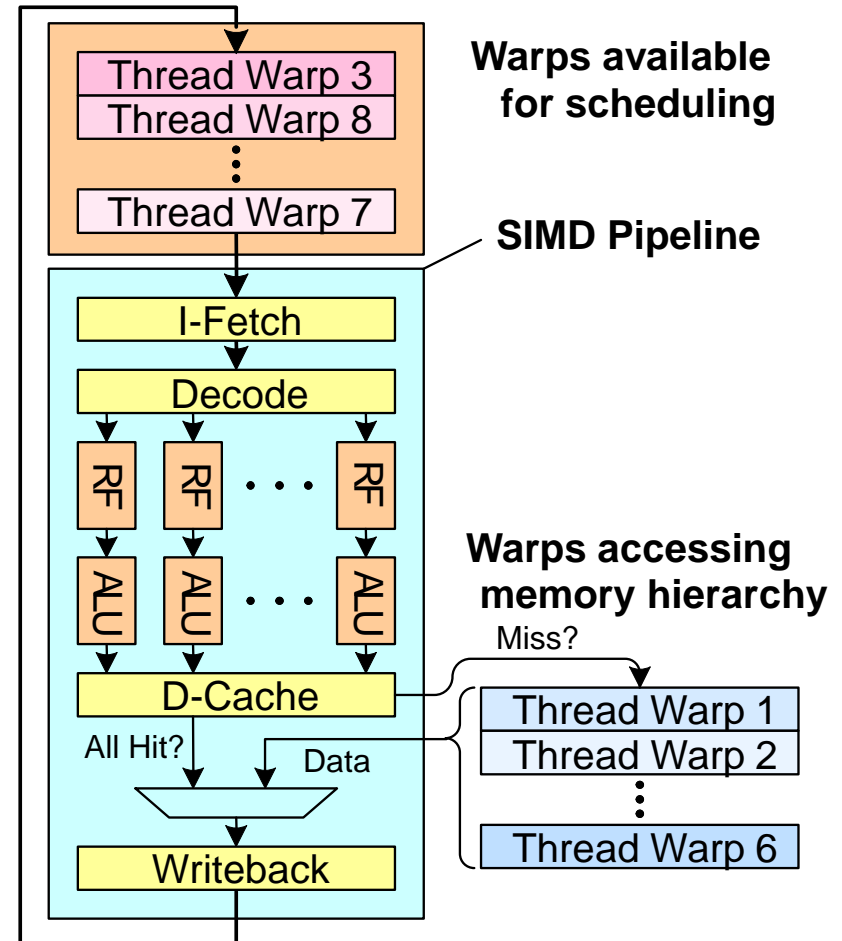
# Food for Thought

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- What is the main bottleneck in GPU programs?
- “Tensor cores”:
  - Can you think about other operations than matrix multiplication?
  - What other applications could benefit from specialized cores?
- Compare and contrast GPUs vs other accelerators (e.g., systolic arrays)
  - Which one is better for machine learning?
  - Which one is better for image/vision processing?
  - What types of parallelism each one exploits?
  - What are the tradeoffs?

# Recall: Latency Hiding via Warp-Level FGMT

- Warp: A set of threads that execute the same instruction (on different data elements)
- Fine-grained multithreading
  - One instruction per thread in pipeline at a time (No interlocking)
  - Interleave warp execution to hide latencies
- Register values of all threads stay in register file
- FGMT enables long latency tolerance
  - Millions of pixels

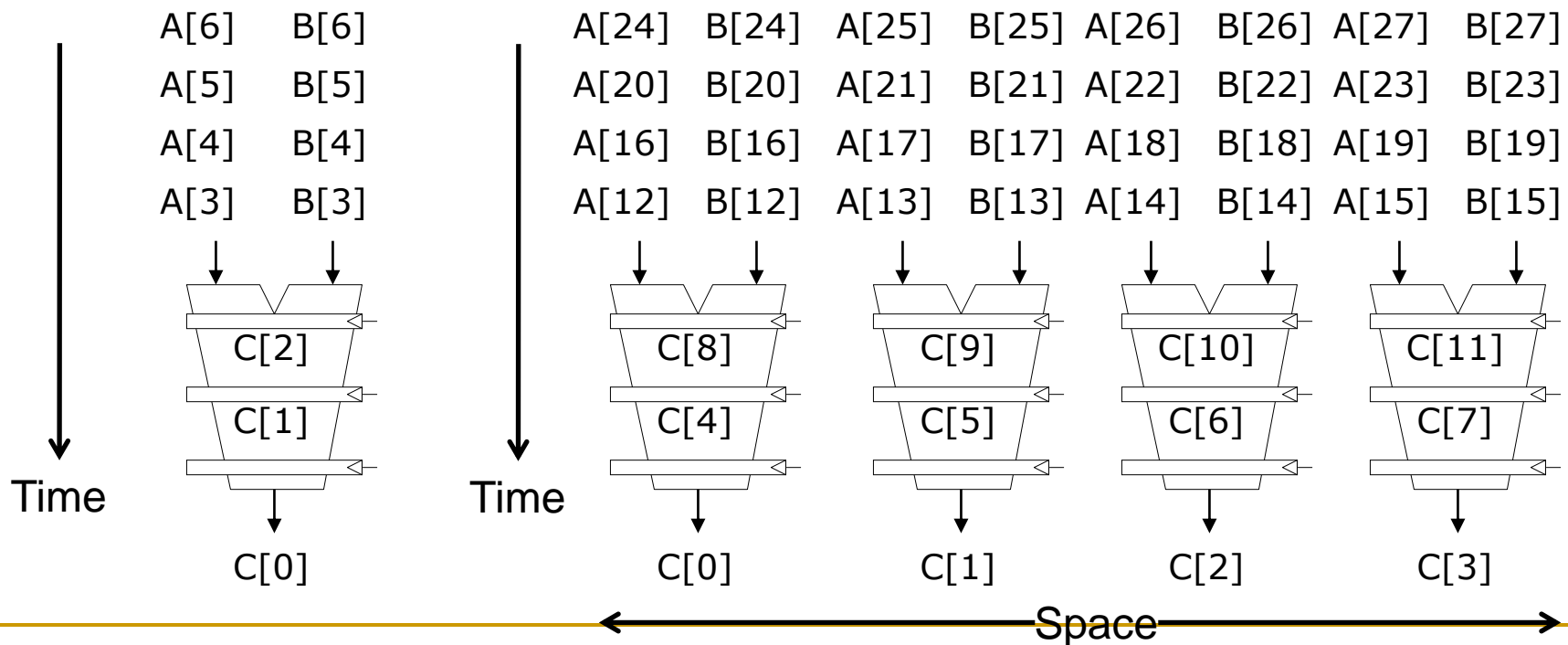


# Recall: Warp Execution

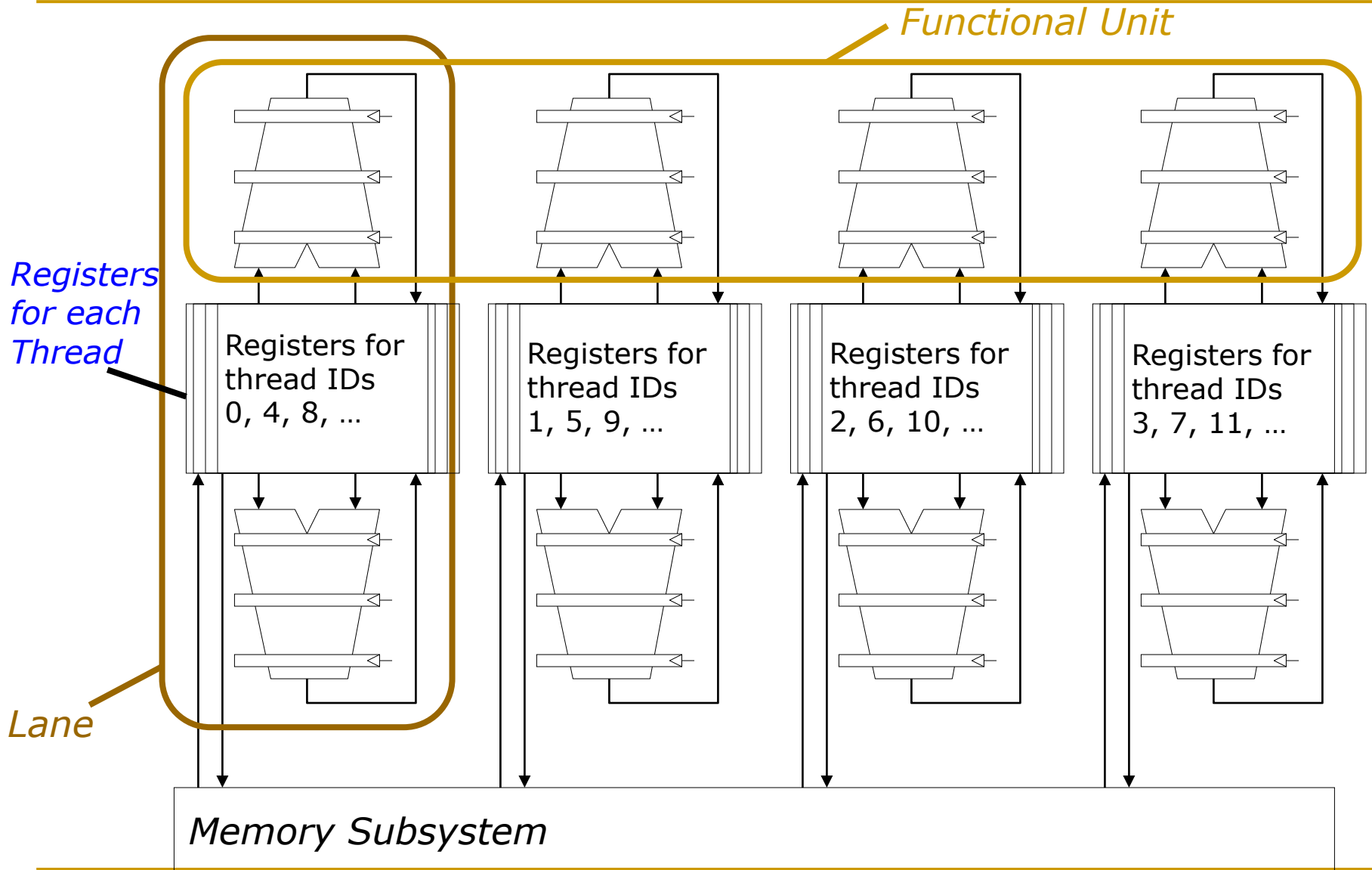
32-thread warp executing  $\text{ADD } A[\text{tid}], B[\text{tid}] \rightarrow C[\text{tid}]$

*Execution using  
one pipelined  
functional unit*

*Execution using  
four pipelined  
functional units*



# Recall: SIMD Execution Unit Structure

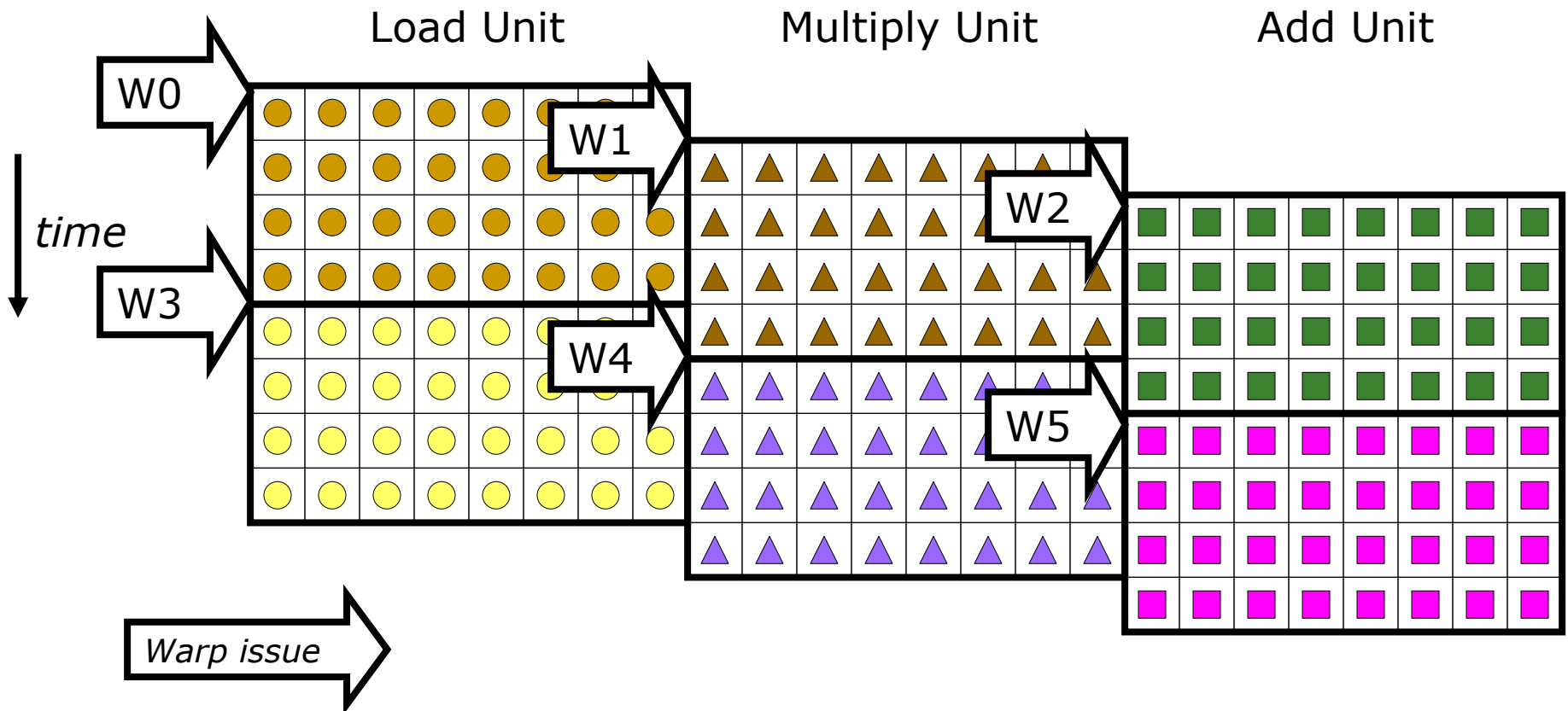




# Recall: Warp Instruction Level Parallelism

Can overlap execution of multiple instructions

- Example machine has 32 threads per warp and 8 lanes
- Completes 24 operations/cycle while issuing 1 warp/cycle



# Clarification of some GPU Terms

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Generic Term	NVIDIA Term	AMD Term	Comments
Vector length	Warp size	Wavefront size	Number of threads that run in parallel (lock-step) on a SIMD functional unit
Pipelined functional unit / Scalar pipeline	Streaming processor / CUDA core	-	Functional unit that executes instructions for one GPU thread
SIMD functional unit / SIMD pipeline	Group of N streaming processors (e.g., N=8 in GTX 285, N=16 in Fermi)	Vector ALU	SIMD functional unit that executes instructions for an entire warp
GPU core	Streaming multiprocessor	Compute unit	It contains one or more warp schedulers and one or several SIMD pipelines

# GPU Programming

# Recall: Vector Processor Disadvantages

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- Works (only) if parallelism is regular (data/SIMD parallelism)
  - ++ Vector operations
  - Very inefficient if parallelism is irregular
    - How about searching for a key in a linked list?

To program a vector machine, the compiler or hand coder must make the data structures in the code fit nearly exactly the regular structure built into the hardware. That's hard to do in first place, and just as hard to change. One tweak, and the low-level code has to be rewritten by a very smart and dedicated programmer who knows the hardware and often the subtleties of the application area. Often the rewriting is

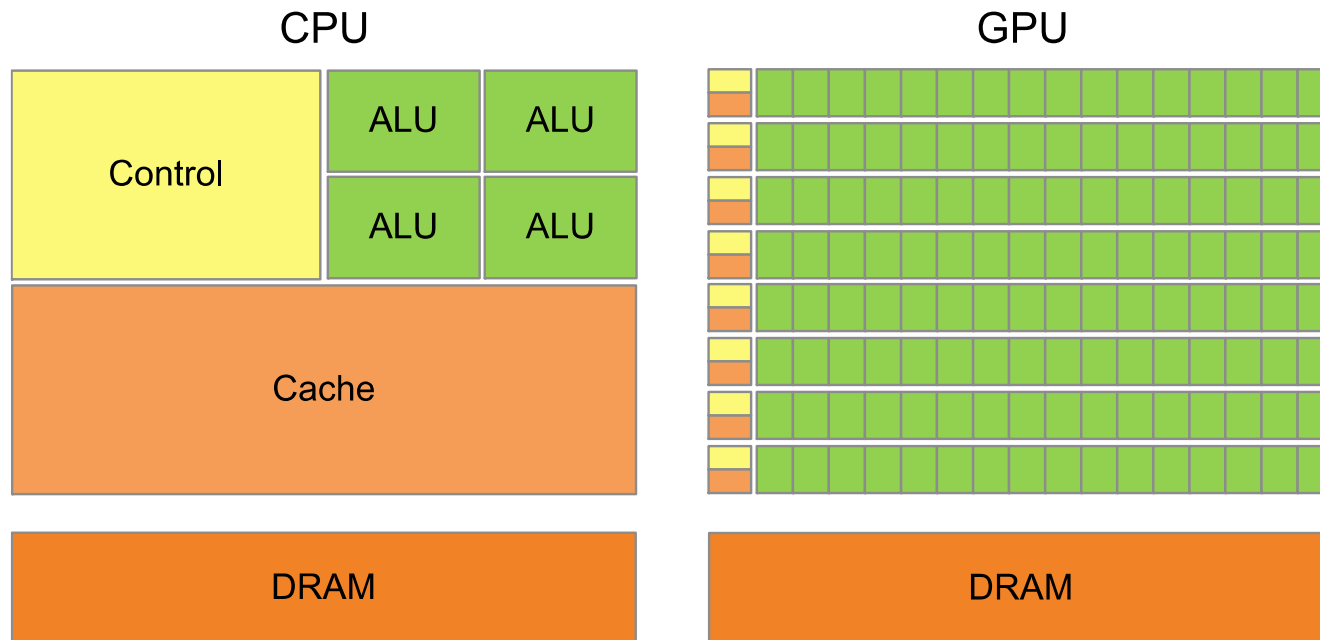
# General Purpose Processing on GPU

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- Easier programming of SIMD processors with SPMD
  - GPUs have democratized High Performance Computing (HPC)
  - Great FLOPS/\$, massively parallel chip on a commodity PC
- Many workloads exhibit inherent parallelism
  - Matrices
  - Image processing
  - Deep neural networks
- However, this is not for free
  - New programming model
  - Algorithms need to be re-implemented and rethought
- Still some bottlenecks
  - CPU-GPU data transfers (PCIe, NVLINK)
  - DRAM memory bandwidth (GDDR5, GDDR6, HBM2)
    - Data layout

# CPU vs. GPU

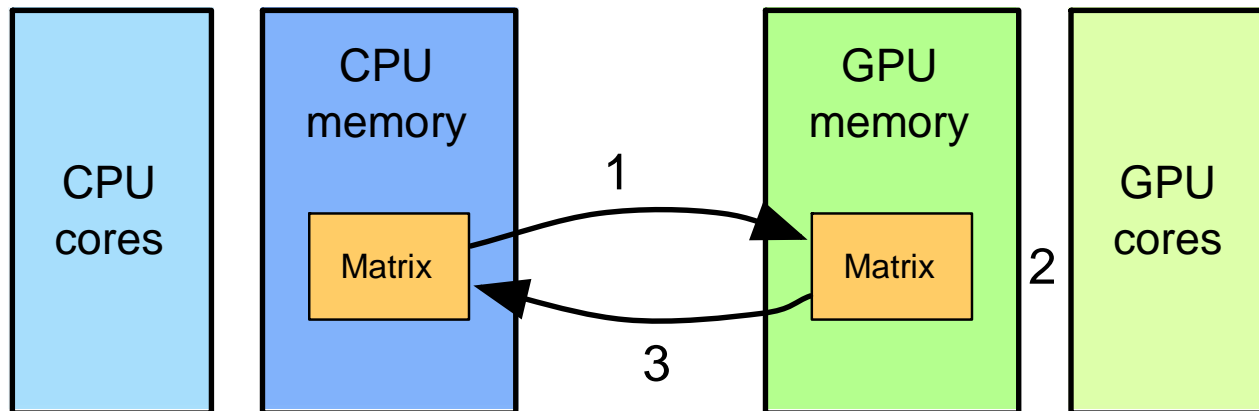
- Different design philosophies
  - ❑ CPU: A few out-of-order cores
  - ❑ GPU: Many in-order FGMT cores



# GPU Computing

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- Computation is **offloaded to the GPU**
- Three steps
  - ❑ CPU-GPU data transfer (1)
  - ❑ GPU kernel execution (2)
  - ❑ GPU-CPU data transfer (3)



# Traditional Program Structure

- CPU threads and GPU kernels
  - ▣ Sequential or modestly parallel sections on CPU
  - ▣ Massively parallel sections on GPU

Serial Code (host)

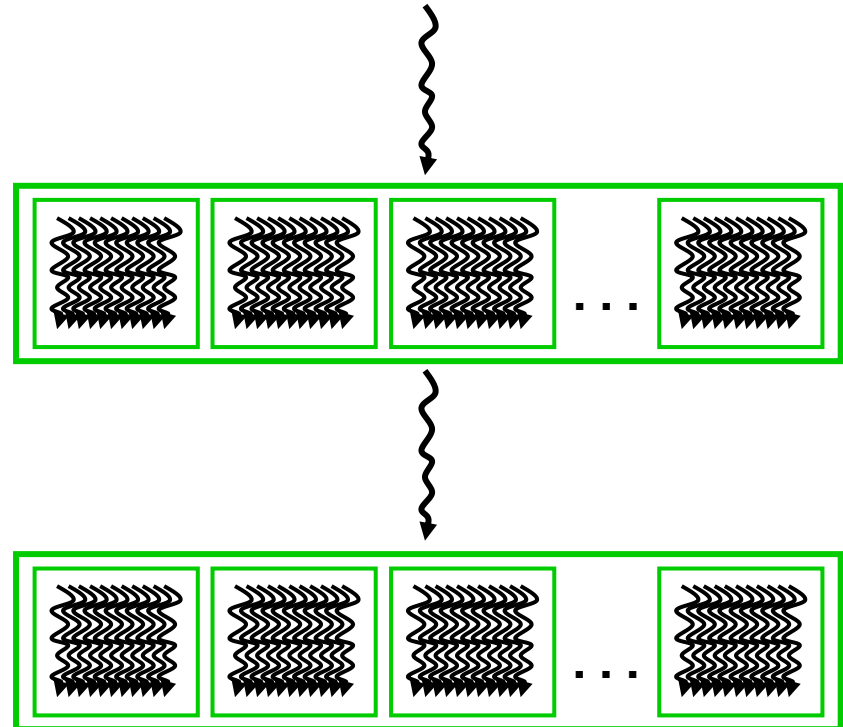
Parallel Kernel (device)

```
KernelA<<< nBlk, nThr >>>(args);
```

Serial Code (host)

Parallel Kernel (device)

```
KernelB<<< nBlk, nThr >>>(args);
```





# Recall: SPMD

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- Single procedure/program, multiple data
  - This is a programming model rather than computer organization
- Each processing element executes the same procedure, except on different data elements
  - Procedures can synchronize at certain points in program, e.g. barriers
- Essentially, multiple instruction streams execute the same program
  - Each program/procedure 1) works on different data, 2) can execute a different control-flow path, at run-time
  - Many scientific applications are programmed this way and run on MIMD hardware (multiprocessors)
  - Modern GPUs programmed in a similar way on a SIMD hardware

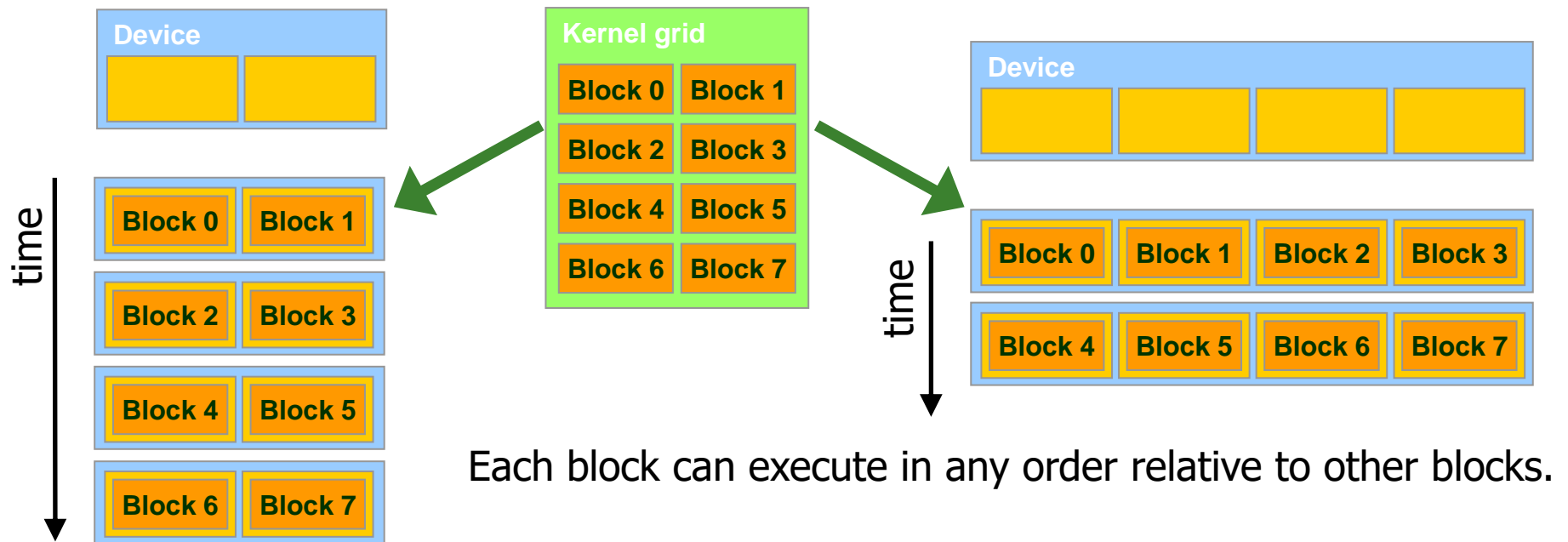
# CUDA/OpenCL Programming Model

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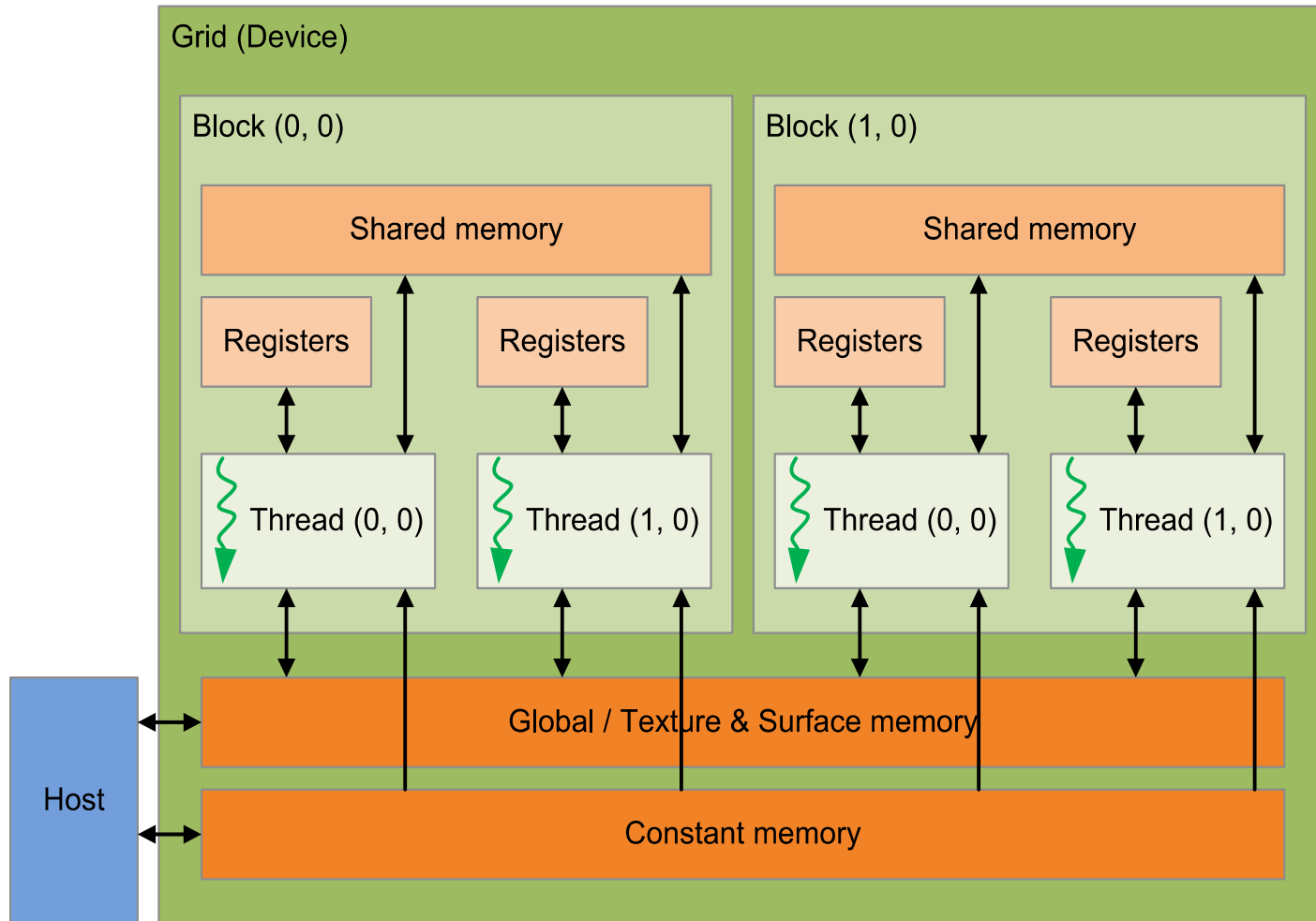
- SIMT or SPMD
- Bulk synchronous programming
  - Global (coarse-grain) synchronization between kernels
- The host (typically CPU) allocates memory, copies data, and launches kernels
- The device (typically GPU) executes kernels
  - Grid (NDRange)
  - Block (work-group)
    - Within a block, shared memory, and synchronization
  - Thread (work-item)

# Transparent Scalability

- Hardware is **free to schedule** thread blocks



# Memory Hierarchy



# Traditional Program Structure in CUDA


## ■ Function prototypes

```
float serialFunction(...);  
__global__ void kernel(...);
```

## ■ main()

- ❑ 1) **Allocate memory** space on the device – `cudaMalloc(&d_in, bytes);`
- ❑ 2) Transfer data from **host to device** – `cudaMemcpy(d_in, h_in, ...);`
- ❑ 3) Execution configuration setup: **#blocks** and **#threads**
- ❑ 4) **Kernel call** – `kernel<<<execution configuration>>>(args...);`
- ❑ 5) Transfer results from **device to host** – `cudaMemcpy(h_out, d_out, ...);`

repeat  
as needed



## ■ Kernel – `__global__ void kernel(type args,...)`

- ❑ Automatic variables transparently assigned to **registers**
- ❑ **Shared memory**: `__shared__`
- ❑ Intra-block **synchronization**: `__syncthreads();`

# CUDA Programming Language

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- **Memory allocation**

```
cudaMalloc((void**)&d_in, #bytes);
```

- **Memory copy**

```
cudaMemcpy(d_in, h_in, #bytes, cudaMemcpyHostToDevice);
```

- **Kernel launch**

```
kernel<<< #blocks, #threads >>>(args);
```

- **Memory deallocation**

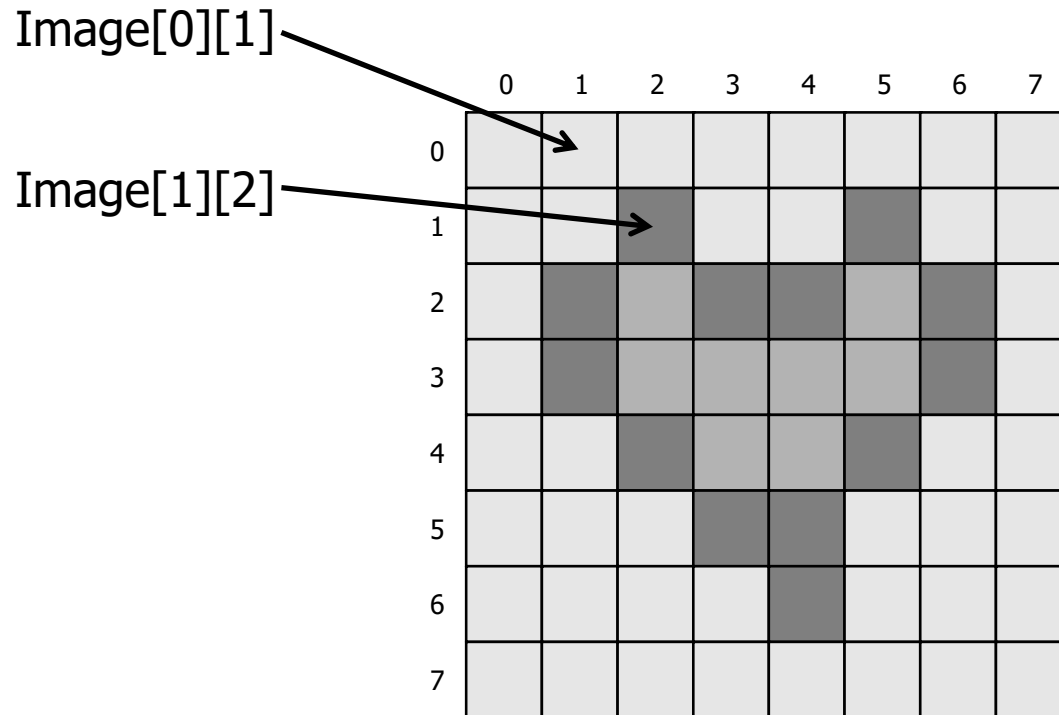
```
cudaFree(d_in);
```

- **Explicit synchronization**

```
cudaDeviceSynchronize();
```

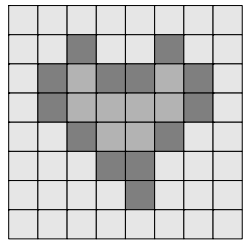
# Indexing and Memory Access

- Images are 2D data structures
  - height x width
  - $\text{Image}[j][i]$ , where  $0 \leq j < \text{height}$ , and  $0 \leq i < \text{width}$



# Image Layout in Memory

- Row-major layout
- $\text{Image}[j][i] = \text{Image}[j \times \text{width} + i]$



Stride = width

$$\text{Image}[0][1] = \text{Image}[0 \times 8 + 1]$$

$$\text{Image}[1][2] = \text{Image}[1 \times 8 + 2]$$



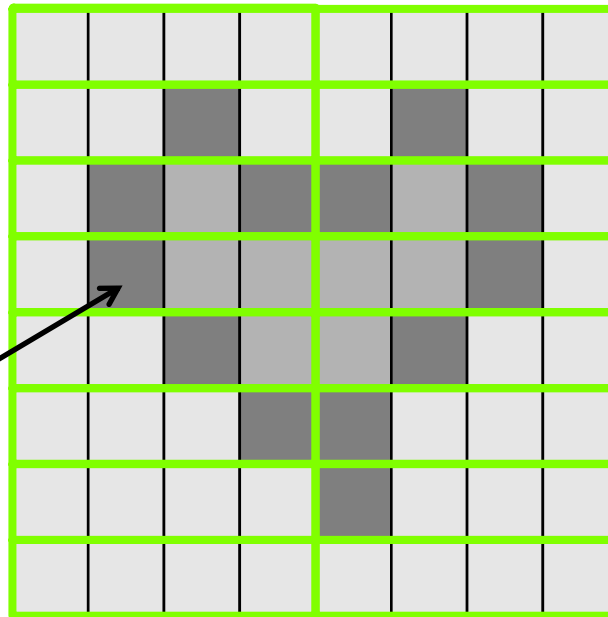
# Indexing and Memory Access: 1D Grid

- One GPU thread per pixel
- Grid of Blocks of Threads
  - `gridDim.x, blockDim.x`
  - `blockIdx.x, threadIdx.x`

`blockIdx.x`

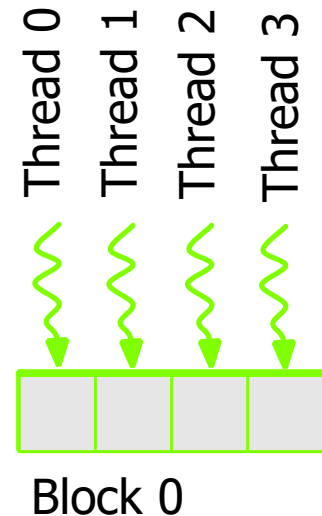
`threadIdx.x`

Block 0



$$6 * 4 + 1 = 25$$

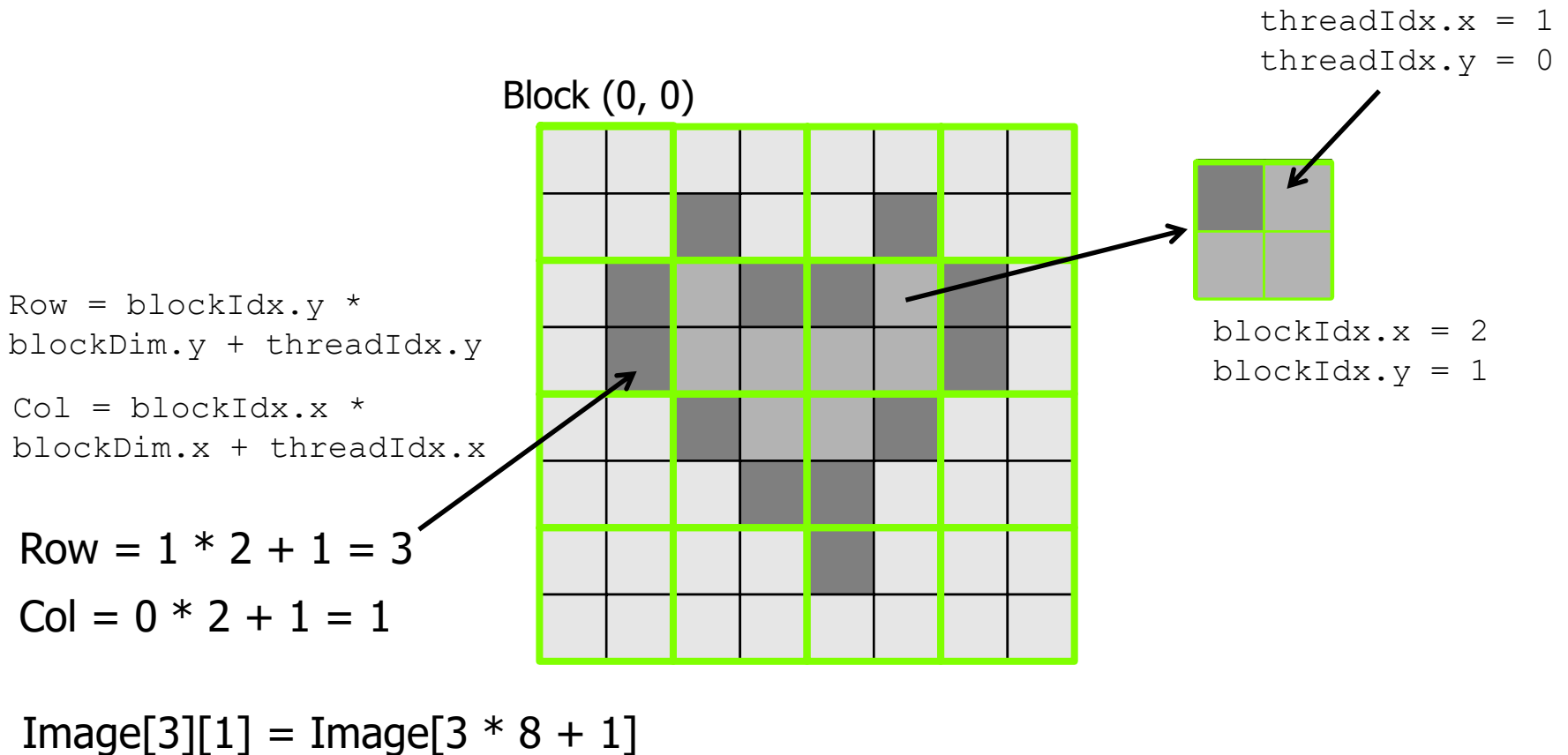
`blockIdx.x * blockDim.x + threadIdx.x`



# Indexing and Memory Access: 2D Grid

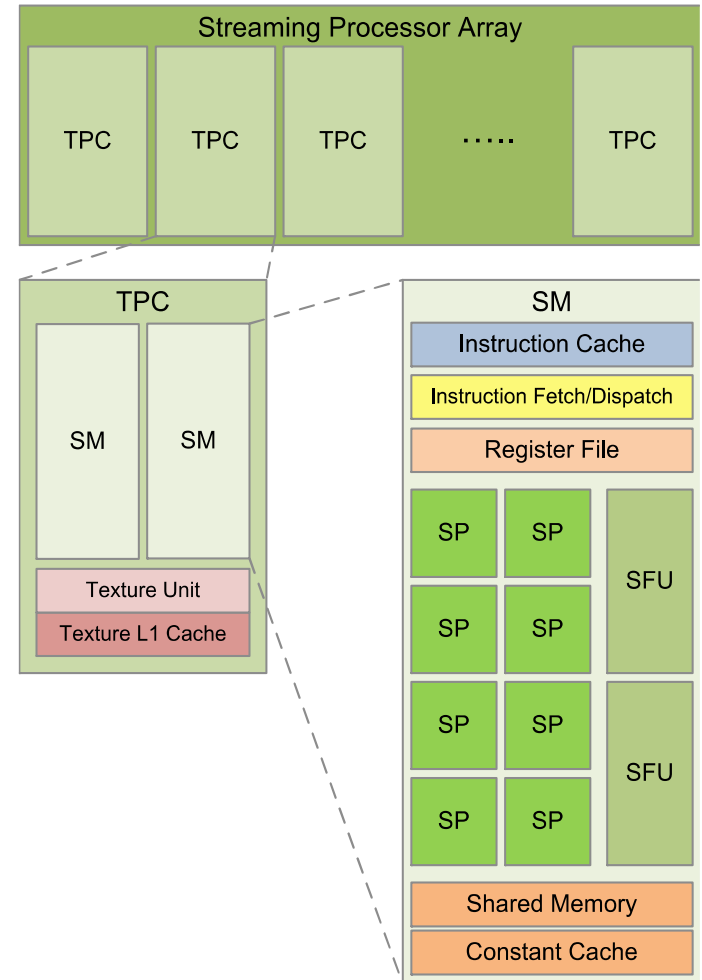
## ■ 2D blocks

□ `gridDim.x`, `gridDim.y`



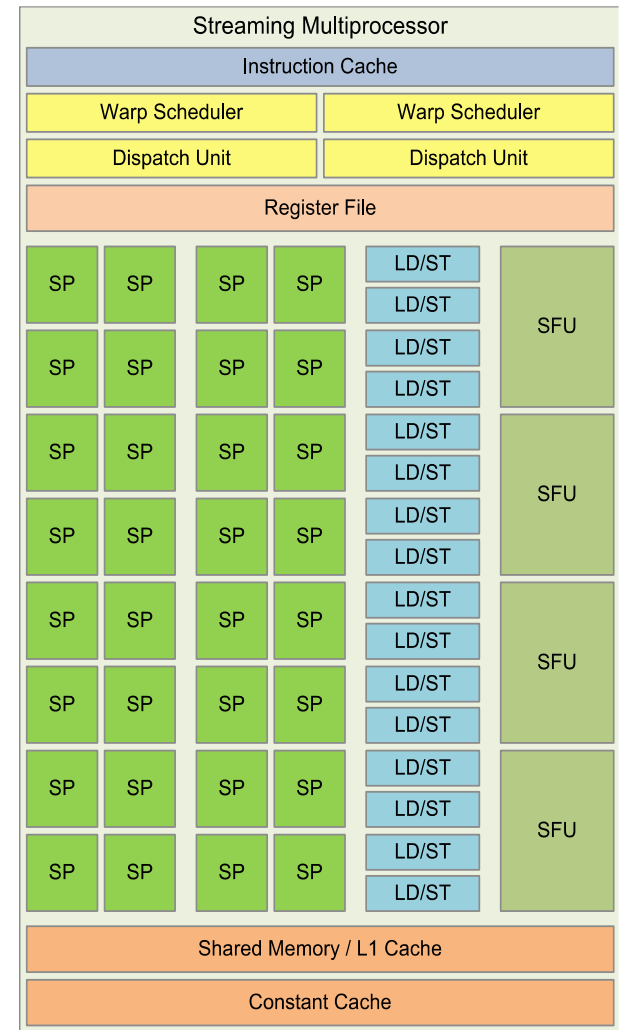
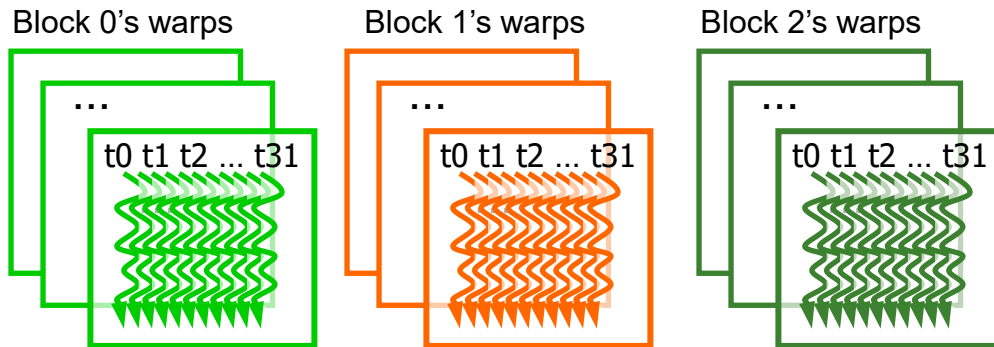
# Brief Review of GPU Architecture (I)

- Streaming Processor Array
  - Tesla architecture (G80/GT200)



# Brief Review of GPU Architecture (II)

- Streaming Multiprocessors (SM)
  - Streaming Processors (SP)
- Blocks are divided into warps
  - SIMD unit (32 threads)



NVIDIA Fermi architecture

# Brief Review of GPU Architecture (III)

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- Streaming Multiprocessors (SM) or Compute Units (CU)
  - SIMD pipelines
- Streaming Processors (SP) or CUDA "cores"
  - Vector lanes
- Number of SMs x SPs across generations
  - Tesla (2007): 30 x 8
  - Fermi (2010): 16 x 32
  - Kepler (2012): 15 x 192
  - Maxwell (2014): 24 x 128
  - Pascal (2016): 56 x 64
  - Volta (2017): 80 x 64

# Performance Considerations

# Performance Considerations

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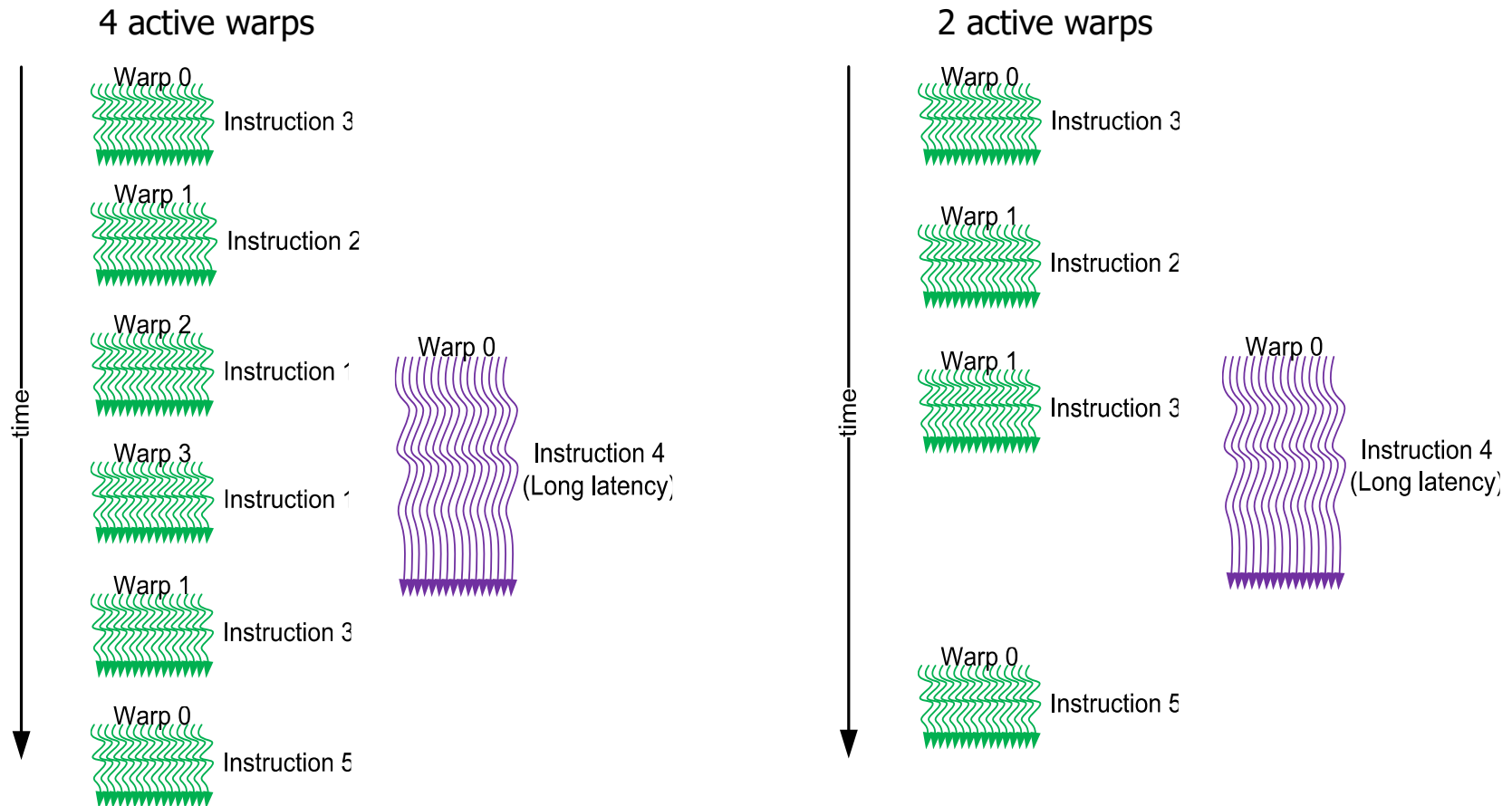
- Main bottlenecks
  - ❑ Global memory access
  - ❑ CPU-GPU data transfers
- Memory access
  - ❑ Latency hiding
    - Occupancy
  - ❑ Memory coalescing
  - ❑ Data reuse
    - Shared memory usage
- SIMD (Warp) Utilization: Divergence
- Atomic operations: Serialization
- Data transfers between CPU and GPU
  - ❑ Overlap of communication and computation

# Memory Access



# Latency Hiding

- FGMT can hide **long latency operations** (e.g., memory accesses)
- **Occupancy**: ratio of active warps



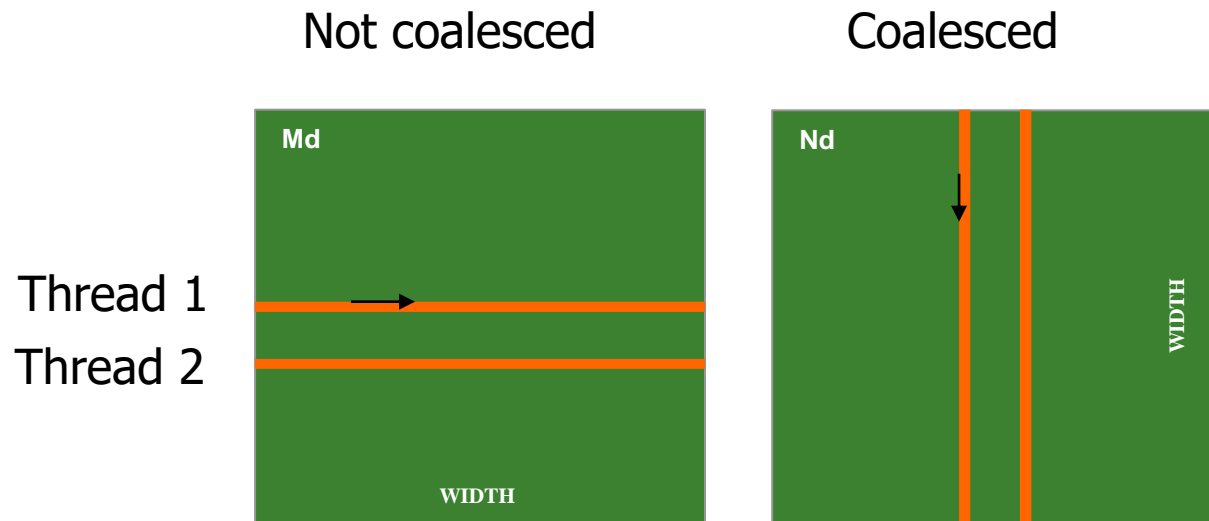
# Occupancy

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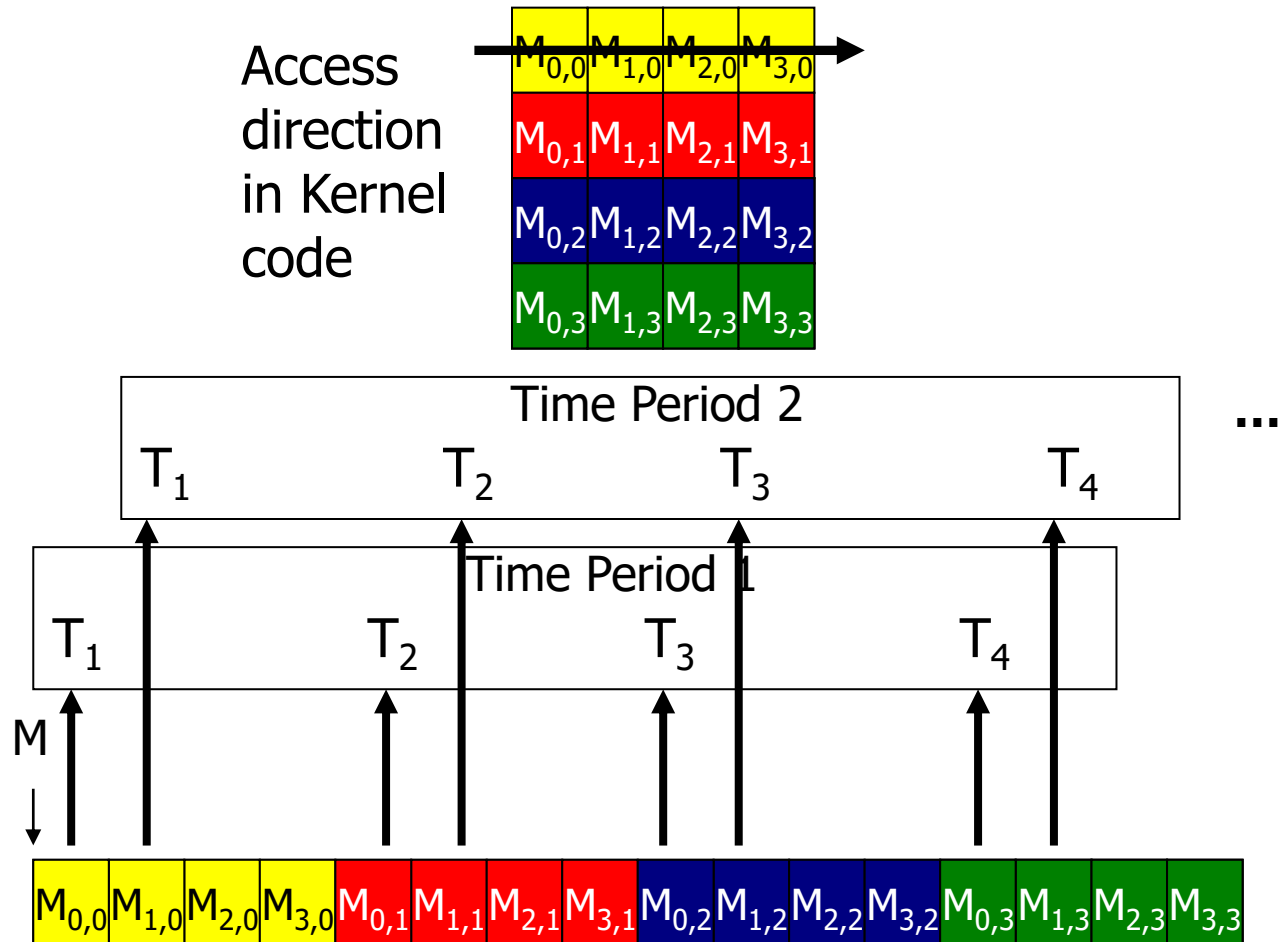
- SM resources (typical values)
  - ❑ Maximum number of warps per SM (64)
  - ❑ Maximum number of blocks per SM (32)
  - ❑ Register usage (256KB)
  - ❑ Shared memory usage (64KB)
- Occupancy calculation
  - ❑ Number of threads per block (defined by the programmer)
  - ❑ Registers per thread (known at compile time)
  - ❑ Shared memory per block (defined by the programmer)

# Memory Coalescing

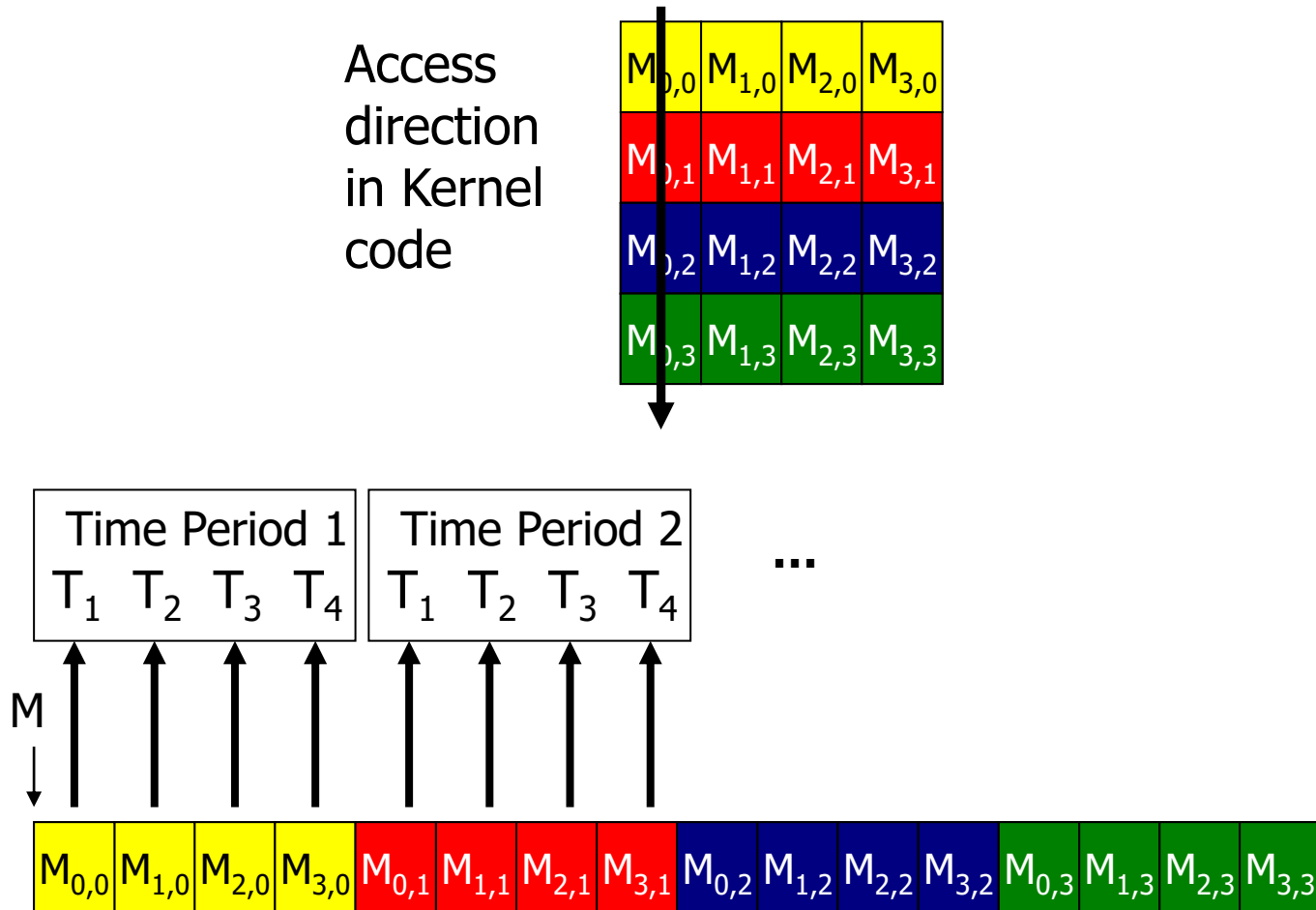
- When accessing global memory, we want to make sure that **concurrent threads access nearby memory locations**
- **Peak bandwidth** utilization occurs when all threads in a warp access **one cache line**



# Uncoalesced Memory Accesses



# Coalesced Memory Accesses

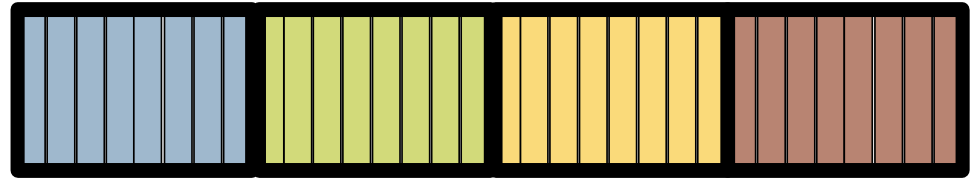


# AoS vs. SoA

## ■ Array of Structures vs. Structure of Arrays

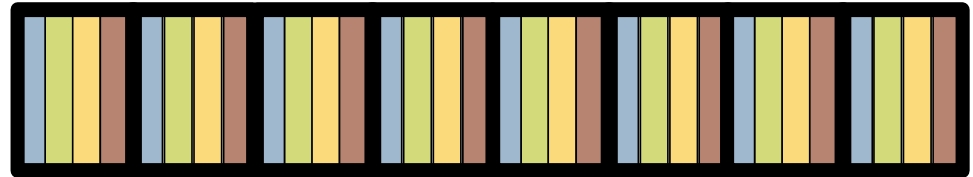
Structure of  
Arrays  
(SoA)

```
struct foo{  
    float a[8];  
    float b[8];  
    float c[8];  
    int d[8];  
} A;
```



Array of  
Structures  
(AoS)

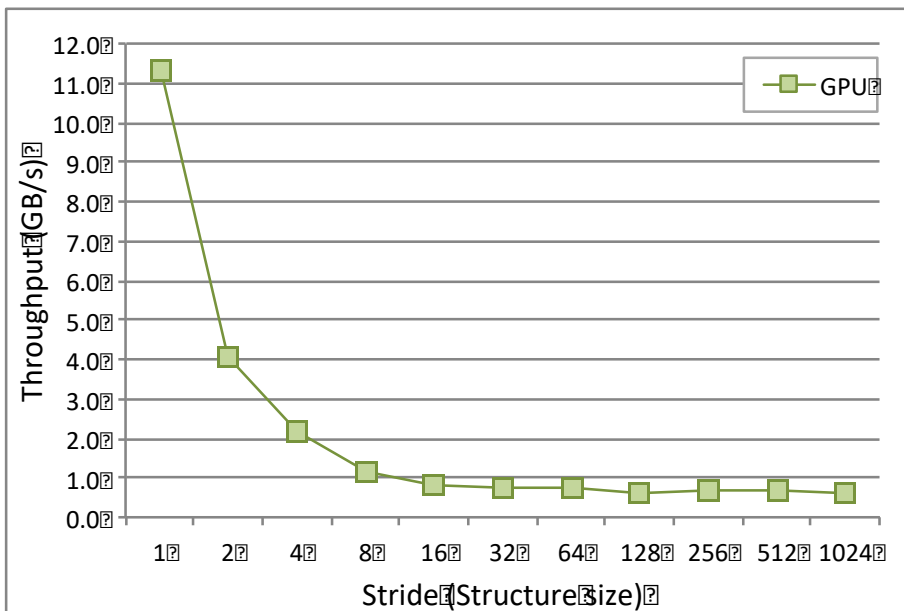
```
struct foo{  
    float a;  
    float b;  
    float c;  
    int d;  
} A[8];
```



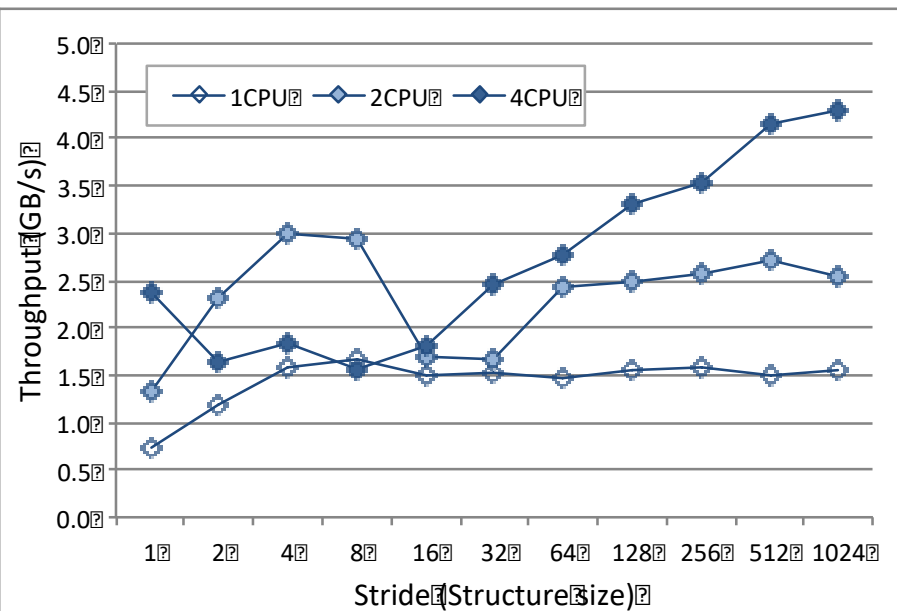
# CPUs Prefer AoS, GPUs Prefer SoA

## Linear and strided accesses

GPU



CPU

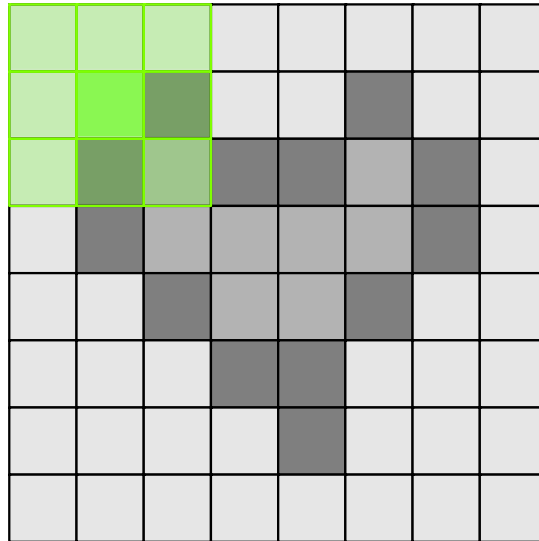


AMD Kaveri A10-7850K

Sung+, “[DL: A data layout transformation system for heterogeneous computing](#),” INPAR 2012

# Data Reuse

- Same memory locations accessed by neighboring threads

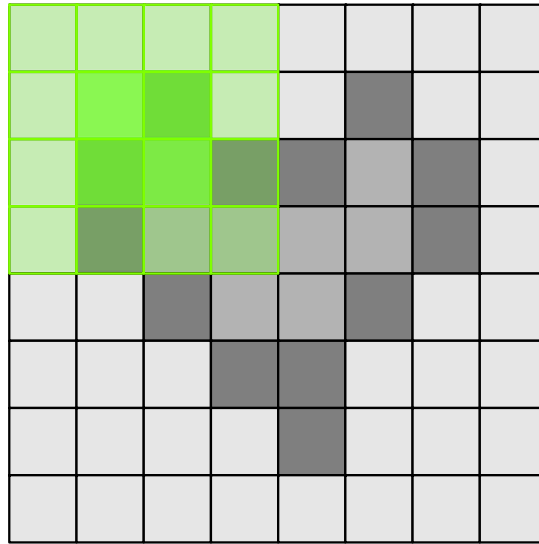


```
for (int i = 0; i < 3; i++){  
    for (int j = 0; j < 3; j++){  
        sum += gauss[i][j] * Image[(i+row-1)*width + (j+col-1)];  
    }  
}
```



# Data Reuse: Tiling

- To take advantage of data reuse, we divide the input into **tiles** that can be loaded into **shared memory**



```
__shared__ int l_data[(L_SIZE+2)*(L_SIZE+2)];  
...  
Load tile into shared memory  
__syncthreads();  
for (int i = 0; i < 3; i++){  
    for (int j = 0; j < 3; j++){  
        sum += gauss[i][j] * l_data[(i+l_row-1)*(L_SIZE+2)+j+l_col-1];  
    }  
}
```

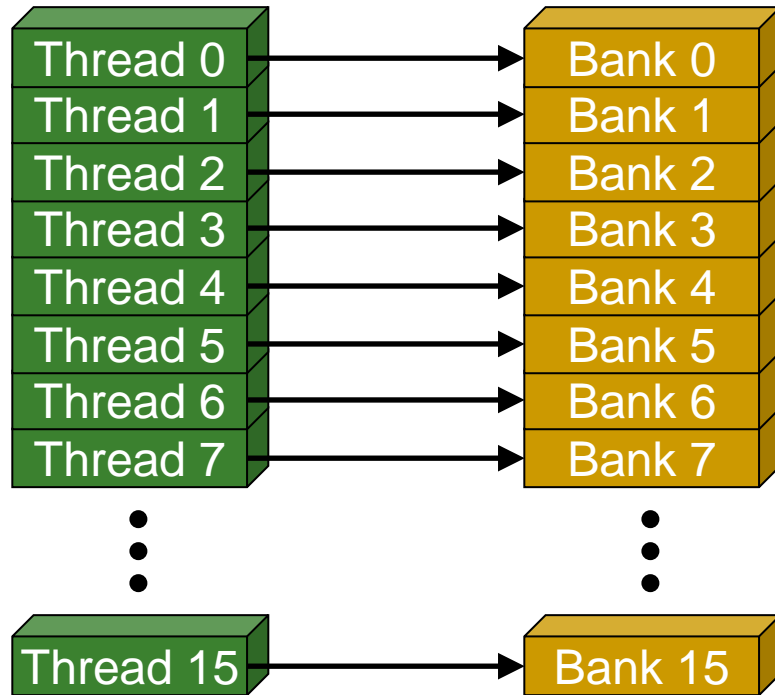
# Shared Memory

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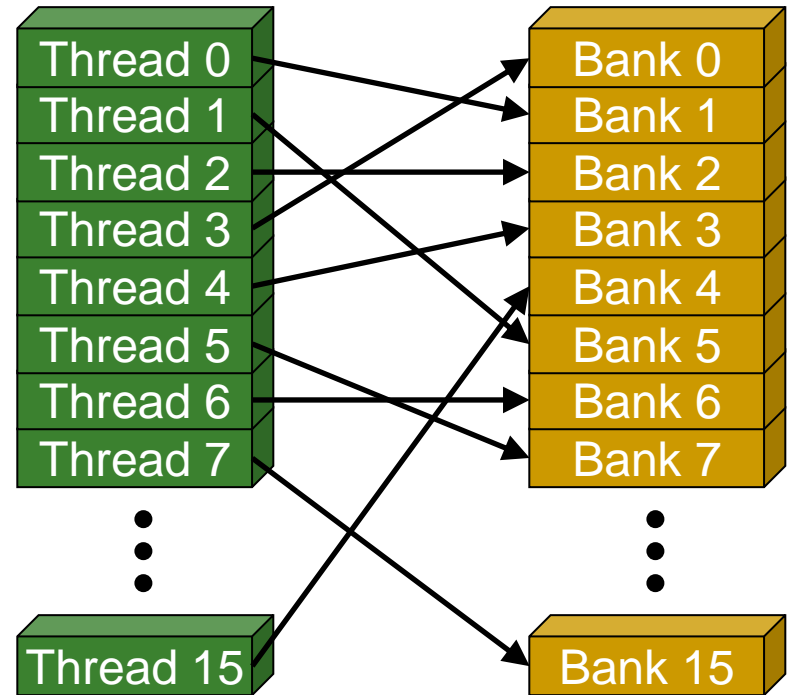
- Shared memory is an **interleaved (banked) memory**
  - Each bank can service one address per cycle
- Typically, 32 banks in NVIDIA GPUs
  - Successive 32-bit words are assigned to successive banks
    - $\text{Bank} = \text{Address} \% 32$
- Bank conflicts are **only possible within a warp**
  - No bank conflicts between different warps

# Shared Memory Bank Conflicts (I)

- Bank conflict free



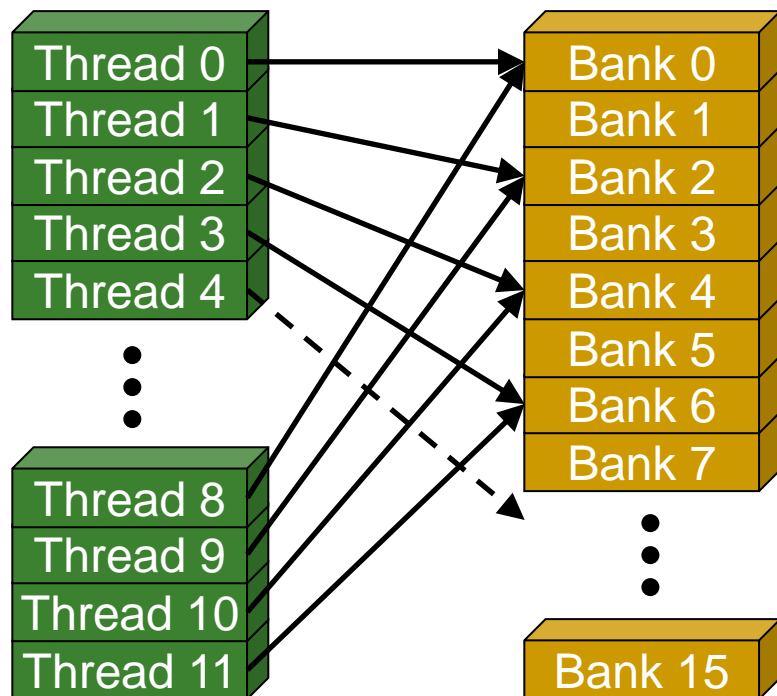
Linear addressing: stride = 1



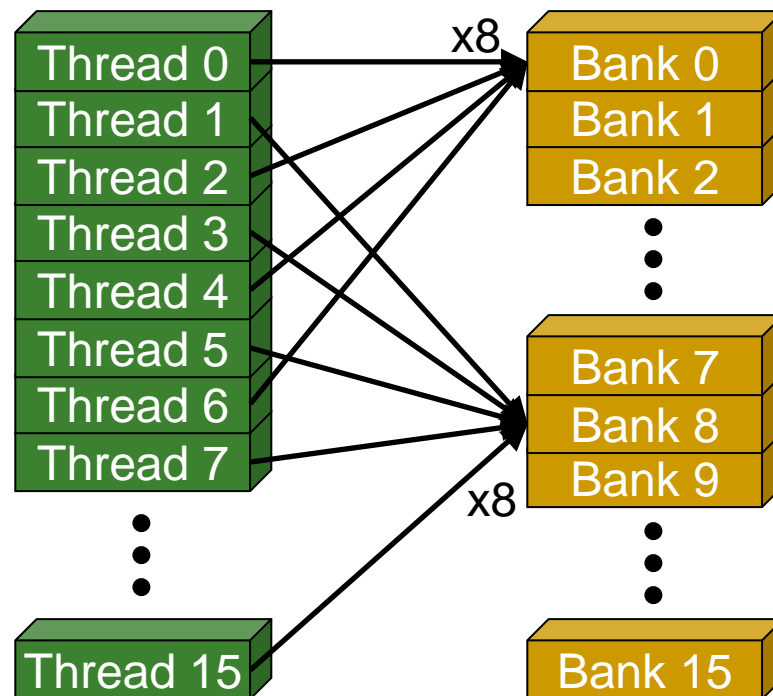
Random addressing 1:1

# Shared Memory Bank Conflicts (II)

## ■ N-way bank conflicts



2-way bank conflict: stride = 2



8-way bank conflict: stride = 8

# Reducing Shared Memory Bank Conflicts

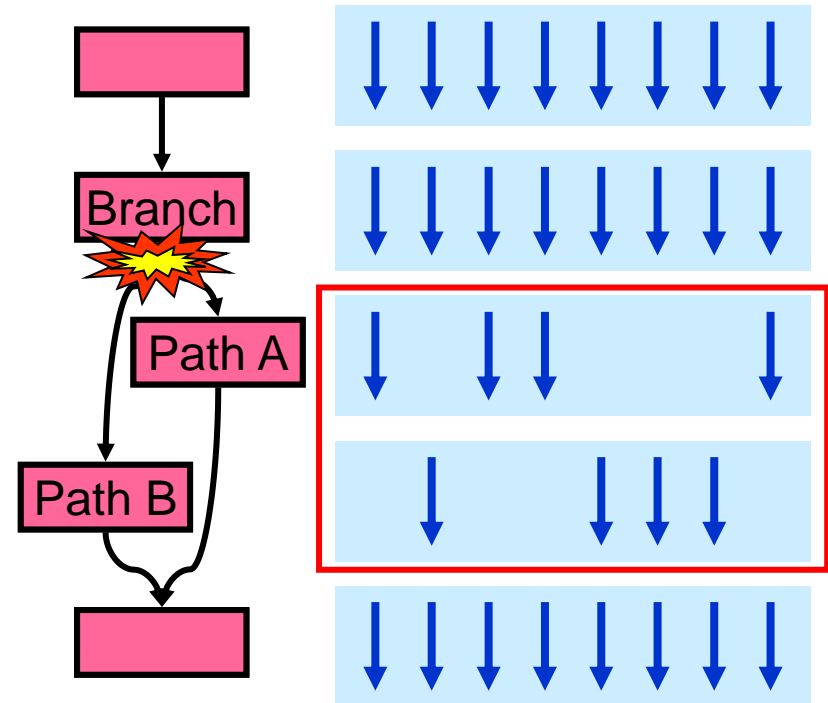
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- Bank conflicts are only possible within a warp
  - No bank conflicts between different warps
- If strided accesses are needed, some optimization techniques can help
  - Padding
  - Randomized mapping
    - Rau, “Pseudo-randomly interleaved memory,” ISCA 1991
  - Hash functions
    - V.d.Braak+, “Configurable XOR Hash Functions for Banked Scratchpad Memories in GPUs,” IEEE TC, 2016

# SIMD Utilization

# Control Flow Problem in GPUs/SIMT

- A GPU uses a SIMD pipeline to save area on control logic
  - Groups scalar threads into warps
- **Branch divergence** occurs when threads inside warps branch to different execution paths

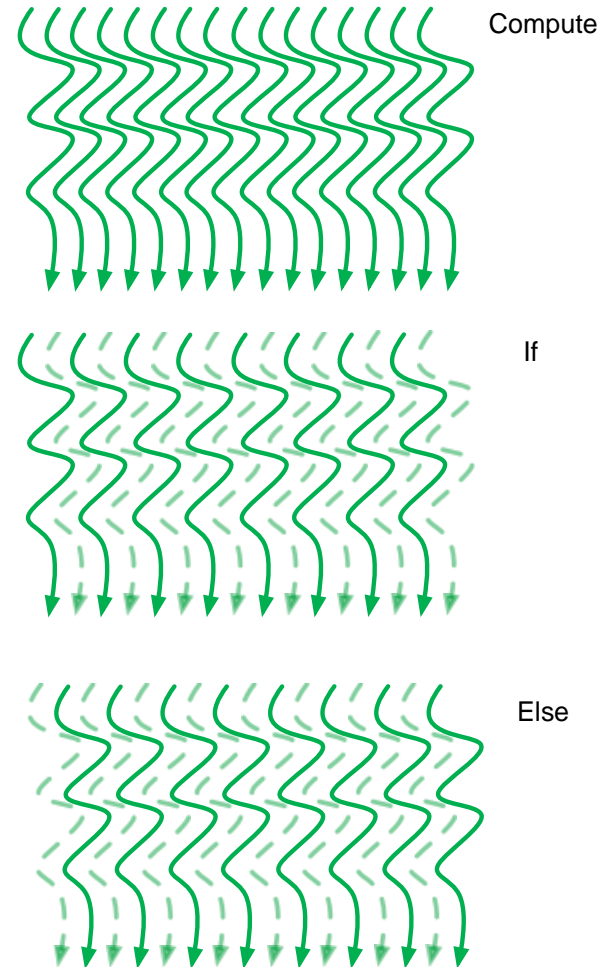


**This is the same as conditional/predicated/masked execution. Recall the Vector Mask and Masked Vector Operations?**

# SIMD Utilization

## ■ Intra-warp divergence

```
Compute(threadIdx.x);  
if (threadIdx.x % 2 == 0) {  
    Do_this(threadIdx.x);  
}  
else {  
    Do_that(threadIdx.x);  
}
```

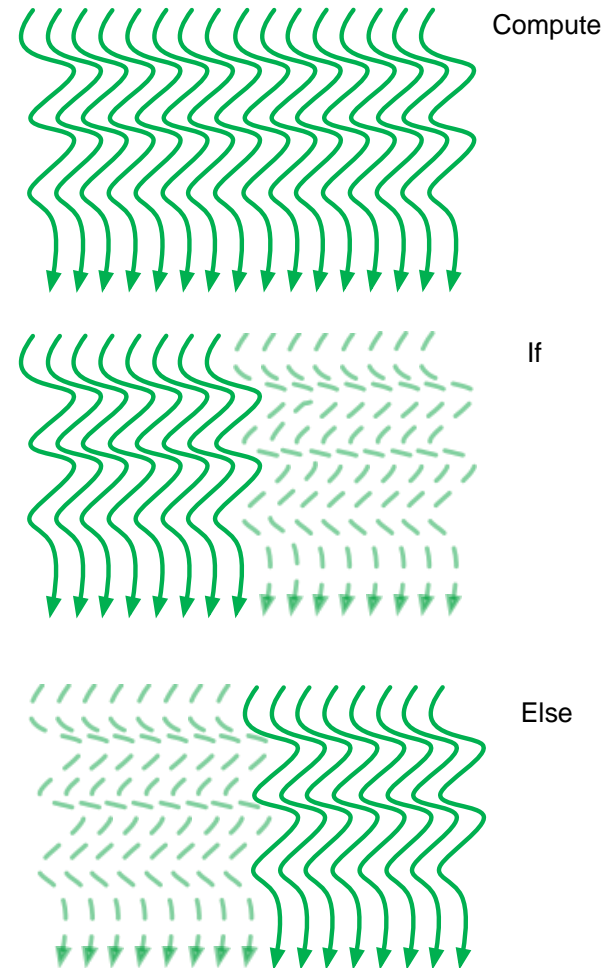




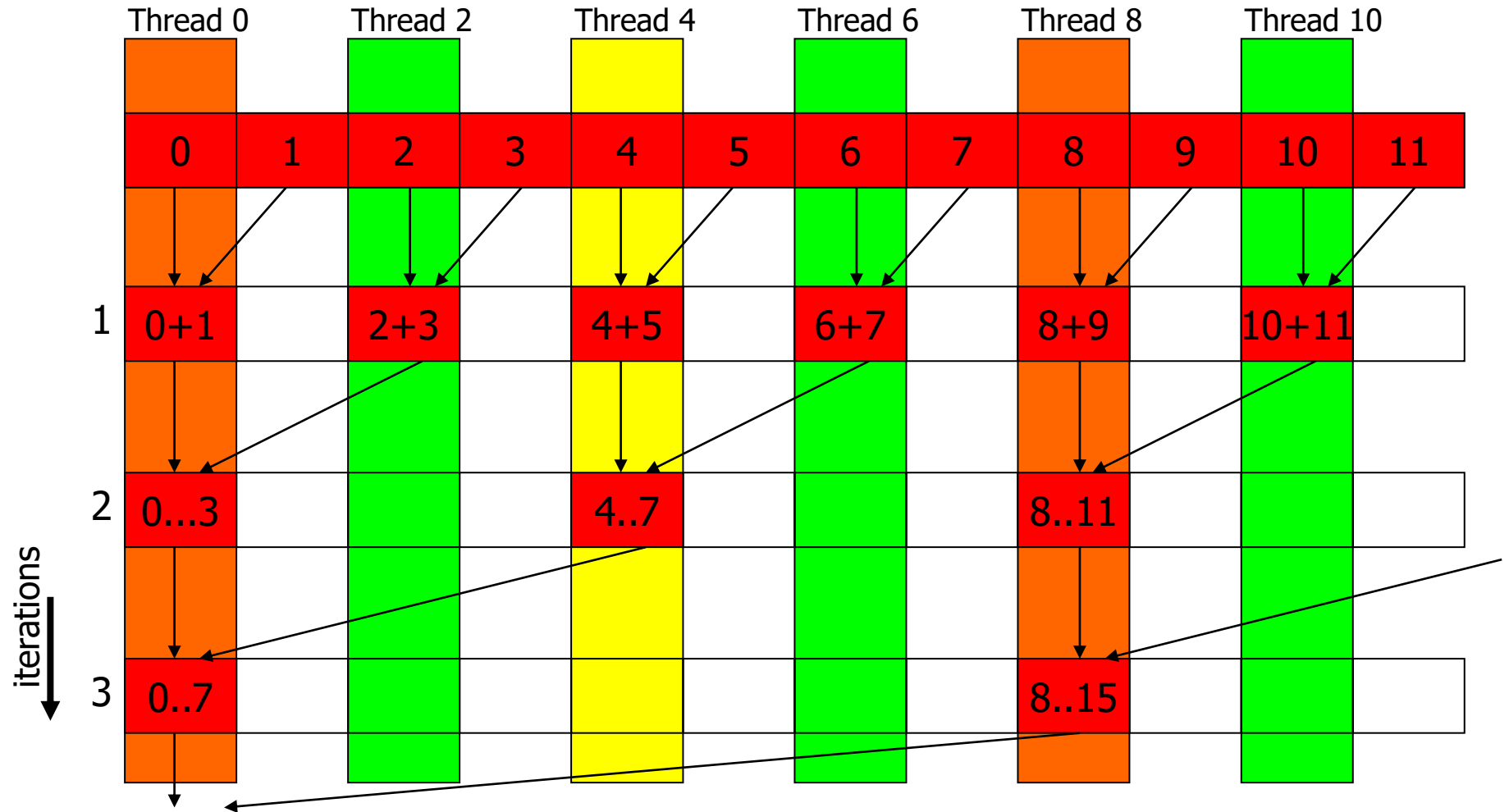
# Increasing SIMD Utilization

## ■ Divergence-free execution

```
Compute(threadIdx.x);  
if (threadIdx.x < 32){  
    Do_this(threadIdx.x * 2);  
}  
else{  
    Do_that((threadIdx.x%32)*2+1);  
}
```



# Vector Reduction: Naïve Mapping (I)



Slide credit: Hwu & Kirk

# Vector Reduction: Naïve Mapping (II)

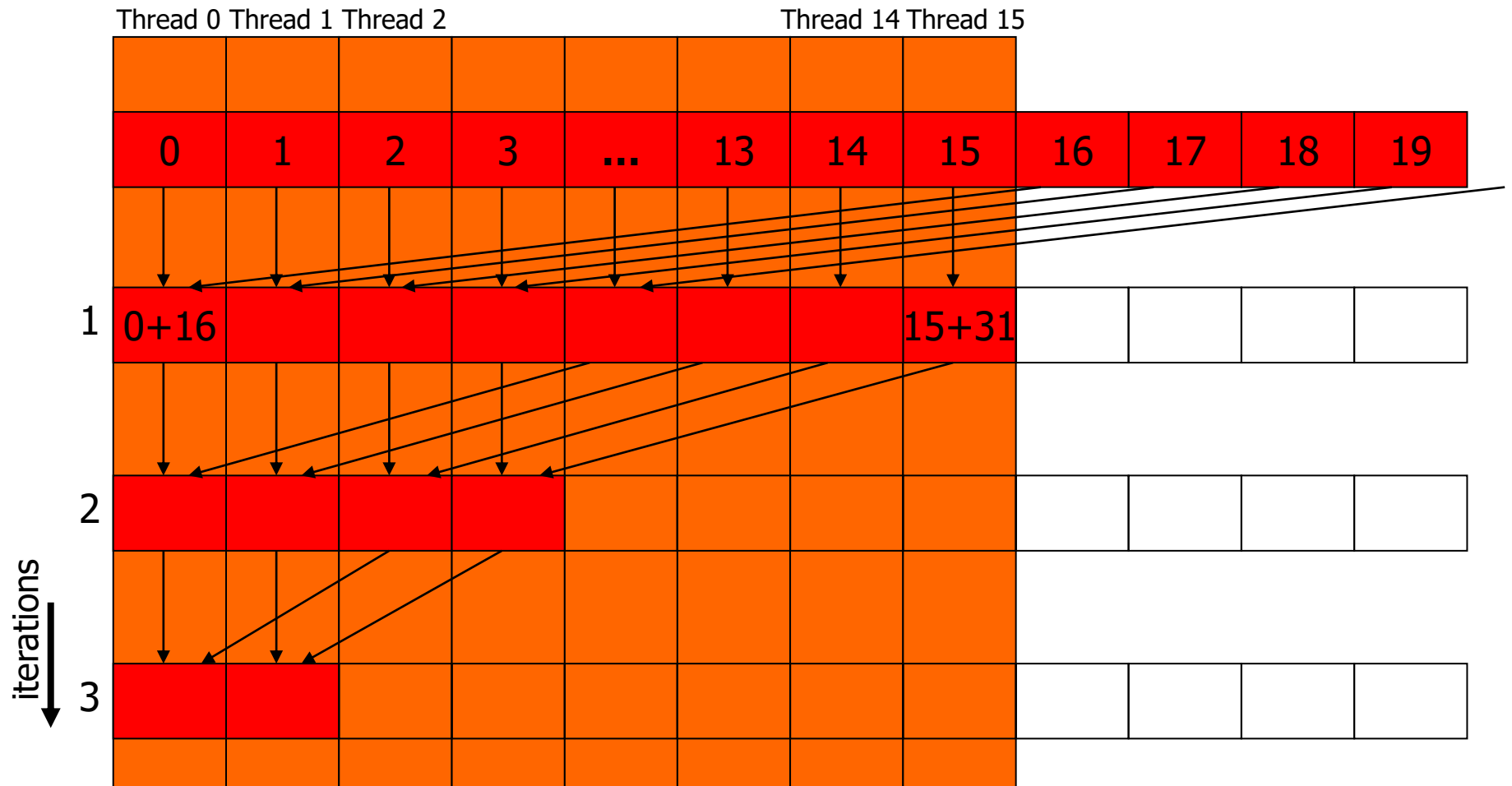
---

- Program with **low SIMD utilization**

```
__shared__ float partialSum[]  
  
unsigned int t = threadIdx.x;  
  
for (int stride = 1; stride < blockDim.x; stride *= 2) {  
  
    __syncthreads();  
  
    if (t % (2*stride) == 0)  
        partialSum[t] += partialSum[t + stride];  
  
}
```

# Divergence-Free Mapping (I)

- All active threads belong to the same warp



Slide credit: Hwu & Kirk

# Divergence-Free Mapping (II)

---

- Program with high SIMD utilization

```
__shared__ float partialSum[]  
  
unsigned int t = threadIdx.x;  
  
for (int stride = blockDim.x; stride > 1; stride >> 1){  
  
    __syncthreads();  
  
    if (t < stride)  
        partialSum[t] += partialSum[t + stride];  
  
}
```

# Atomic Operations

# Shared Memory Atomic Operations

---

- Atomic Operations are needed when threads might **update the same memory locations at the same time**
- **CUDA:** `int atomicAdd(int*, int);`
- **PTX:** `atom.shared.add.u32 %r25, [%rd14], %r24;`
- **SASS:**

## Tesla, Fermi, Kepler

```
/*00a0*/ LDSLK P0, R9, [R8];  
/*00a8*/ @P0 IADD R10, R9, R7;  
/*00b0*/ @P0 STSCUL P1, [R8], R10;  
/*00b8*/ @!P1 BRA 0xa0;
```

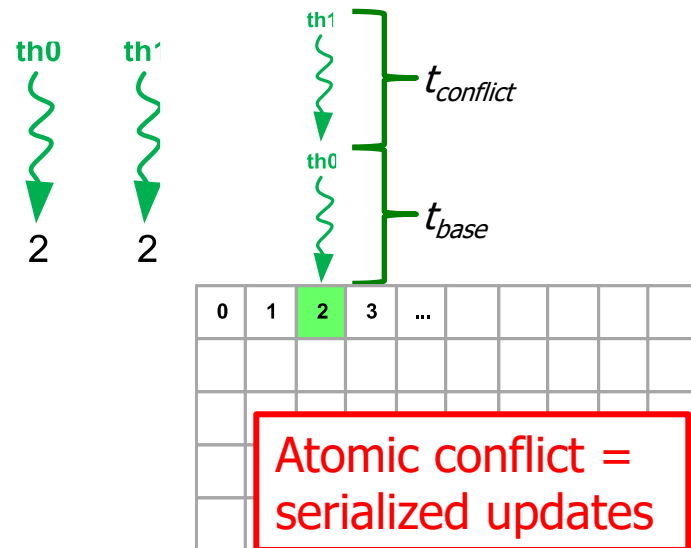
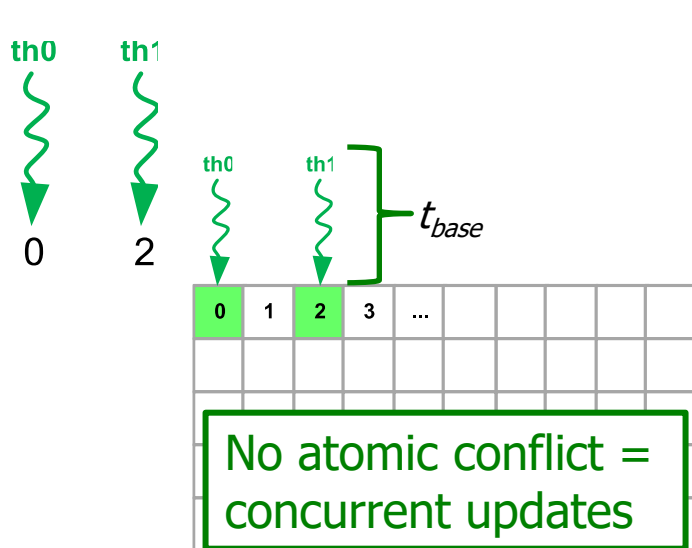
## Maxwell, Pascal, Volta

```
/*01f8*/ ATOMS.ADD RZ, [R7], R11;
```

Native atomic operations for  
32-bit integer, and 32-bit and  
64-bit atomicCAS

# Atomic Conflicts

- We define the intra-warp **conflict degree** as the number of threads in a warp that update the same memory position
- The conflict degree can be between 1 and 32

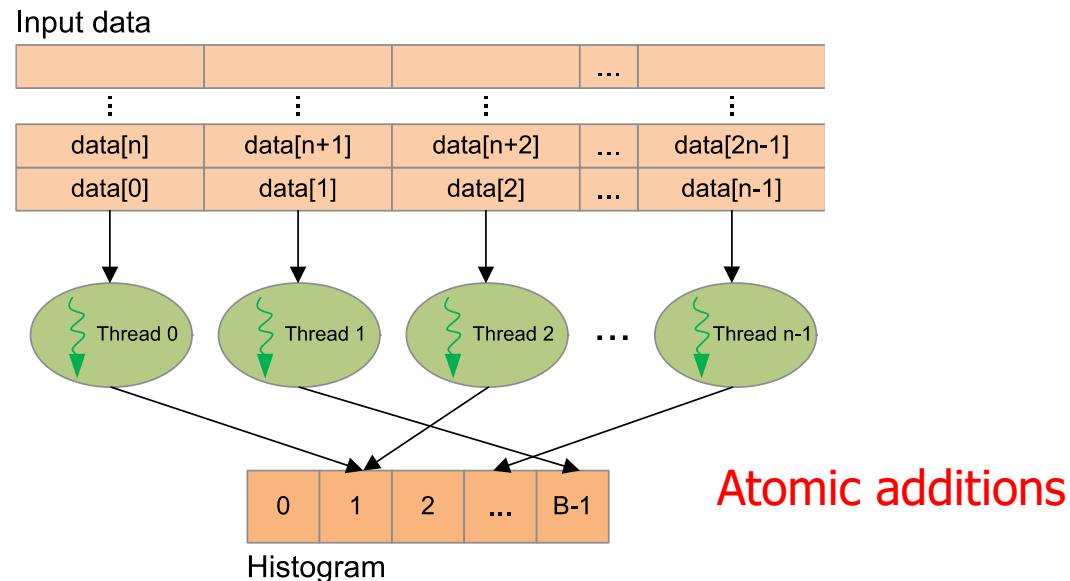




# Histogram Calculation

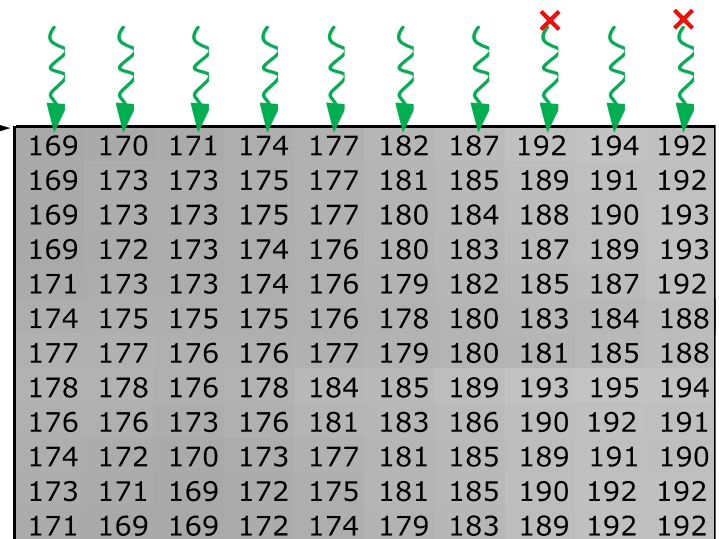
- Histograms count the number of data instances in disjoint categories (bins)

```
for (each pixel i in image I){  
    Pixel = I[i]                // Read pixel  
    Pixel' = Computation(Pixel) // Optional computation  
    Histogram[Pixel']++         // Vote in histogram bin  
}
```



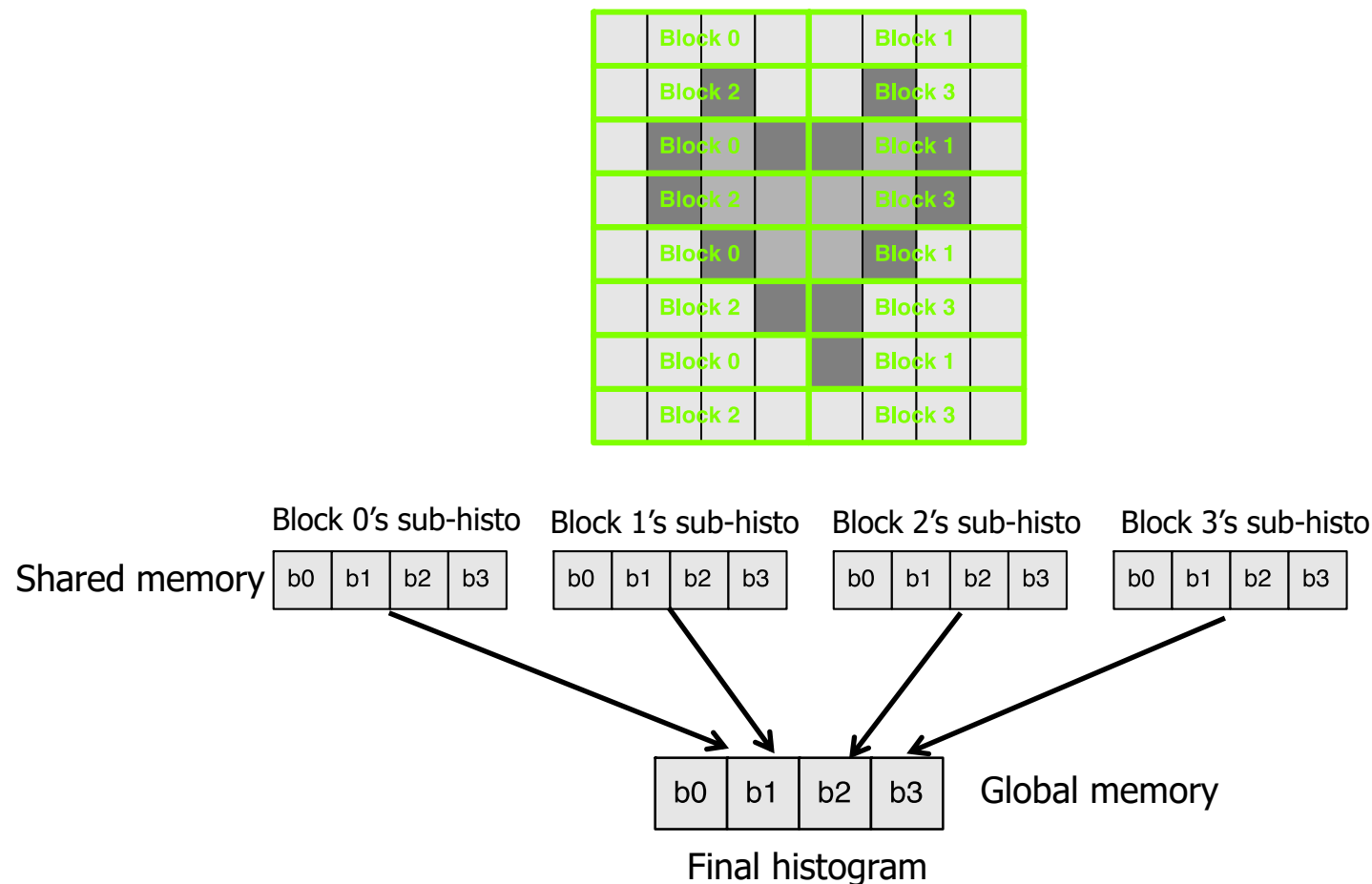
# Histogram Calculation of Natural Images

- Frequent conflicts in natural images



# Optimizing Histogram Calculation

- **Privatization:** Per-block sub-histograms in shared memory

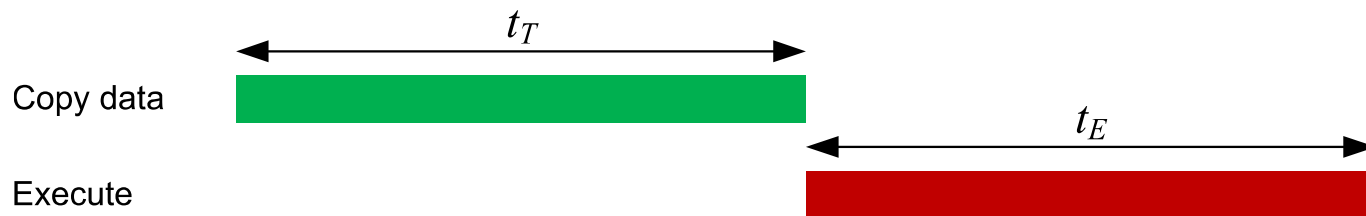


# Data Transfers between CPU and GPU

# Data Transfers

---

- Synchronous and asynchronous transfers
- Streams (Command queues)
  - Sequence of operations that are performed in order
    - CPU-GPU data transfer
    - Kernel execution
      - D input data instances, B blocks
    - GPU-CPU data transfer
  - Default stream



# Asynchronous Transfers

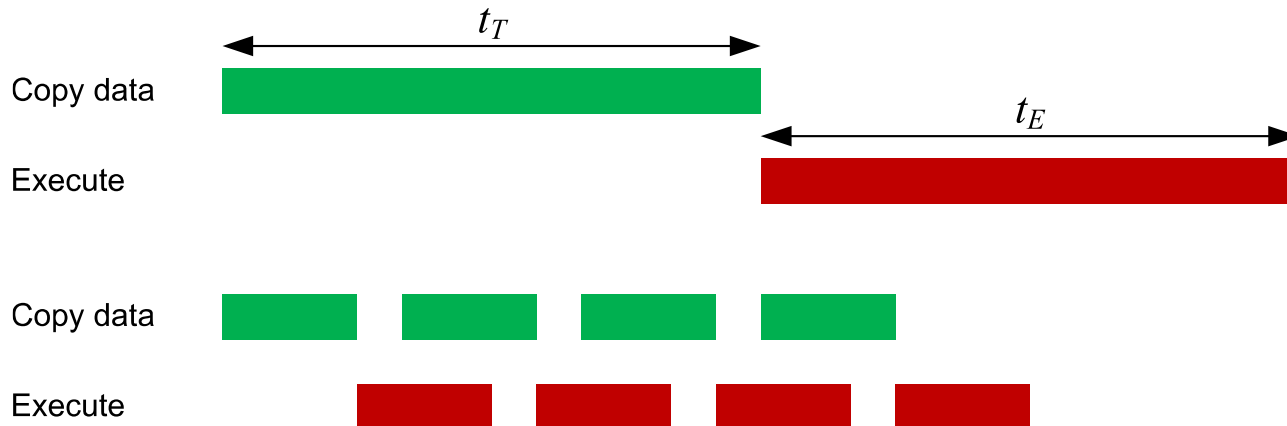
- Computation **divided into nStreams**

- D input data instances, B blocks

- nStreams

- D/nStreams data instances

- B/nStreams blocks



- Estimates

$$t_E + \frac{t_T}{nStreams}$$

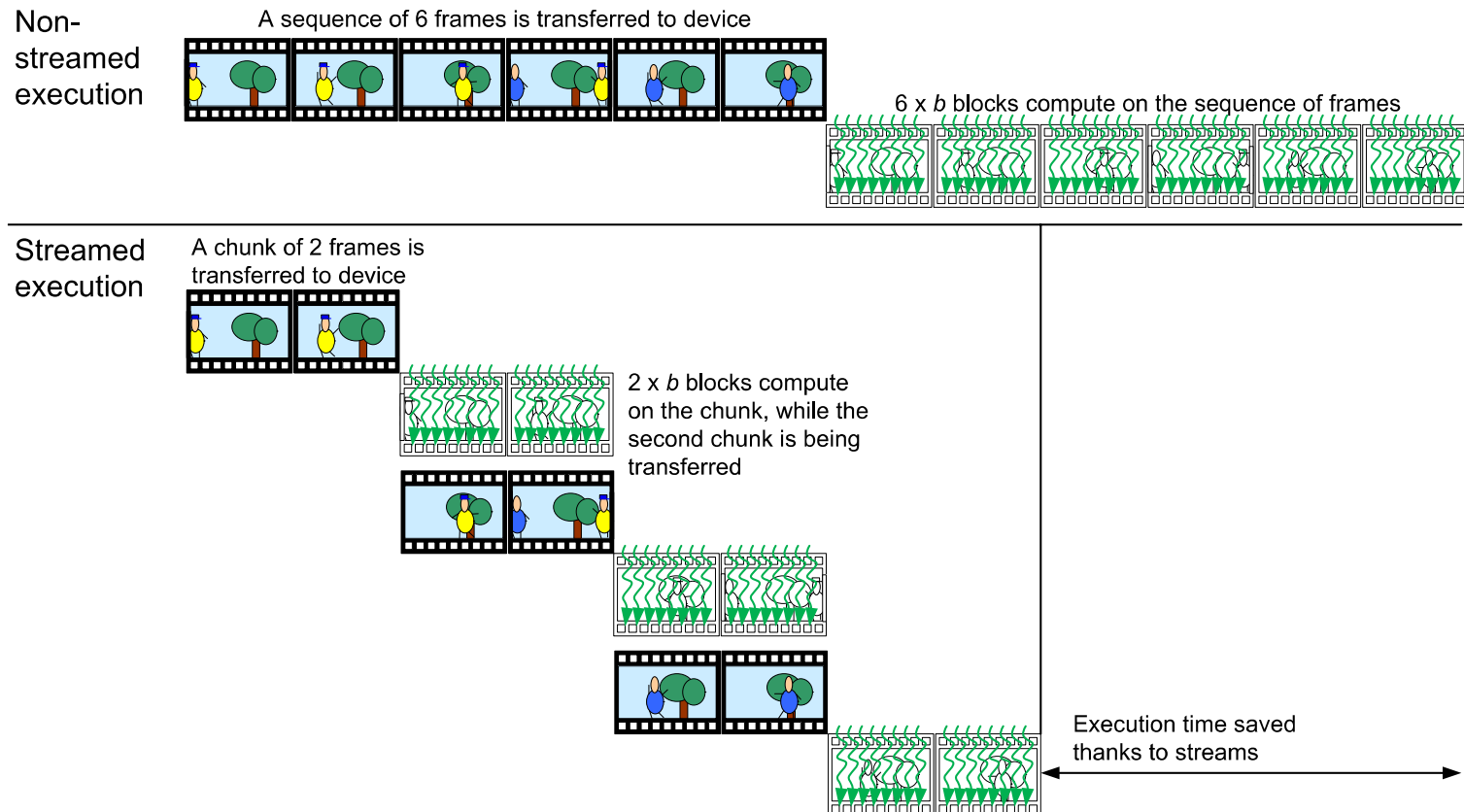
$t_E \geq t_T$  (dominant kernel)

$$t_T + \frac{t_E}{nStreams}$$

$t_T > t_E$  (dominant transfers)

# Overlap of Communication and Computation

- Applications with independent computation on different data instances can benefit from asynchronous transfers
- For instance, **video processing**



# Summary

---

- GPU as an accelerator
  - Program structure
    - Bulk synchronous programming model
  - Memory hierarchy and memory management
  - Performance considerations
    - Memory access
      - Latency hiding: occupancy (TLP)
      - Memory coalescing
      - Data reuse: shared memory
    - SIMD utilization
    - Atomic operations
    - Data transfers



# Collaborative Computing

# Review

---

- Device allocation, CPU-GPU transfer, and GPU-CPU transfer
  - ❑ `cudaMalloc()`;
  - ❑ `cudaMemcpy()`;

```
// Allocate input
malloc(input, ...);
cudaMalloc(d_input, ...);
cudaMemcpy(d_input, input, ..., HostToDevice); // Copy to device memory

// Allocate output
malloc(output, ...);
cudaMalloc(d_output, ...);

// Launch GPU kernel
gpu_kernel<<<blocks, threads>>> (d_output, d_input, ...);

// Synchronize
cudaDeviceSynchronize();

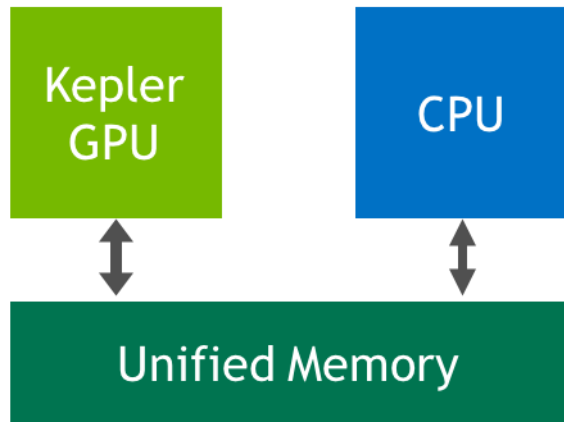
// Copy output to host memory
cudaMemcpy(output, d_output, ..., DeviceToHost);
```

# Unified Memory (I)

---

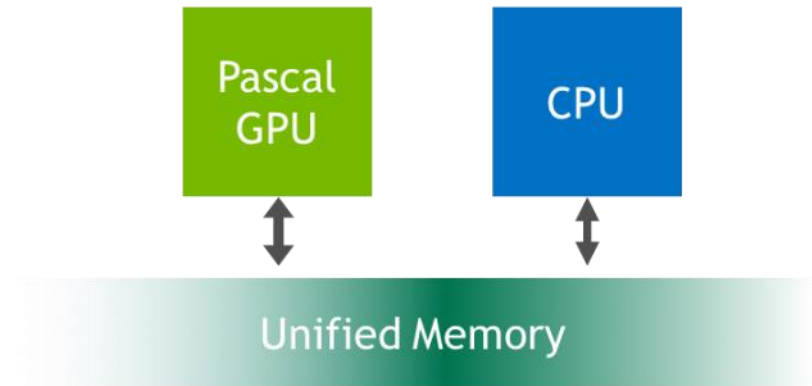
- Unified Virtual Address
- Since CUDA 6.0: **Unified memory**
- Since CUDA 8.0 + Pascal: **GPU page faults**

CUDA 6 Unified Memory



(Limited to GPU Memory Size)

Pascal Unified Memory



(Limited to System Memory Size)

# Unified Memory (II)

---

- Easier programming with **Unified Memory**
  - `cudaMallocManaged()`;

```
// Allocate input
malloc(input, ...);
cudaMallocManaged(d_input, ...);
memcpy(d_input, input, ...); // Copy to managed memory

// Allocate output
cudaMallocManaged(d_output, ...);

// Launch GPU kernel
gpu_kernel<<<blocks, threads>>> (d_output, d_input, ...);

// Synchronize
cudaDeviceSynchronize();
```

# Collaborative Computing Algorithms

---

- Case studies using CPU and GPU
- Kernel launches are asynchronous
  - CPU can work while waits for GPU to finish
  - Traditionally, this is the most efficient way to exploit heterogeneity

```
// Allocate input
malloc(input, ...);
cudaMalloc(d_input, ...);
cudaMemcpy(d_input, input, ..., HostToDevice); // Copy to device memory

// Allocate output
malloc(output, ...);
cudaMalloc(d_output, ...);

// Launch GPU kernel
gpu_kernel<<<blocks, threads>>> (d_output, d_input, ...);

// CPU can do things here

// Synchronize
cudaDeviceSynchronize();

// Copy output to host memory
cudaMemcpy(output, d_output, ..., DeviceToHost);
```

# Fine-Grained Heterogeneity

---

- Fine-grain heterogeneity becomes possible with Pascal/Volta architecture
- Pascal/Volta Unified Memory
  - CPU-GPU memory coherence
  - System-wide atomic operations

```
// Allocate input
cudaMallocManaged(input, ...);

// Allocate output
cudaMallocManaged(output, ...);

// Launch GPU kernel
gpu_kernel<<<blocks, threads>>> (output, input, ...);

// CPU can do things here
output[x] = input[y];

output[x+1].fetch_add(1);
```

# Since CUDA 8.0

---

- Unified memory

```
cudaMallocManaged(&h_in, in_size);
```

- System-wide atomics

```
old = atomicAdd_system(&h_out[x], inc);
```

# Since OpenCL 2.0

---

## ■ Shared virtual memory

```
XYZ * h_in = (XYZ *)clSVMAlloc(  
    ocl.clContext, CL_MEM_SVM_FINE_GRAIN_BUFFER, in_size, 0);
```

## ■ More flags:

```
CL_MEM_READ_WRITE  
CL_MEM_SVM_ATOMICS
```

## ■ C++11 atomic operations

(memory\_scope\_all\_svm\_devices)

```
old = atomic_fetch_add(&h_out[x], inc);
```



# C++AMP (HCC)

---

- Unified memory space (HSA)

```
XYZ *h_in = (XYZ *)malloc(in_size);
```

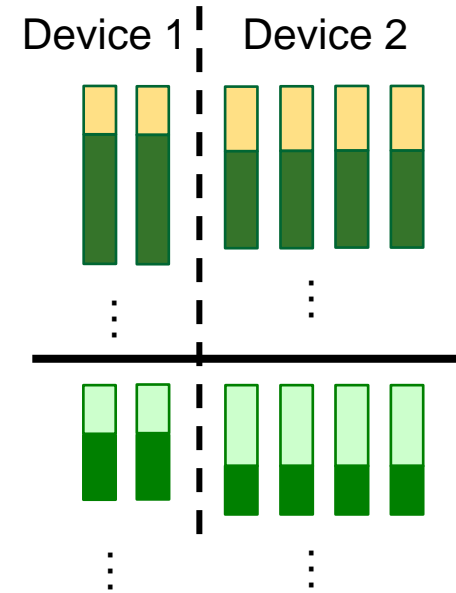
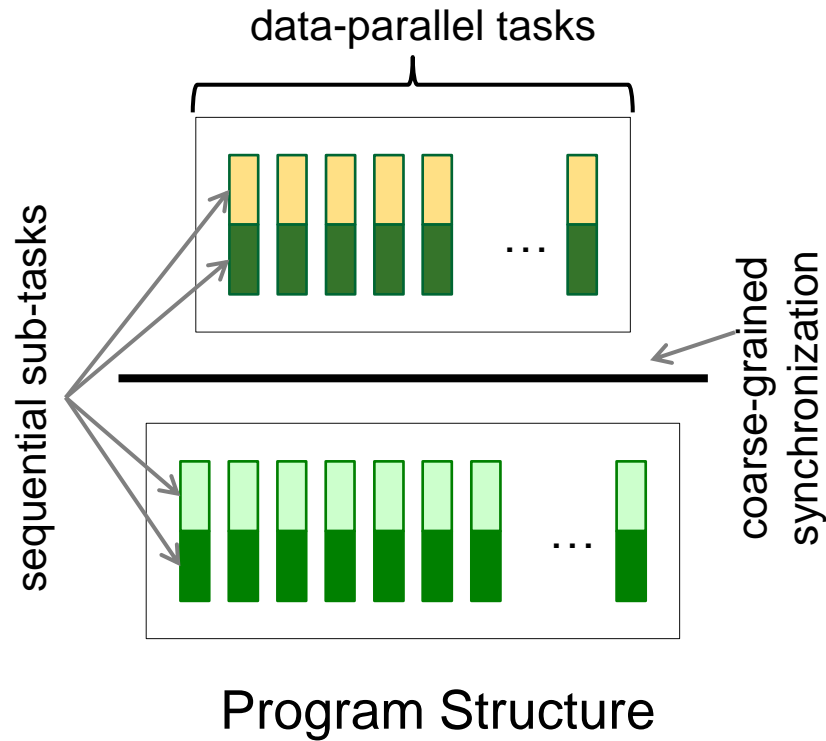
- C++11 atomic operations

```
(memory_scope_all_svm_devices)
```

- Platform atomics (HSA)

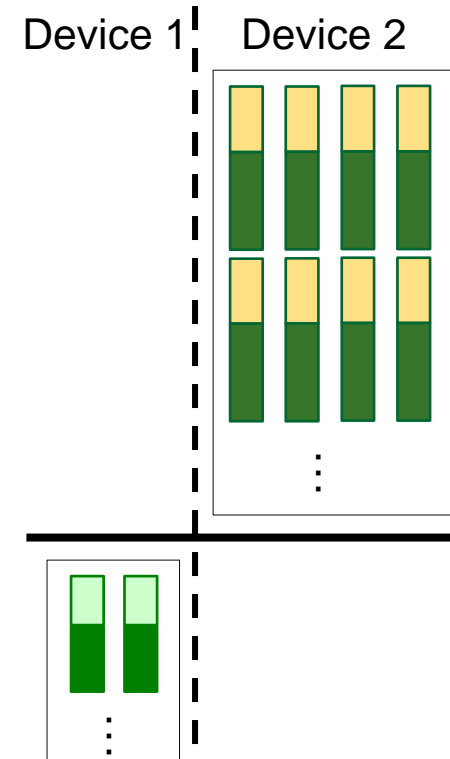
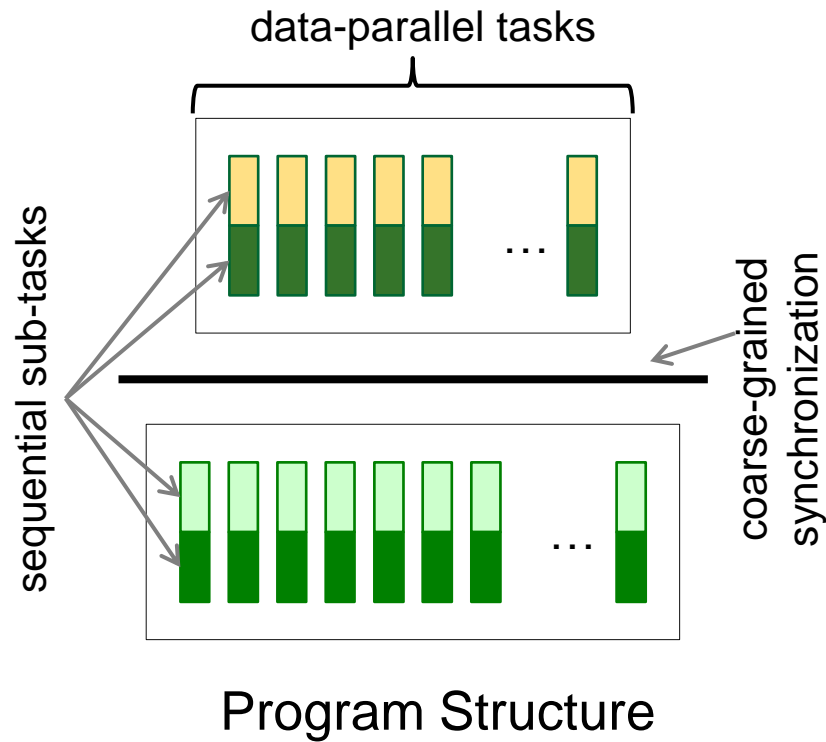
```
old = atomic_fetch_add(&h_out[x], inc);
```

# Collaborative Patterns (I)



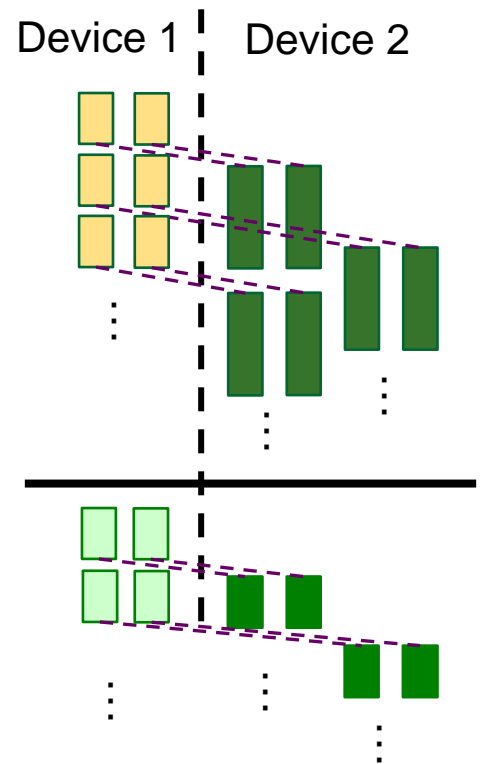
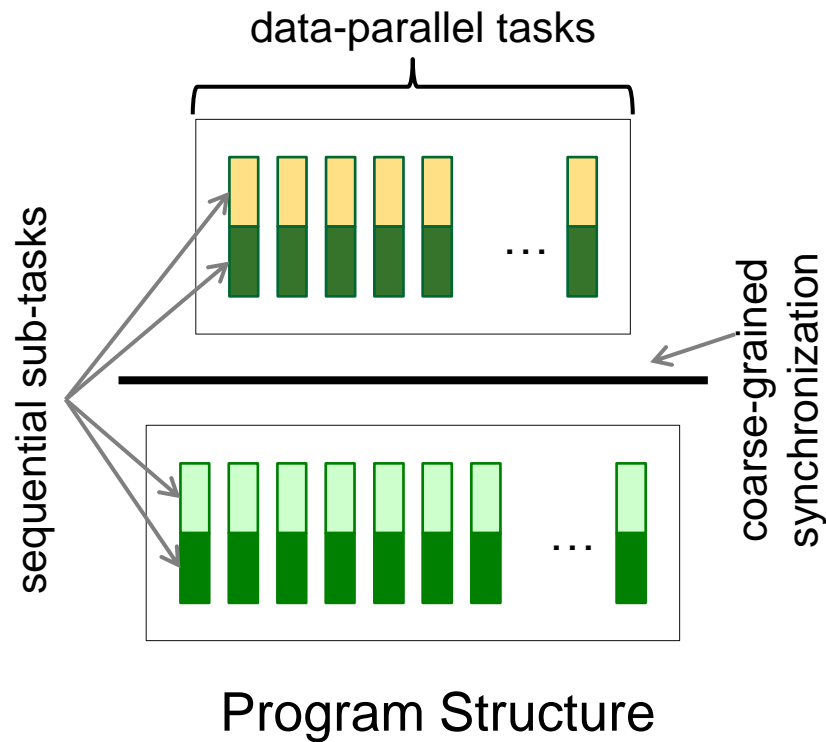
Data Partitioning

# Collaborative Patterns (II)



Coarse-grained Task Partitioning

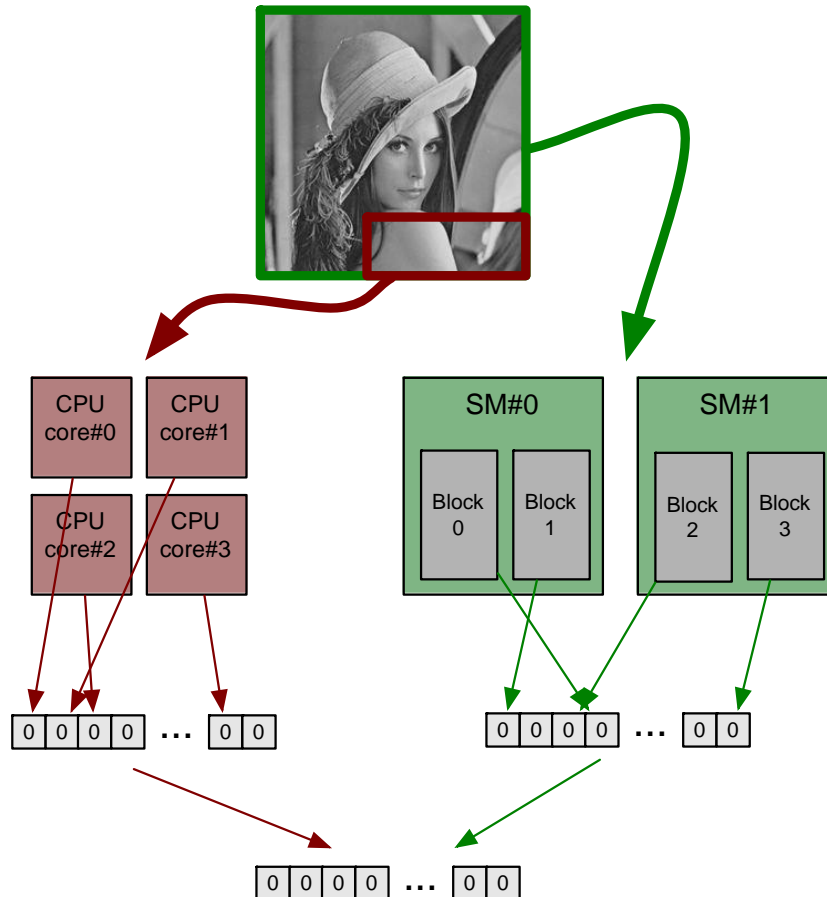
# Collaborative Patterns (III)



**Fine-grained Task Partitioning**

# Histogram (I)

- Previous generations: **separate CPU and GPU histograms** are merged at the end



```
malloc(CPU image);
cudaMalloc(GPU image);
cudaMemcpy(GPU image, CPU image, ...,
           Host to Device);
malloc(CPU histogram);
memset(CPU histogram, 0);
cudaMalloc(GPU histogram);
cudaMemset(GPU histogram, 0);

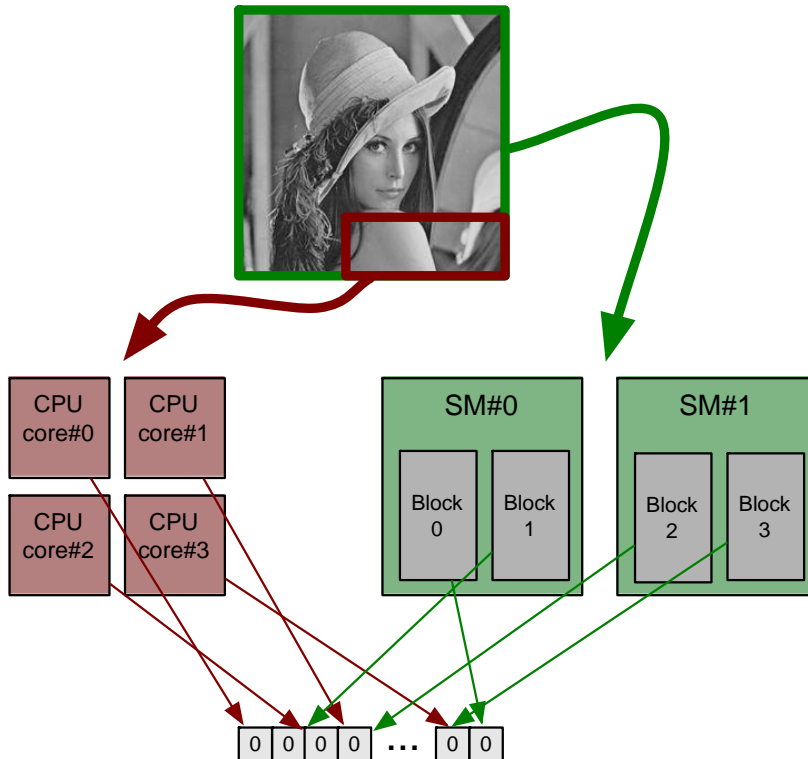
// Launch CPU threads
// Launch GPU kernel

cudaMemcpy(GPU histogram, DeviceToHost);

// Launch CPU threads for merging
```

# Histogram (II)

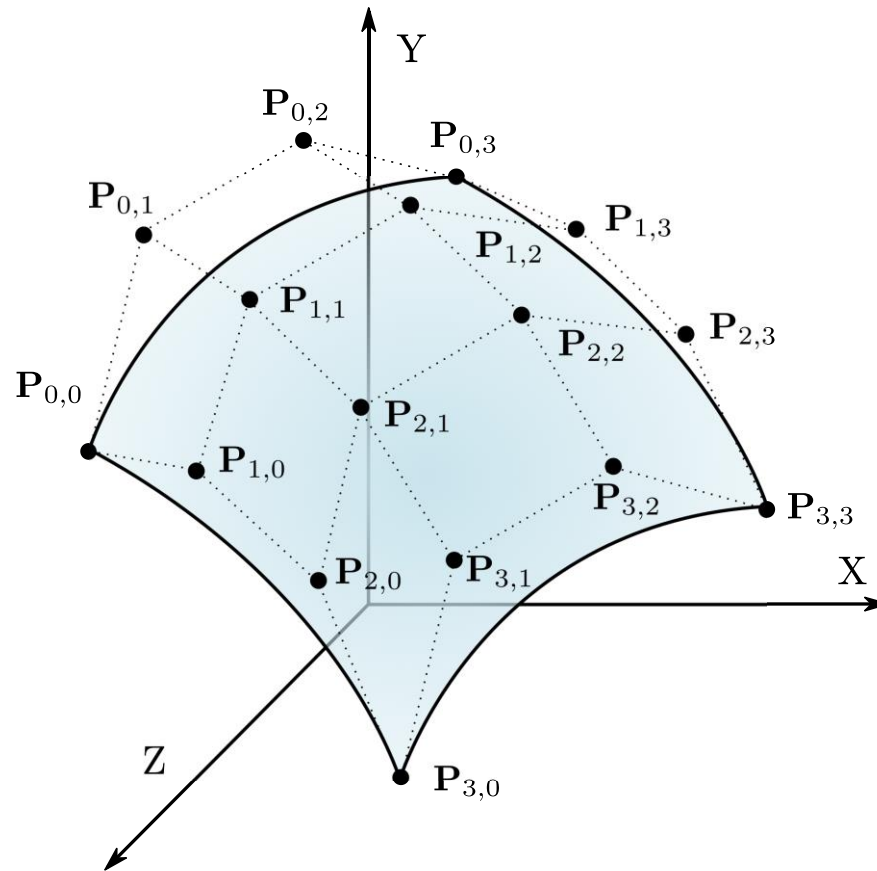
- System-wide atomic operations: **one single histogram**



```
cudaMallocManaged(Histogram);  
cudaMemset(Histogram, 0);  
  
// Launch CPU threads  
// Launch GPU kernel (atomicAdd_system)
```

# Bézier Surfaces (I)

- Bézier surface: 4x4 net of control points



# Bézier Surfaces (II)

---

- Parametric non-rational formulation
  - Bernstein polynomials
  - Bi-cubic surface  $m = n = 3$

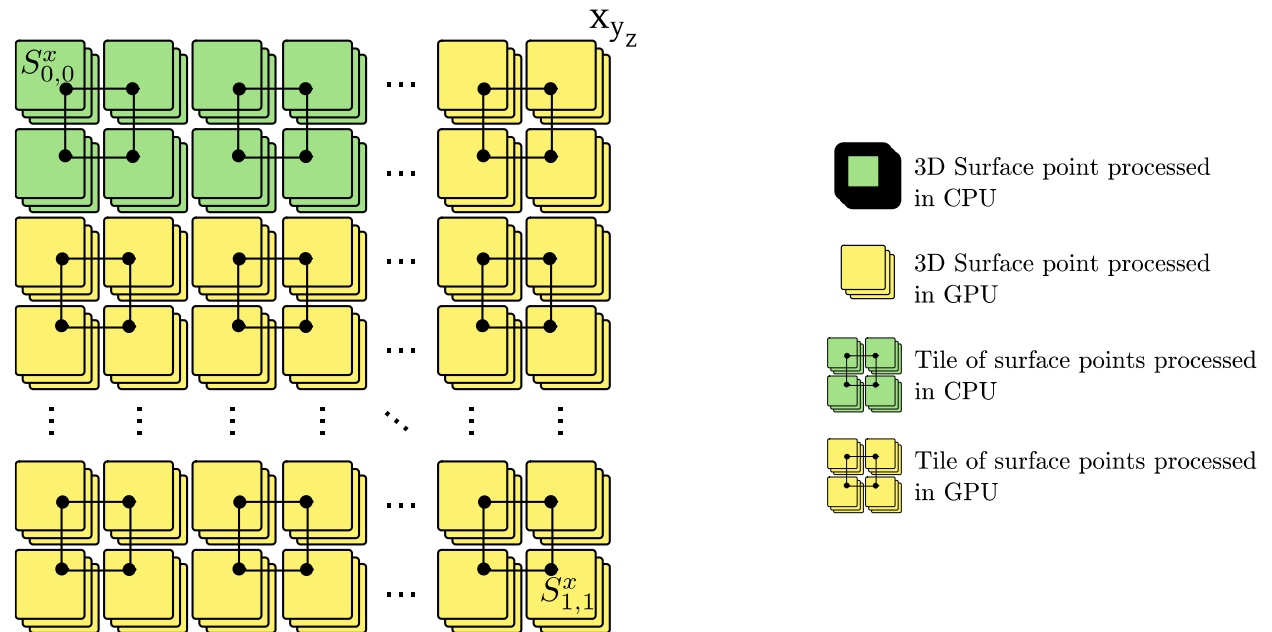
$$\mathbf{S}(u, v) = \sum_{i=0}^m \sum_{j=0}^n \mathbf{P}_{i,j} B_{i,m}(u) B_{j,n}(v), \quad (1)$$

$$B_{i,m}(u) = \binom{m}{i} (1-u)^{(m-i)} u^i, \quad (2)$$



# Bézier Surfaces (III)

- Collaborative implementation
  - Tiles calculated by GPU blocks or CPU threads
  - **Static distribution**



# Bézier Surfaces (IV)

---

## ■ Without Unified Memory

```
// Allocate control points
malloc(control_points, ...);
generate_cp(control_points);
cudaMalloc(d_control_points, ...);
cudaMemcpy(d_control_points, control_points, ..., HostToDevice); // Copy to device memory

// Allocate surface
malloc(surface, ...);
cudaMalloc(d_surface, ...);

// Launch CPU threads
std::thread main_thread (run_cpu_threads, control_points, surface, ...);

// Launch GPU kernel
gpu_kernel<<<blocks, threads>>> (d_surface, d_control_points, ...);

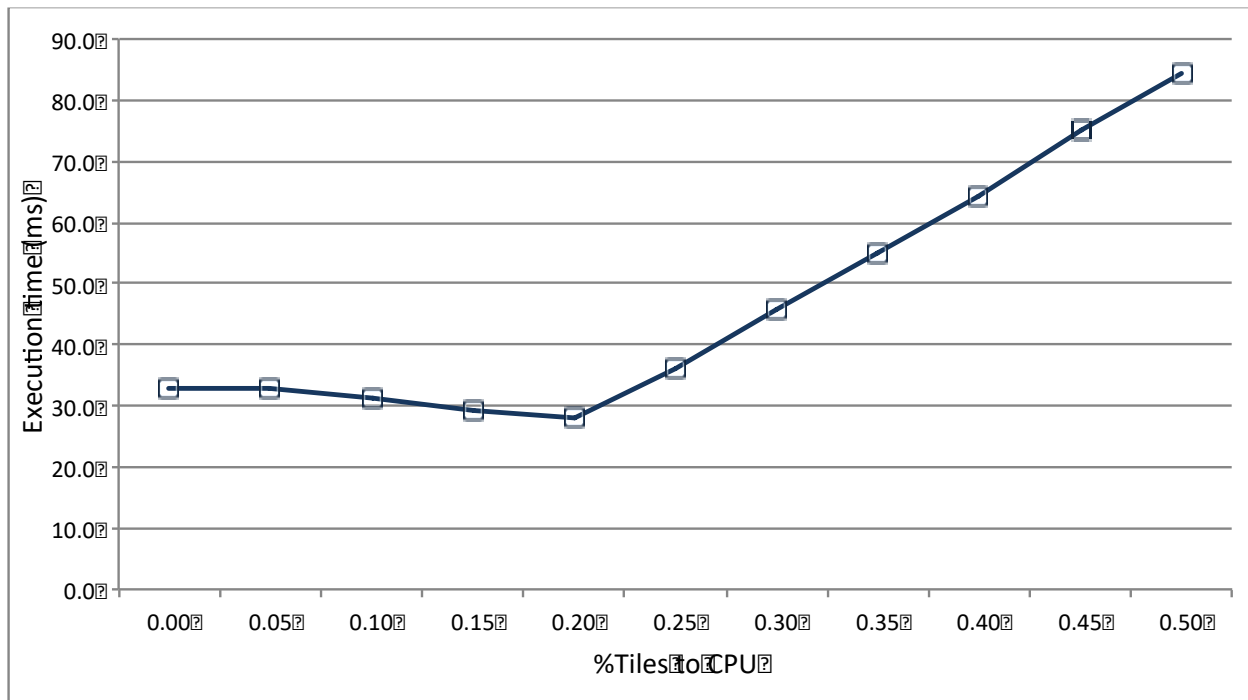
// Synchronize
main_thread.join();
cudaDeviceSynchronize();

// Copy gpu part of surface to host memory
cudaMemcpy(&surface[end_of_cpu_part], d_surface, ..., DeviceToHost);
```

# Bézier Surfaces (V)

## ■ Execution results

- Bezier surface: 300x300, 4x4 control points
- %Tiles to CPU
- NVIDIA Jetson TX1 (4 ARMv8 + 2 SMX): 17% speedup wrt GPU only



# Bézier Surfaces (VI)

---

## ■ With Unified Memory (Pascal/Volta)

```
// Allocate control points
malloc(control_points, ...);
generate_cp(control_points);
cudaMalloc(d_control_points, ...);
cudaMemcpy(d_control_points, control_points, ..., HostToDevice); // Copy to device memory

// Allocate surface
cudaMallocManaged(surface, ...);

// Launch CPU threads
std::thread main_thread (run_cpu_threads, control_points, surface, ...);

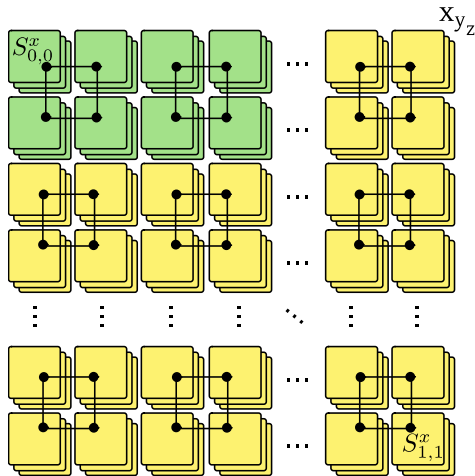
// Launch GPU kernel
gpu_kernel<<<blocks, threads>>> (surface, d_control_points, ...);

// Synchronize
main_thread.join();
cudaDeviceSynchronize();
```

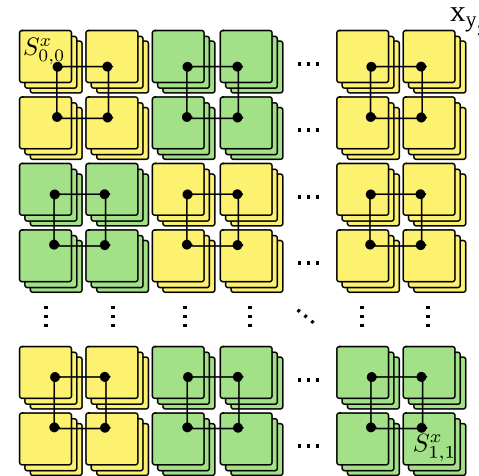
# Bézier Surfaces (VII)





## ■ Static vs. dynamic implementation

(a) Static Distribution



(b) Dynamic Distribution



-  3D Surface point processed in CPU
-  3D Surface point processed in GPU
-  Tile of surface points processed in CPU
-  Tile of surface points processed in GPU

## □ Pascal/Volta Unified Memory: **system-wide atomic operations**

```
while(true){
    if(threadIdx.x == 0)
        my_tile = atomicAdd_system(tile_num, 1); // my_tile in shared memory; tile_num in UM

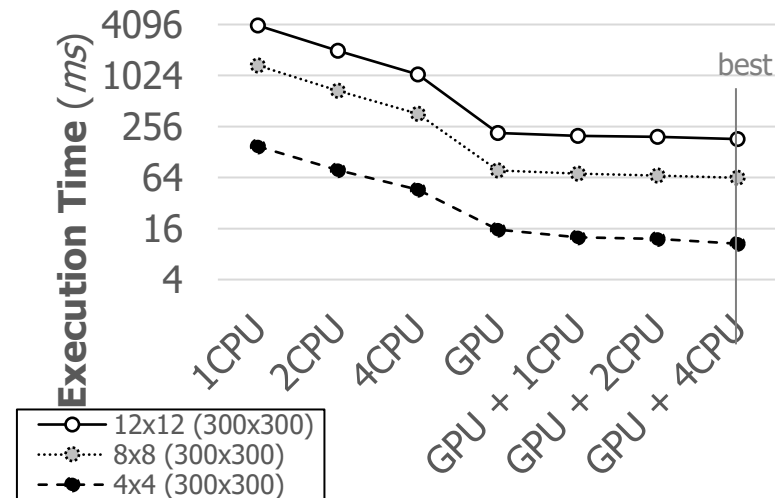
    __syncthreads(); // Synchronization

    if(my_tile >= number_of_tiles) break; // Break when all tiles processed

    ...
}
```

# Benefits of Collaboration

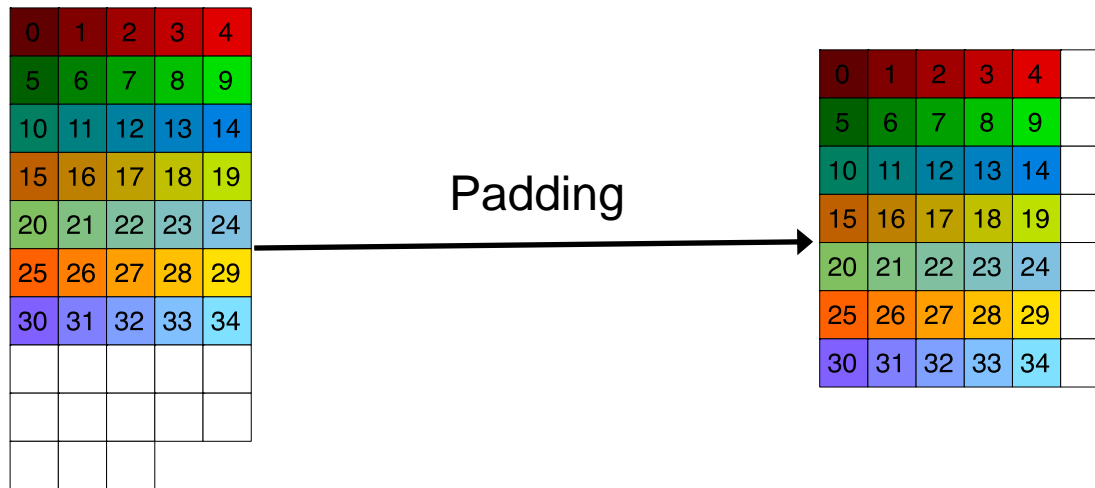
- Data partitioning improves performance
  - AMD Kaveri (4 CPU cores + 8 GPU CUs)



**Bézier Surfaces**  
(up to 47% improvement over GPU only)

# Padding (I)

- Matrix padding
  - Memory alignment
  - Transposition of near-square matrices

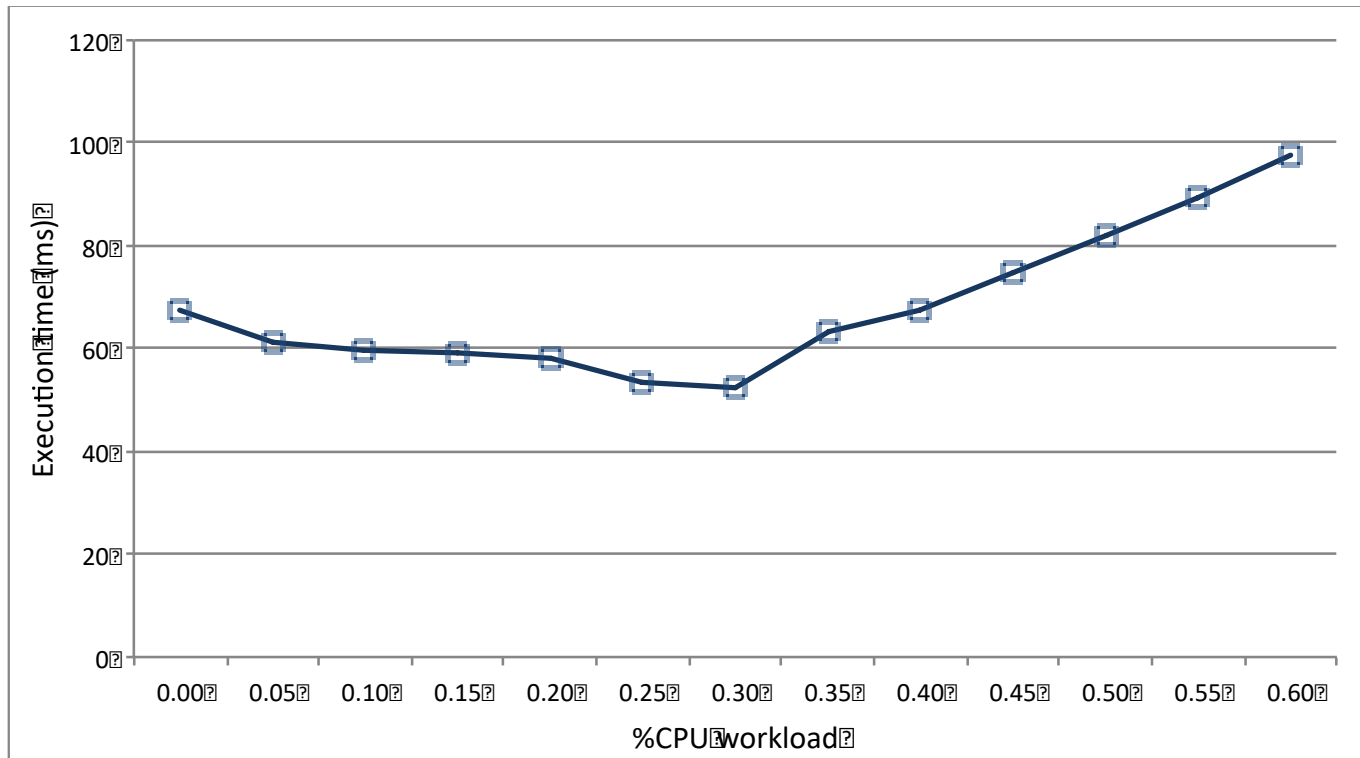


- Traditionally, it can only be performed out-of-place

# Padding (II)

## ■ Execution results

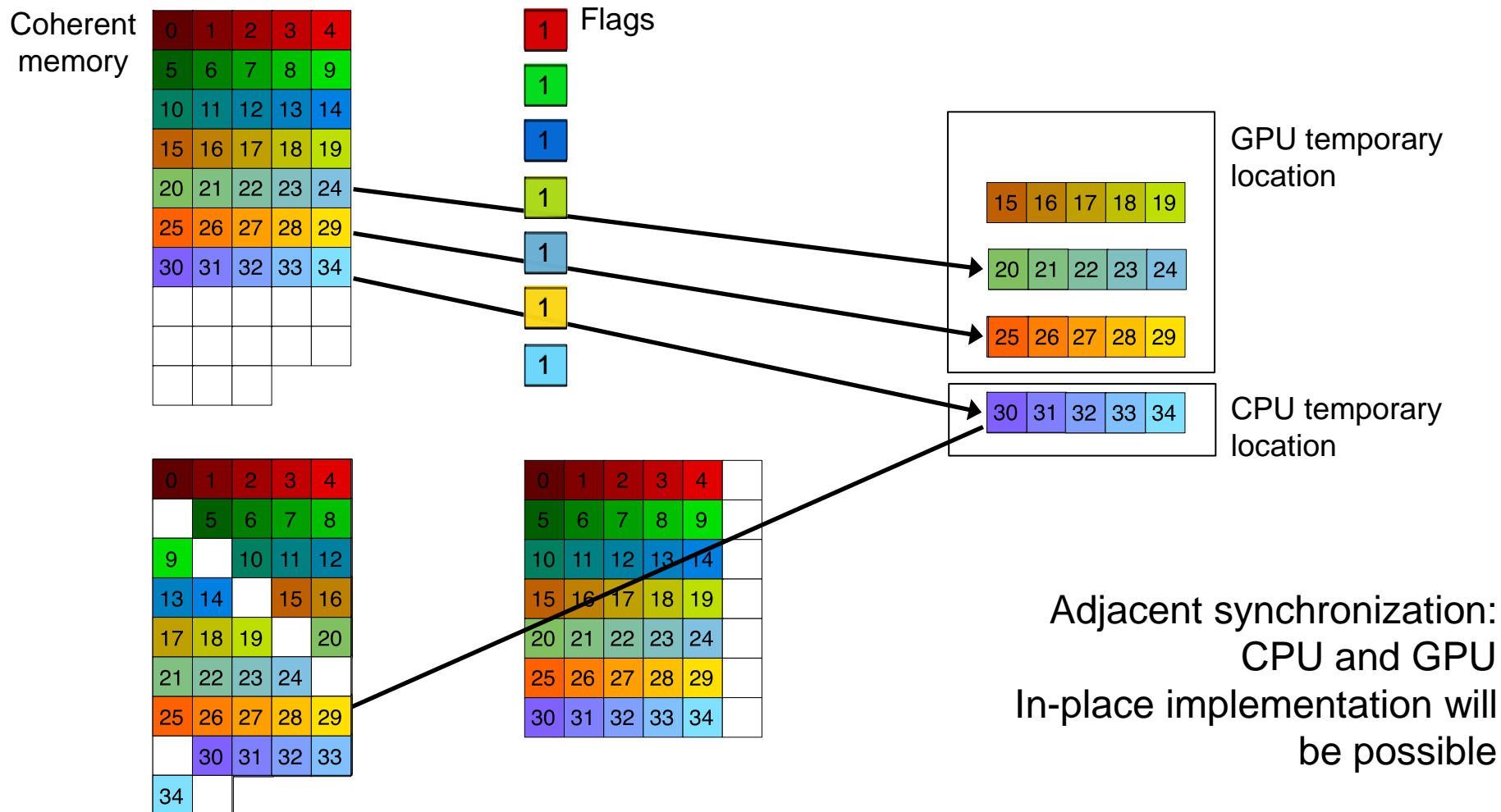
- ❑ Matrix size: 4000x4000, padding = 1
- ❑ NVIDIA Jetson TX1 (4 ARMv8 + 2 SMX): 29% speedup wrt GPU only





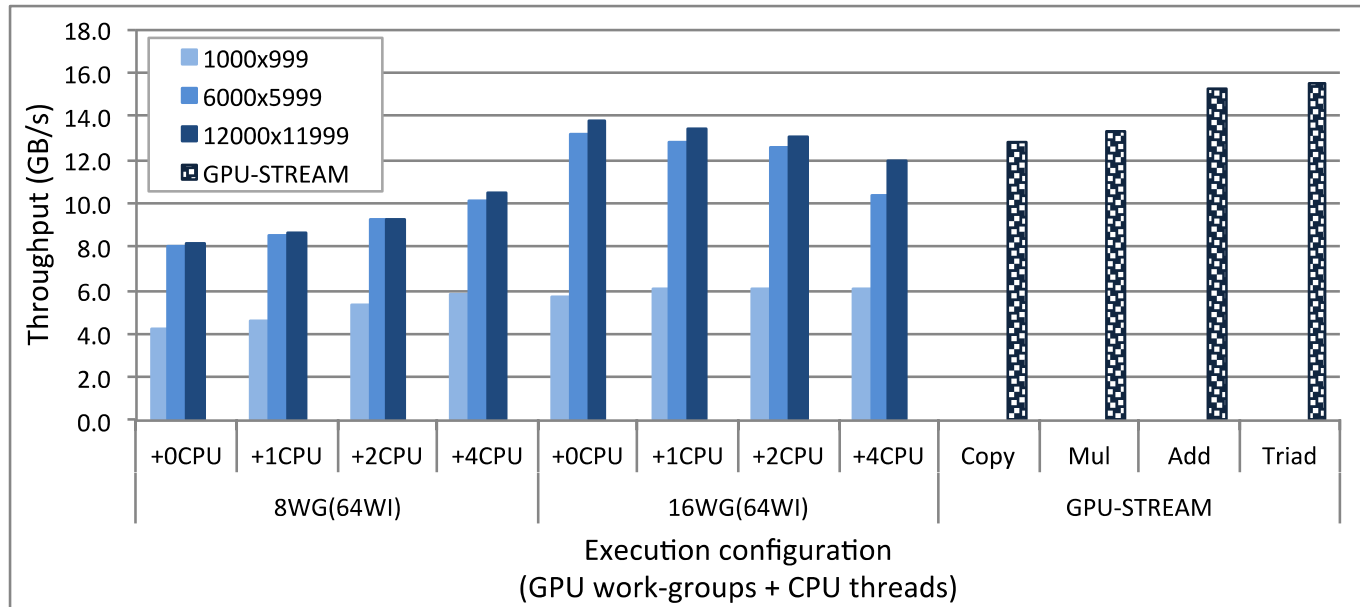
# In-Place Padding

## ■ Pascal/Volta Unified Memory



# Benefits of Collaboration

- Optimal number of devices is not always max
  - AMD Kaveri (4 CPU cores + 8 GPU CUs)

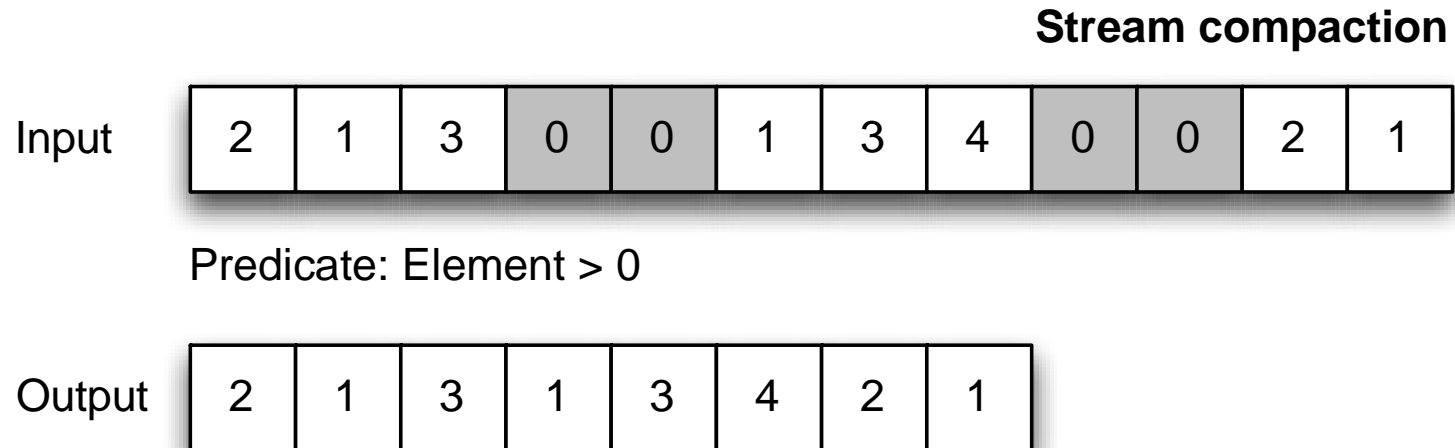


# Stream Compaction (I)

---

## ■ Stream compaction

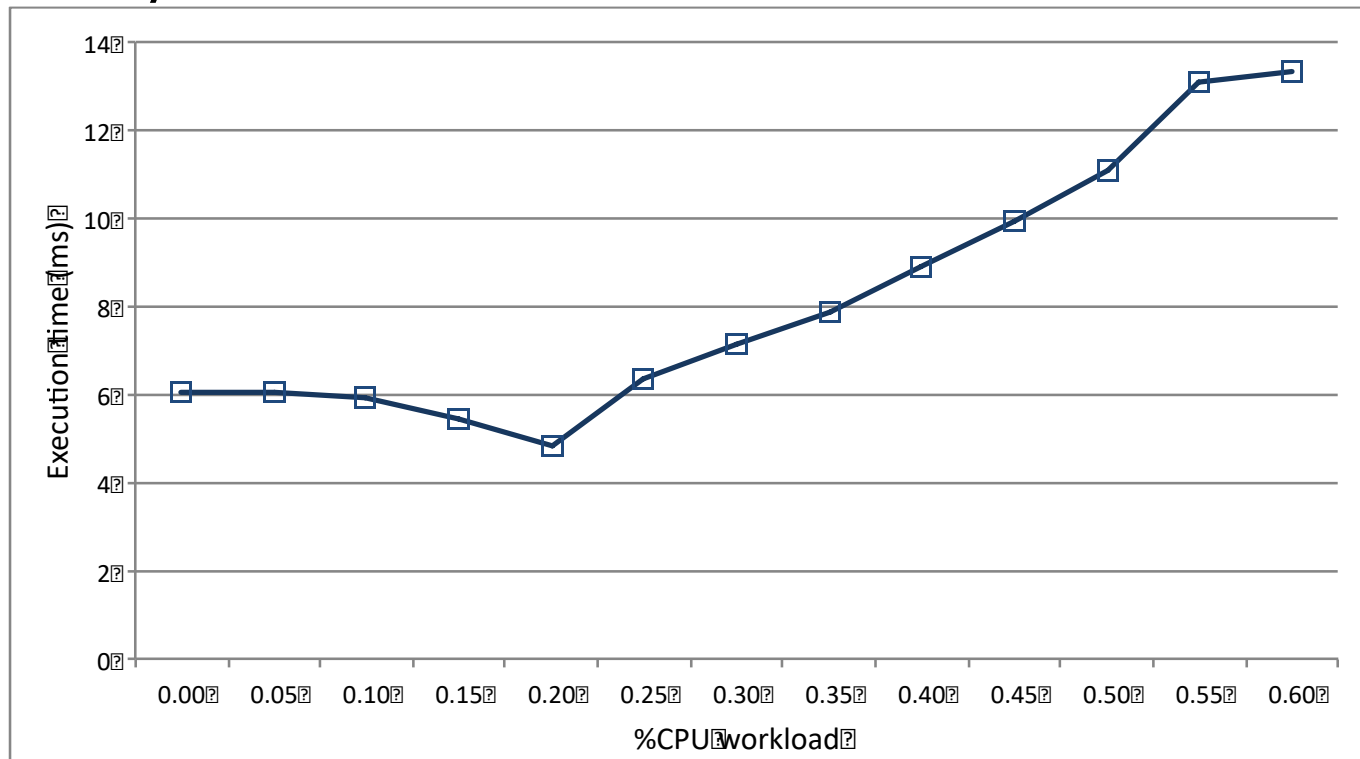
- Saving memory storage in sparse data
- Similar to padding, but local reduction result (non-zero element count) is propagated



# Stream Compaction (II)

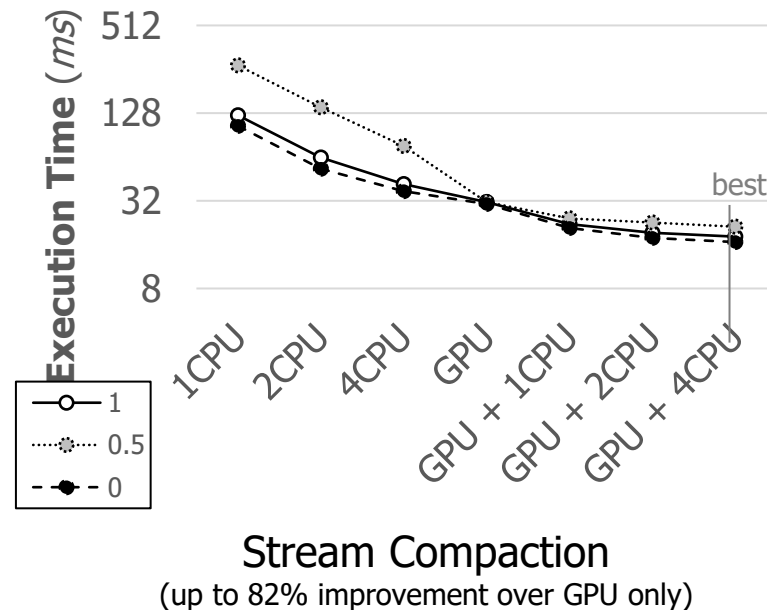
## ■ Execution results

- Array size: 2 MB, Filtered items = 50%
- NVIDIA Jetson TX1 (4 ARMv8 + 2 SMX): 25% speedup wrt GPU only



# Benefits of Collaboration

- Data partitioning improves performance
  - AMD Kaveri (4 CPU cores + 8 GPU CUs)



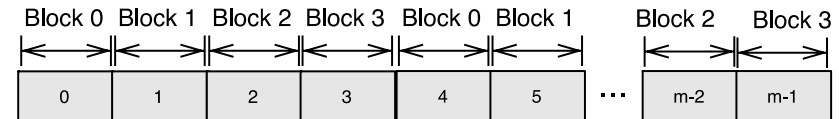
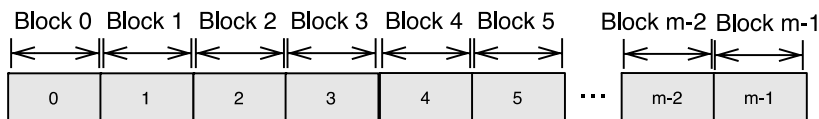
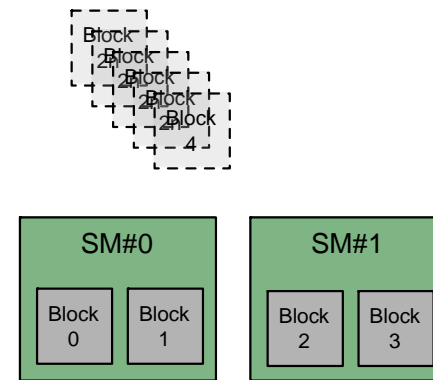
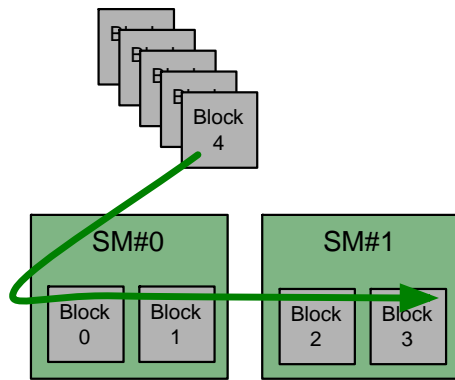
# Breadth-First Search

---

- Small-sized and big-sized frontiers
  - Top-down approach
  - Kernel 1 and Kernel 2
- Atomic-based block synchronization
  - Avoids kernel re-launch
- Very small frontiers
  - Underutilize GPU resources
- Collaborative implementation

# Atomic-Based Block Synchronization (I)

- Combine Kernel 1 and Kernel 2
- We can **avoid kernel re-launch**
- We need to use **persistent thread blocks**
  - Kernel 2 launches ( $\text{frontier\_size} / \text{block\_size}$ ) blocks
  - Persistent blocks: up to ( $\text{number\_SMs} \times \text{max\_blocks\_SM}$ )



# Atomic-Based Block Synchronization (II)

---

## ■ Code (simplified)

```
// GPU kernel
const int gtid = blockIdx.x * blockDim.x + threadIdx.x;

while(frontier_size != 0){

    for(node = gtid; node < frontier_size; node += blockDim.x*gridDim.x){

        // Visit neighbors
        // Enqueue in output queue if needed (global or local queue)

    }

    // Update frontier_size

    // Global synchronization
}
```



# Atomic-Based Block Synchronization (III)

- Global synchronization (simplified)
  - At the end of each iteration

```
const int tid = threadIdx.x;
const int gtid = blockIdx.x * blockDim.x + threadIdx.x;
atomicExch(ptr_threads_run, 0);
atomicExch(ptr_threads_end, 0);
int frontier = 0;
...

frontier++;

if(tid == 0){
    atomicAdd(ptr_threads_end, 1); // Thread block finishes iteration
}

if(gtid == 0){
    while(atomicAdd(ptr_threads_end, 0) != gridDim.x){;} // Wait until all blocks finish

    atomicExch(ptr_threads_end, 0); // Reset
    atomicAdd(ptr_threads_run, 1); // Count iteration
}

if(tid == 0 && gtid != 0){
    while(atomicAdd(ptr_threads_run, 0) < frontier){;} // Wait until ptr_threads_run is updated
}

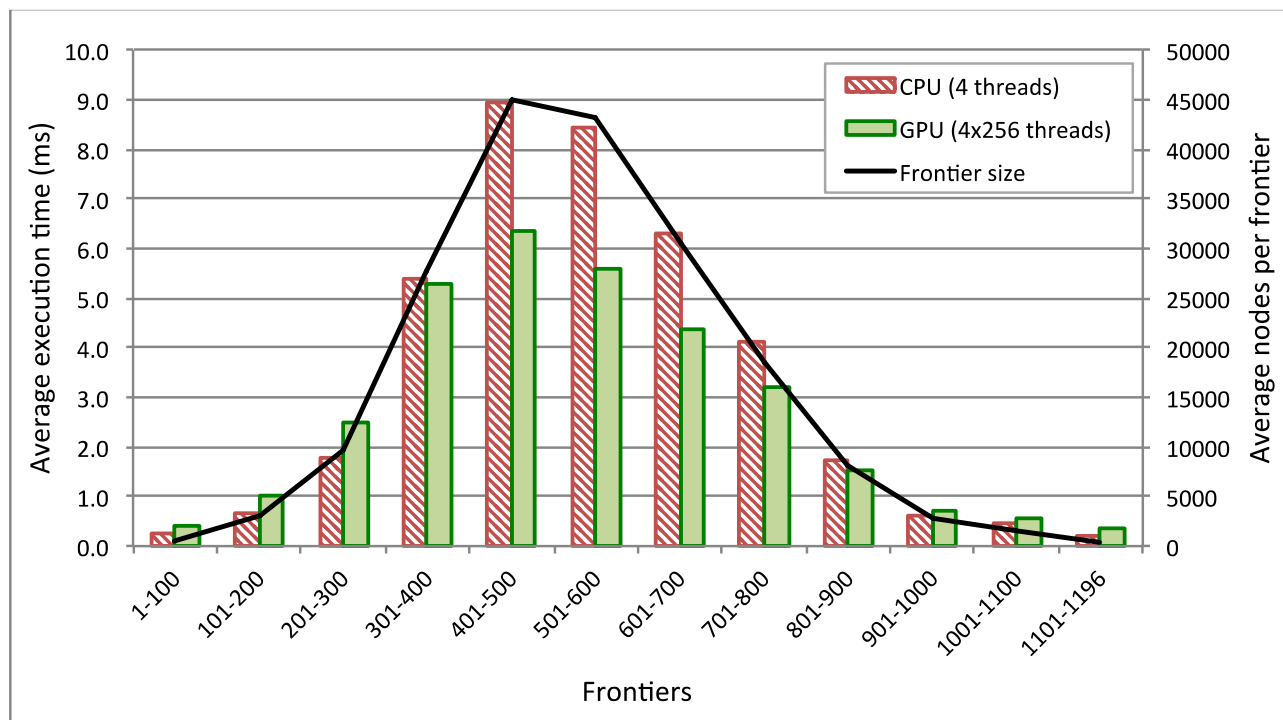
__syncthreads(); // Rest of threads wait here

...
```

# Collaborative Implementation (I)

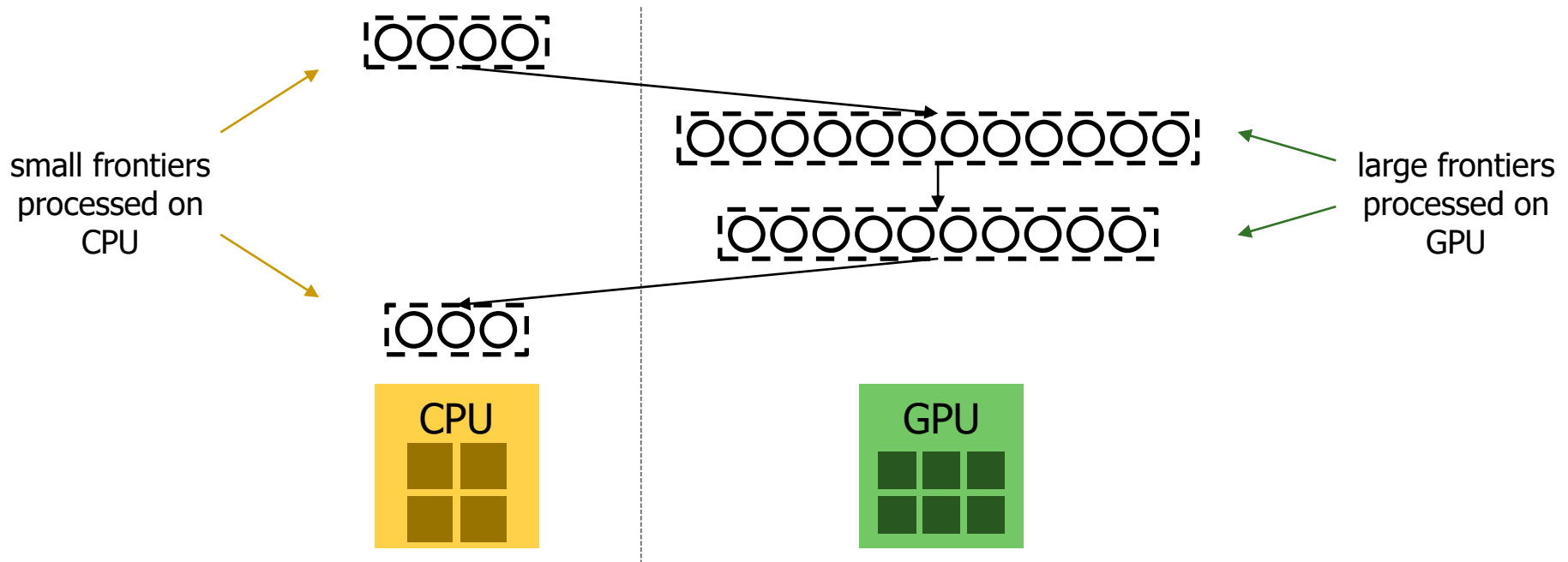
## ■ Motivation

- Small-sized frontiers underutilize GPU resources
  - NVIDIA Jetson TX1 (4 ARMv8 CPUs + 2 SMXs)
  - New York City roads



# Collaborative Implementation (II)

- Choose the most appropriate device



# Collaborative Implementation (III)

---

- Choose CPU or GPU depending on frontier size

```
// Host code
while(frontier_size != 0){

    if(frontier_size < LIMIT){

        // Launch CPU threads

    }
    else{

        // Launch GPU kernel

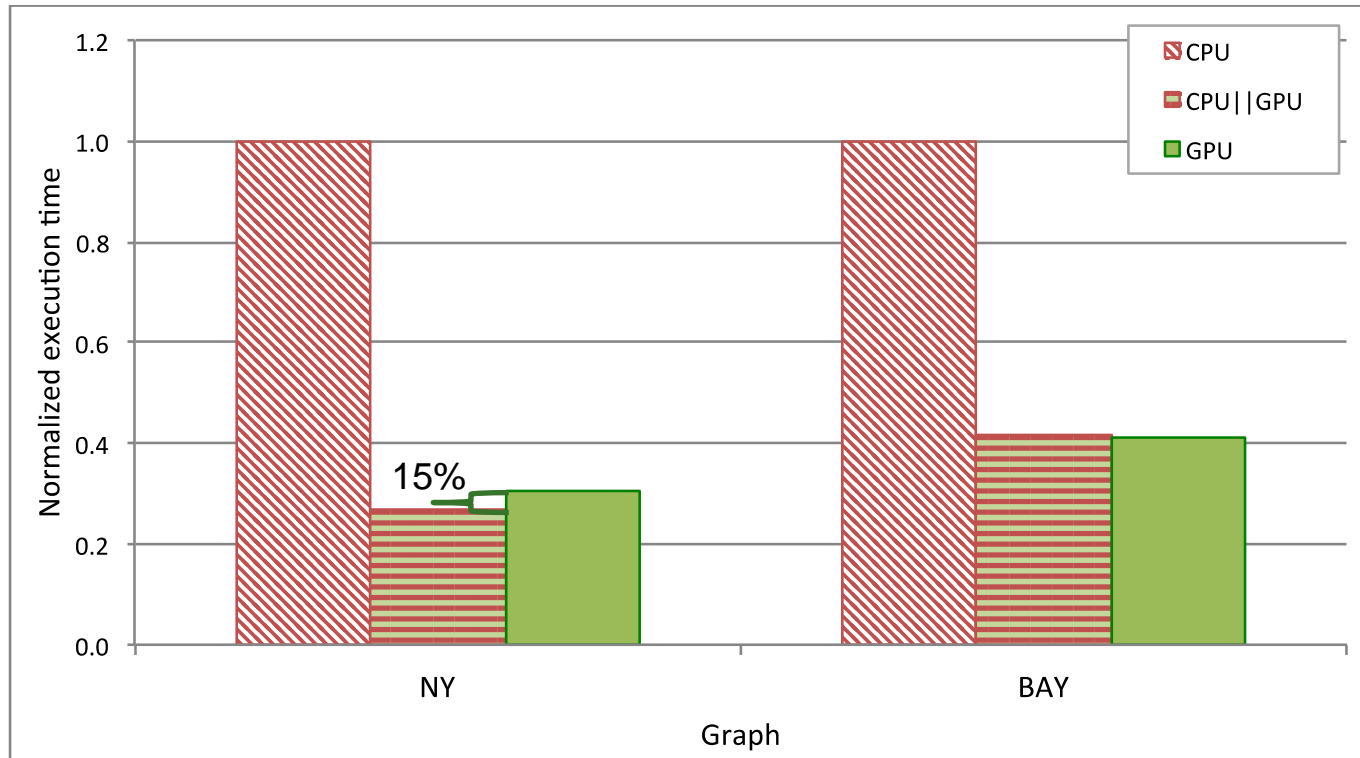
    }

}
```

- CPU threads or GPU kernel keep running while the condition is satisfied

# Collaborative Implementation (IV)

## ■ Execution results



# Collaborative Implementation (V)

---

- **Without** Unified Memory
  - Explicit memory copies

```
// Host code
while(frontier_size != 0){

    if(frontier_size < LIMIT){

        // Launch CPU threads

    }
    else{

        // Copy from host to device (queues and synchronization variables)

        // Launch GPU kernel

        // Copy from device to host (queues and synchronization variables)

    }

}
```

# Collaborative Implementation (VI)

---

## ■ Unified Memory

- ❑ `cudaMallocManaged()` ;
- ❑ Easier programming
- ❑ No explicit memory copies

```
// Host code
while(frontier_size != 0){

    if(frontier_size < LIMIT){

        // Launch CPU threads

    }
    else{

        // Launch GPU kernel

        cudaDeviceSynchronize() ;

    }

}
```

# Collaborative Implementation (VII)

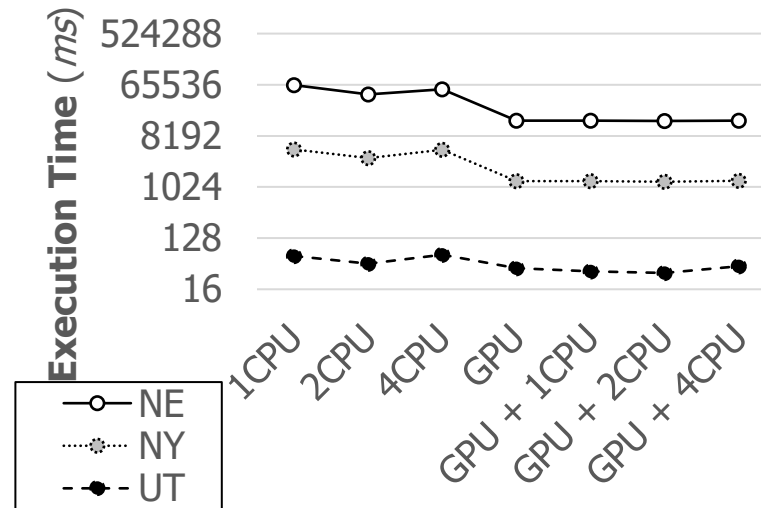
---

- Pascal/Volta Unified Memory
  - ❑ CPU/GPU coherence
  - ❑ System-wide atomic operations
  - ❑ No need to re-launch kernel or CPU threads
  - ❑ Possibility of CPU and GPU working on the same frontier



# Benefits of Collaboration

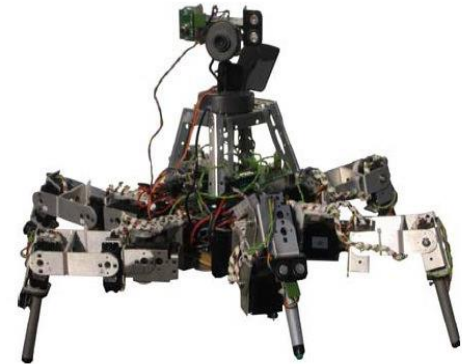
- **SSSP** performs more computation than BFS



Single Source Shortest Path  
(up to 22% improvement over GPU only)

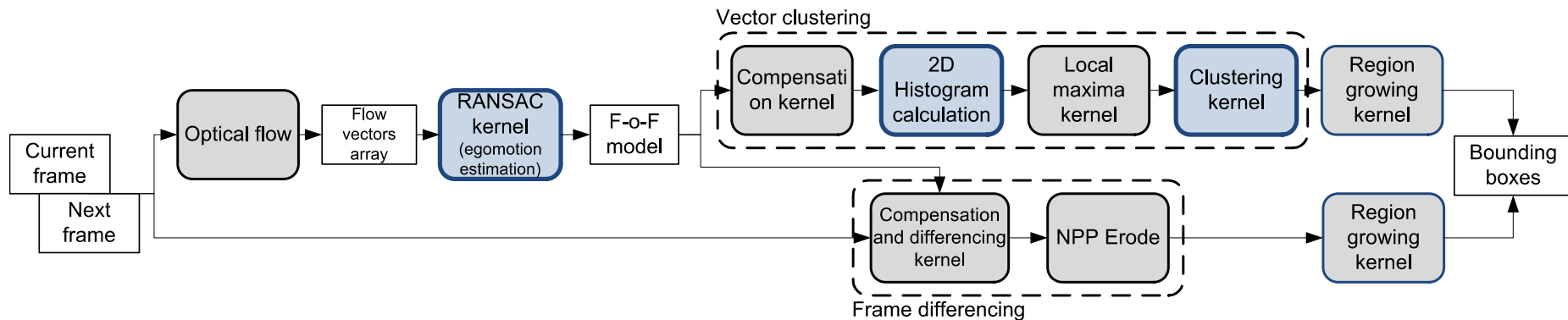
# Egomotion Compensation and Moving Objects Detection (I)

- Hexapod robot OSCAR
  - Rescue scenarios
  - Strong egomotion on uneven terrains



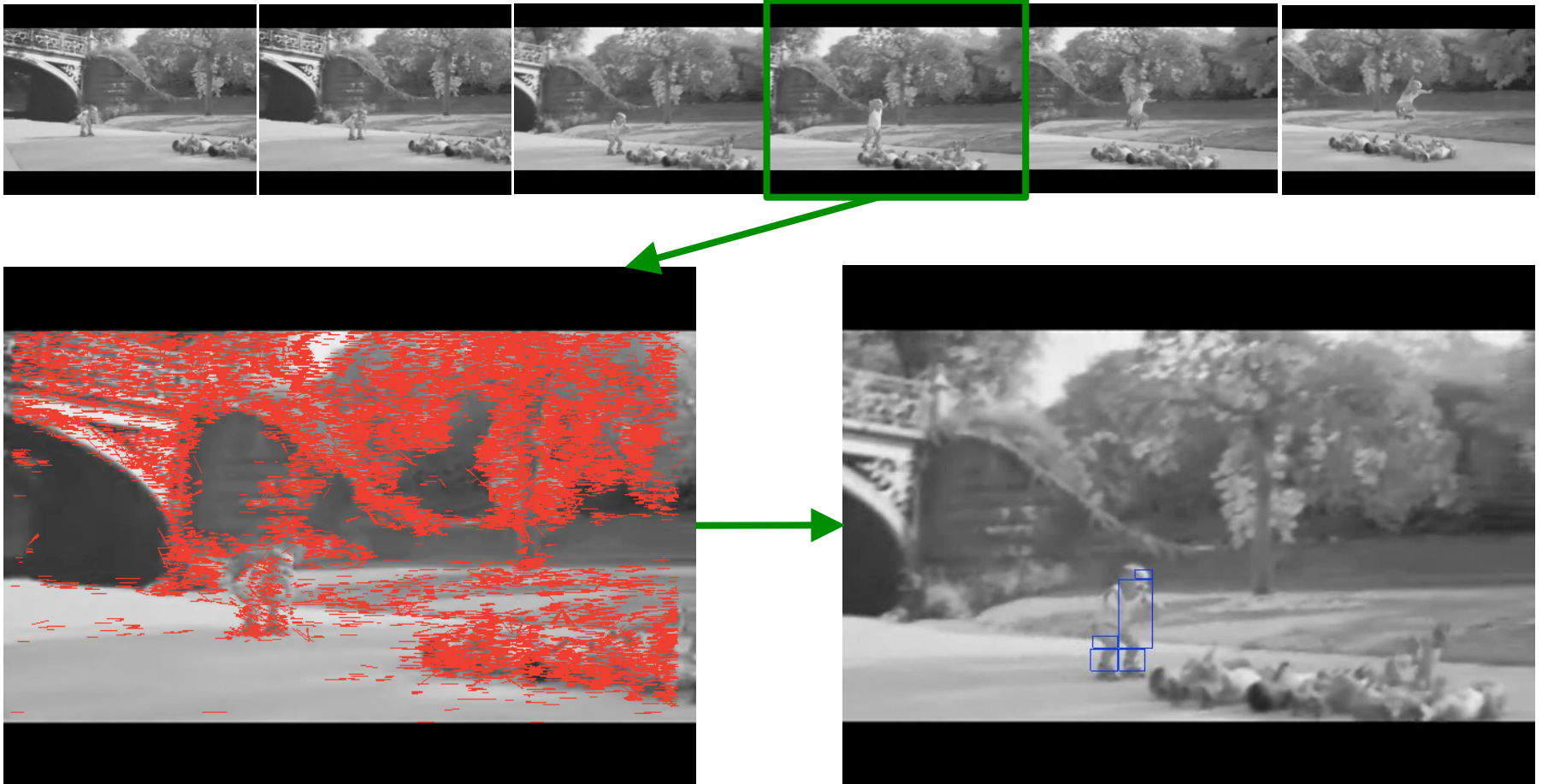
## ■ Algorithm

- Random Sample Consensus (RANSAC): F-o-F model



# Egomotion Compensation and Moving Objects Detection (II)

Fast moving object in strong egomotion scenario detected by vector clustering

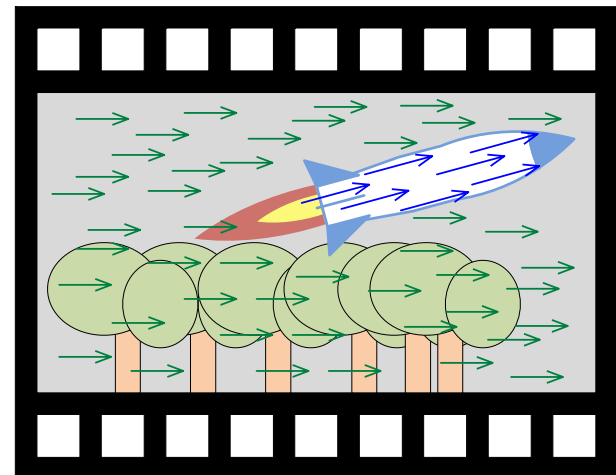


# SISD and SIMD phases

## ■ RANSAC (Fischler *et al.* 1981)

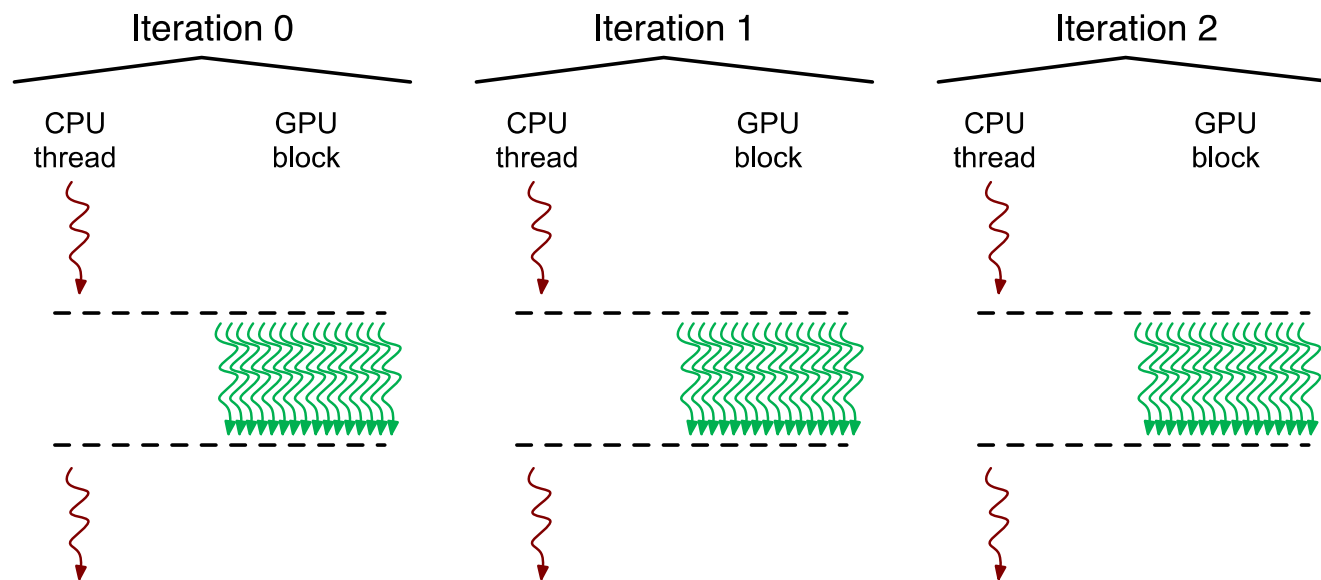
```
While (iteration < MAX_ITER){  
    Fitting stage (Compute F-o-F model)           // SISD phase  
  
    Evaluation stage (Count outliers)             // SIMD phase  
  
    Comparison to best model                     // SISD phase  
  
    Check if best model is good enough and iteration >= MIN_ITER    // SISD phase  
}
```

- ❑ Fitting stage picks two flow vectors randomly
- ❑ Evaluation generates motion vectors from F-o-F model, and compares them to real flow vectors



# Collaborative Implementation

- Randomly picked vectors: **Iterations are independent**
  - We assign one iteration to one CPU thread and one GPU block



# Chai Benchmark Suite (I)

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- Collaboration patterns
  - 8 data partitioning benchmarks
  - 3 coarse-grain task partitioning benchmarks
  - 3 fine-grain task partitioning benchmarks

<https://chai-benchmarks.github.io>



# Chai Benchmark Suite (II)

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Collaboration Pattern		Short Name	Benchmark
Data Partitioning		BS	Bézier Surface
		CEDD	Canny Edge Detection
		HSTI	Image Histogram (Input Partitioning)
		HSTO	Image Histogram (Output Partitioning)
		PAD	Padding
		RSCD	Random Sample Consensus
		SC	Stream Compaction
		TRNS	In-place Transposition
Task Partitioning	Fine-grain	RSCT	Random Sample Consensus
		TQ	Task Queue System (Synthetic)
		TQH	Task Queue System (Histogram)
	Coarse-grain	BFS	Breadth-First Search
		CEDT	Canny Edge Detection
		SSSP	Single-Source Shortest Path

# Computer Architecture

## Lecture 26: GPU Programming

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Fall 2022

06 January 2023