



**CSCI-GA.3033-004**

# **Graphics Processing Units (GPUs): Architecture and Programming**

## **CUDA**

### **Advanced Techniques 1**

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# Some Refreshing Exercises!

- Suppose registers and shared memory capacities were not an issue. When is it still beneficial to put values fetched from memory into the shared memory?

# Some Refreshing Exercises!

- Assume a kernel is launched with 1000 blocks. Each block has 512 threads.
  - If a variable is declared as local in the kernel, how many versions will be created throughout the lifetime of the kernel?
  - How about if the variable is created as shared?

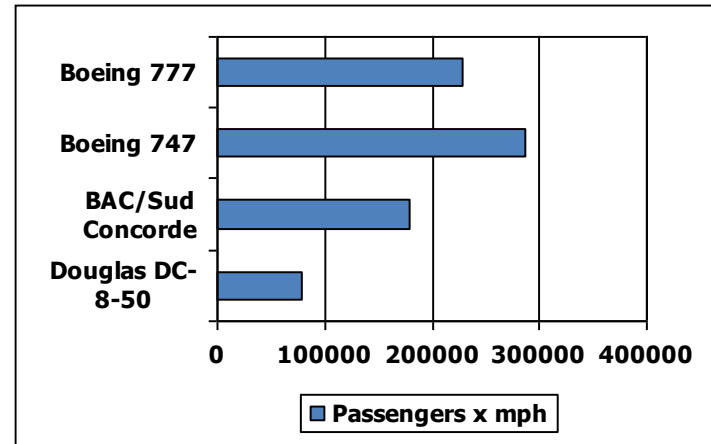
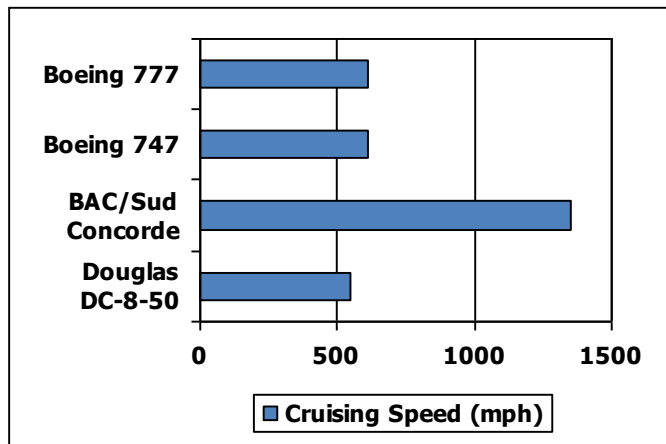
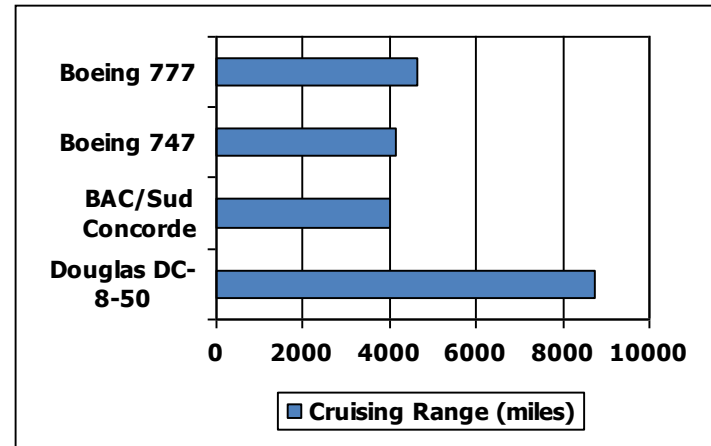
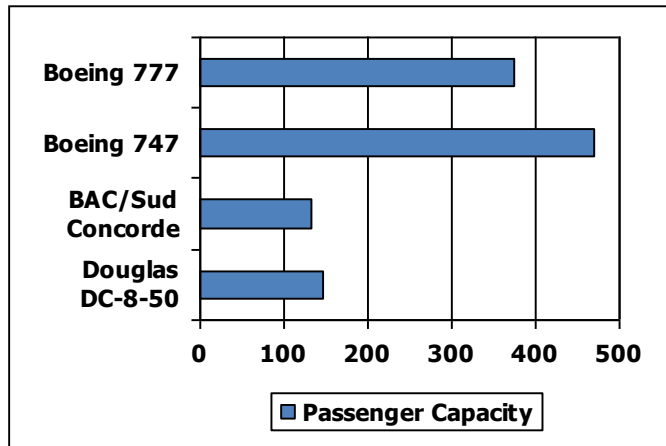
# Some Refreshing Exercises!

- A kernel contains 36 floating point operations and 7 32-bit word global memory accesses per thread. For each of the following device properties, indicate whether this kernel is compute- or memory-bound.
  - Peak FLOPS = 200 GFLOPS, Peak Memory Bandwidth = 100 GB/s
  - Peak FLOPS = 300 GFLOPS, Peak Memory Bandwidth = 250 GB/s

# A Note About Performance

# Defining Performance

- Which airplane has the best performance?

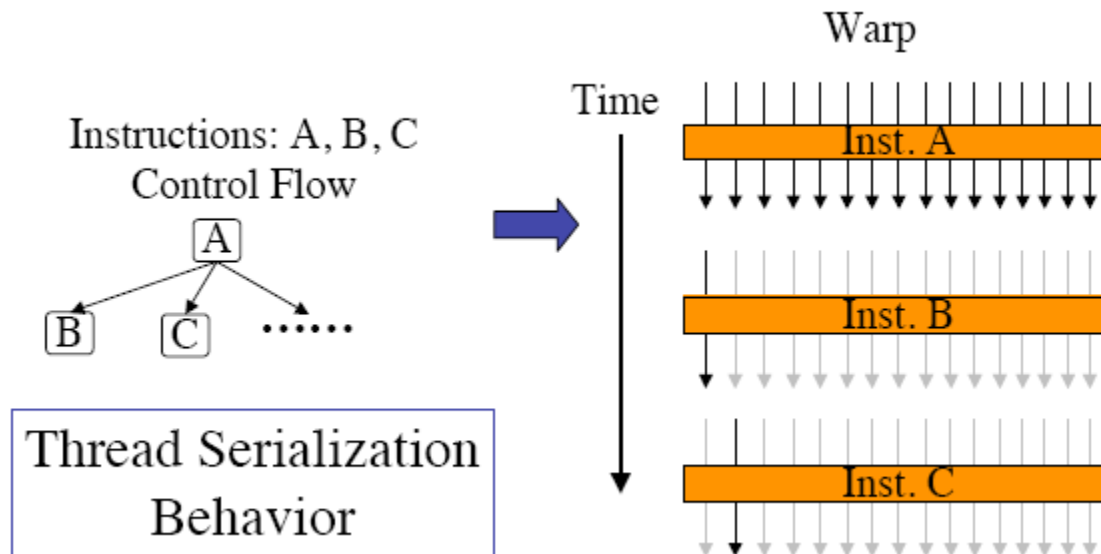


# Performance Considerations

- There are many **hardware constraints**.
- Depending on the application, different constraints may dominate.
- We can improve performance of an application by **trading one resource usage for another**.

# Performance Issue: Thread Diversion

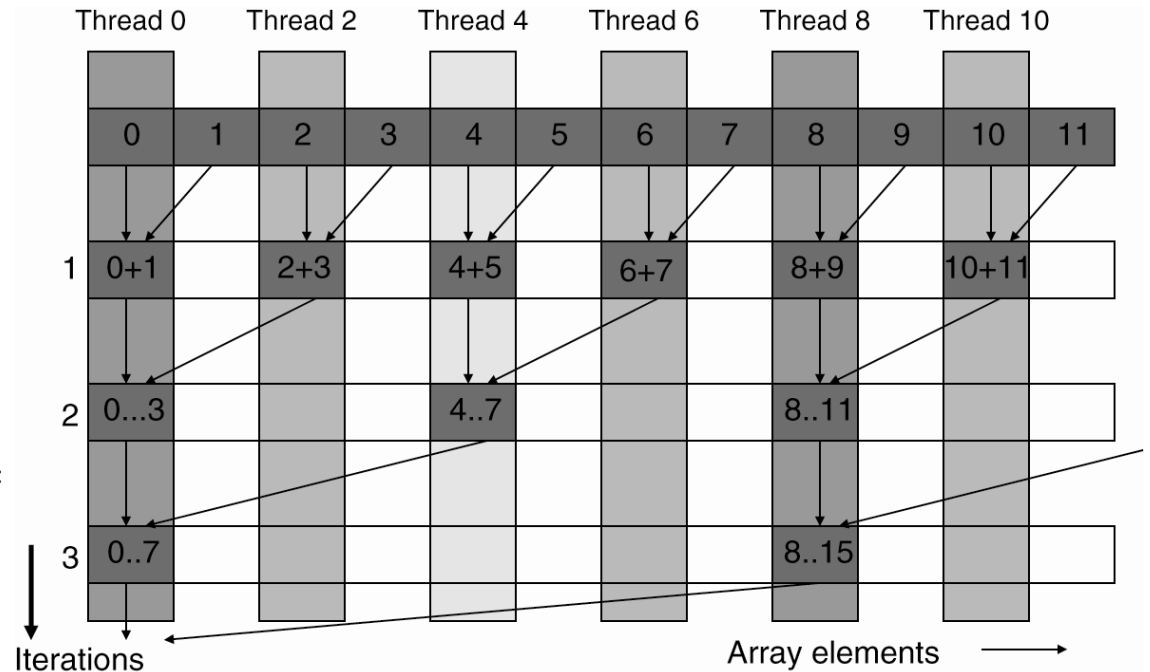
- Works well when all threads in a warp follow the same control-flow
- Performance loss due to thread diversion





# Performance Issue: Thread Diversion

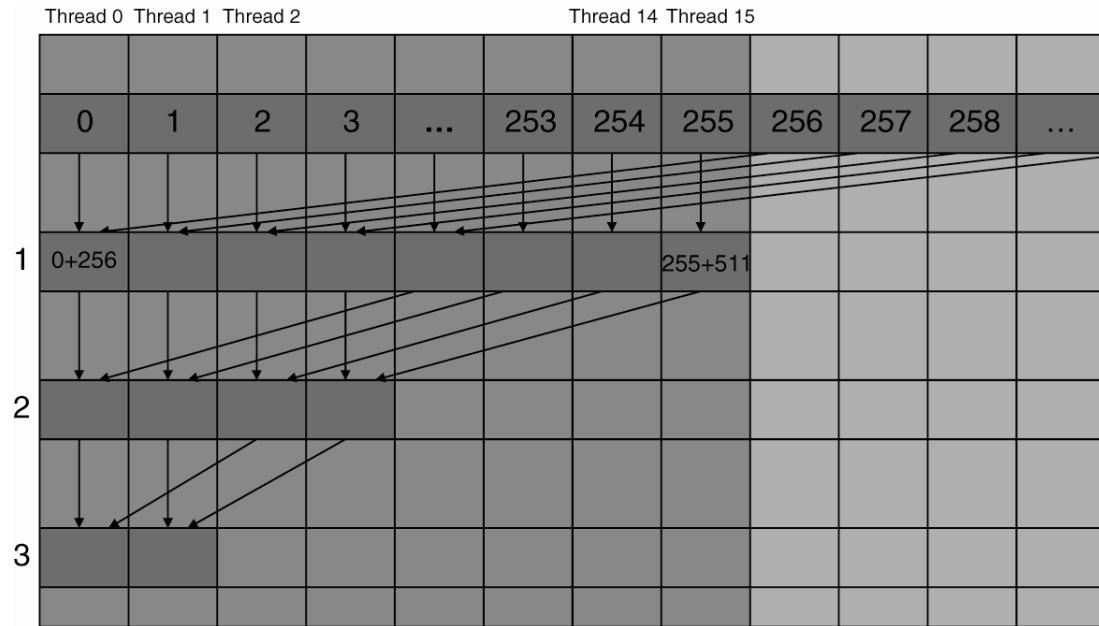
```
1. __shared__ float partialSum[]  
2. unsigned int t = threadIdx.x;  
3. for (unsigned int stride = 1;  
4.     stride < blockDim.x; stride *= 2)  
5. {  
6.     syncthreads();  
7.     if (t % (2*stride) == 0)  
8.         partialSum[t] += partialSum[t+stride];  
9. }
```



Example: Sum Reduction Kernel

# Performance Issue: Thread Diversion

```
1. __shared__ float partialSum[];  
2. unsigned int t = threadIdx.x;  
3. for (unsigned int stride = blockDim.x >> 1;  
4.     stride > 0; stride >>= 1)  
5. {  
6.     __syncthreads();  
7.     if (t < stride)  
8.         partialSum[t] += partialSum[t + stride];  
9. }
```



Why is this version better than the previous one?

Example: Sum Reduction Kernel

# Performance Issue: Global Memory

- Typical application: process massive amount of data within short period of time
  - From global memory
  - large amount + short period = huge bandwidth requirement
- Two main challenges regarding global memory:
  - Long latency
  - Relatively limited bandwidth

# Dealing With Global Memory: **TILING**

- We have seen this before
- Make use of shared memory available in SMs to reduce trips to global memory

# Dealing With Global Memory: Coalescing

- To more effectively move data from global memory to shared memory and registers
- For best results: can be used with tiling
- Global memory:
  - DRAM
  - Reading a bit is slow
  - So memory is implemented to read several bits in parallel

# Dealing With Global Memory: Coalescing

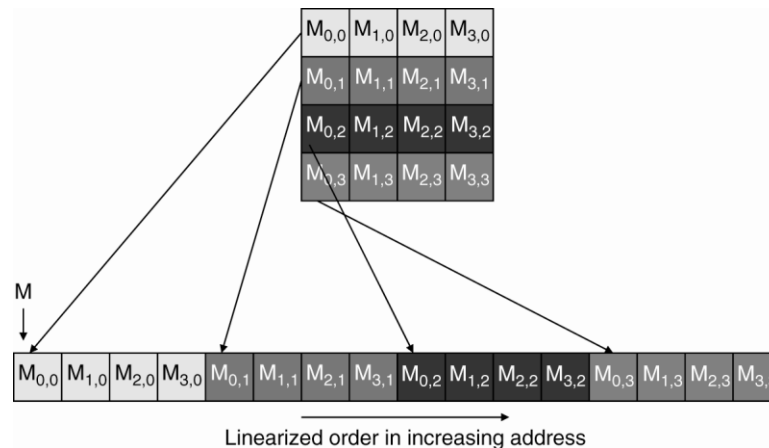
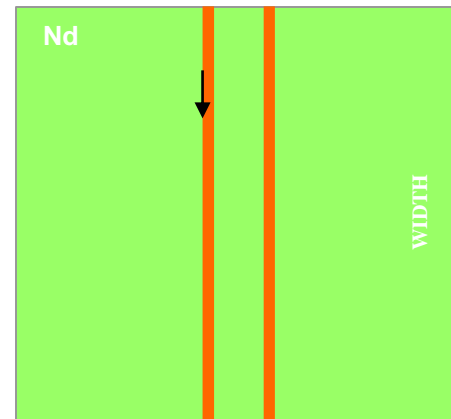
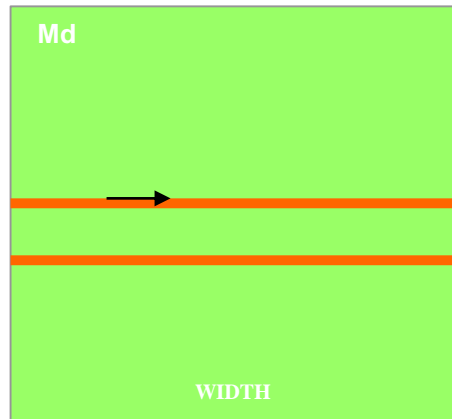
- If an application can make use of data from **multiple consecutive locations**, the DRAM can supply the data in much higher rate.
- Kernel must arrange its data access accordingly
- When all threads in a warp execute a load instruction:
  - The hardware detects whether the addresses are consecutive
  - The hardware combines (coalesces) all accesses in a consolidated access to consecutive DRAM locations

# Dealing With Global Memory: Coalescing

Not coalesced

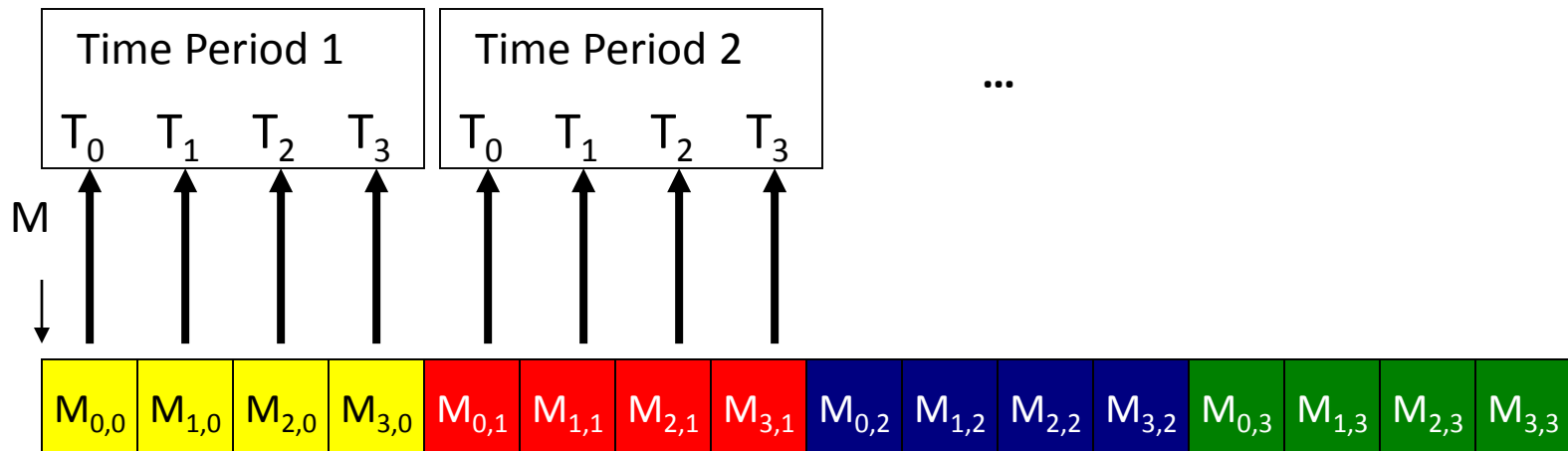
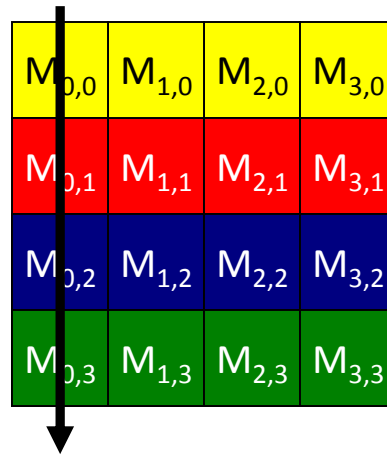
coalesced

Same warp { Thread 1  
Thread 2



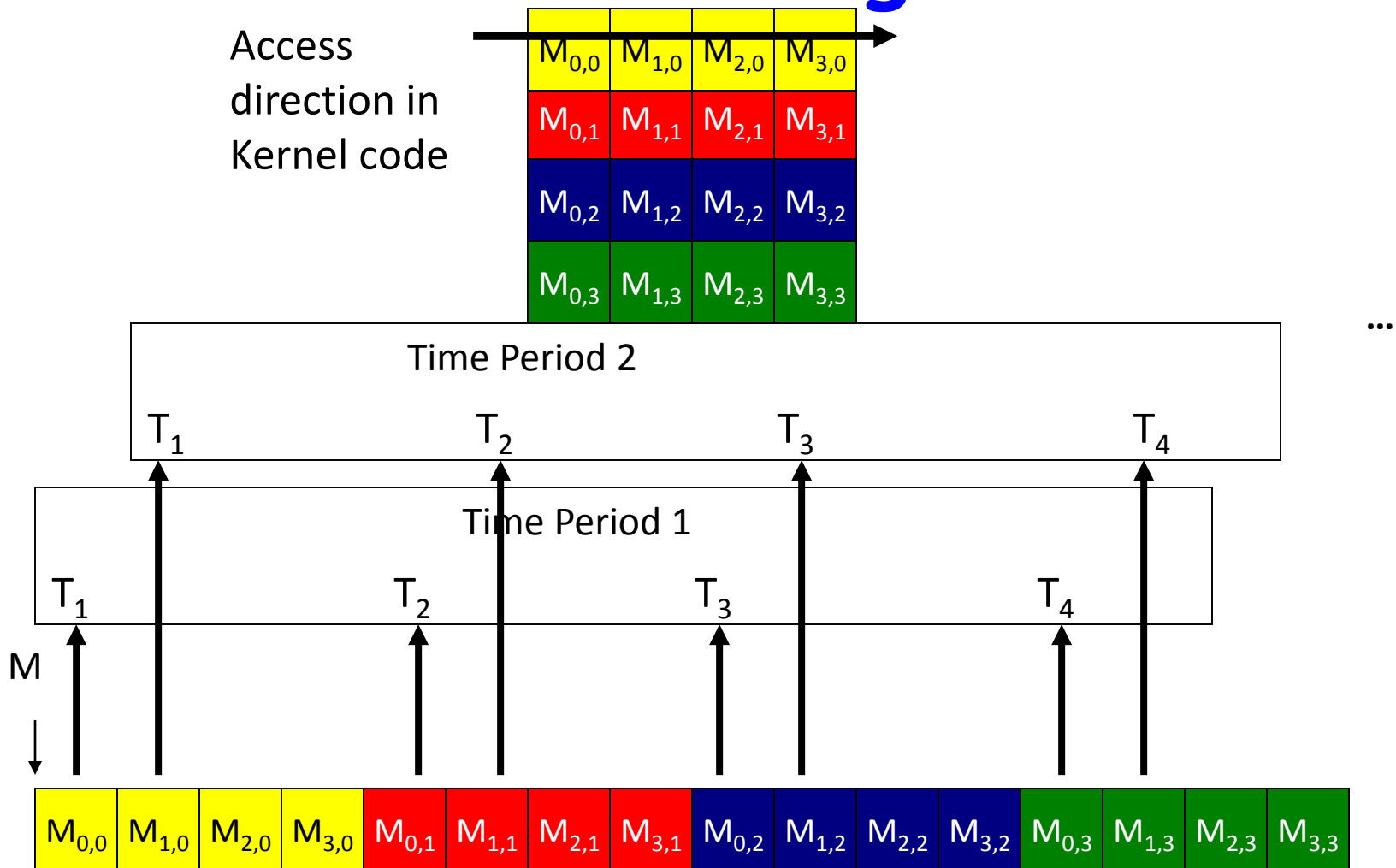
# Dealing With Global Memory: Coalescing

**Favorable** access  
direction in  
Kernel code  
of one thread



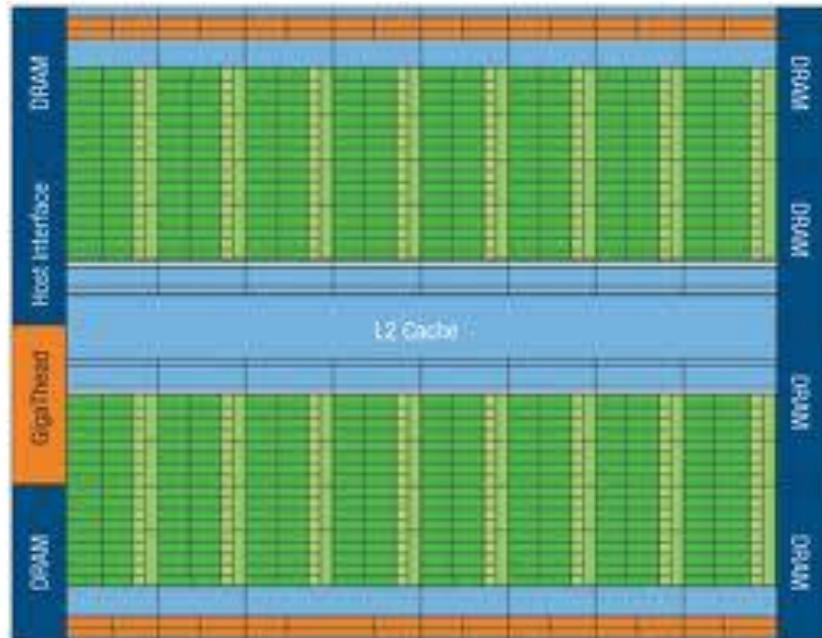


# Dealing With Global Memory: Coalescing



# Dealing With Global Memory: Cache

- Fermi (and later GPUs) have cache for global memory
- Caches automatically coalesce most of kernel access patterns



# Dealing With Global Memory: Prefetching

Prefetch next data elements while consuming the current data elements.

This increases the number of independent instructions between memory accesses and consumers

```
Loop {  
  
  Load current tile to shared  
    memory  
  
  __syncthreads()  
  
  Compute current tile  
  
  __syncthreads()  
}
```

A Without prefetching

```
Load first tile from global memory into  
  registers  
  
Loop {  
  Deposit tile from registers to shared  
    memory  
  __syncthreads()  
  
  Load next tile from global memory into  
    registers  
  
  Compute current tile  
  
  __syncthreads()  
}
```

B With prefetching

# Performance Issue: SM Resources

- Execution resources in SM include:
  - registers
  - block slots
  - thread slots
- There is an interaction among the resources that you must take into account.

# Performance Issue: SM Resources

**Example:** Assume G80 executing the matrix multiplication with 16x16 thread blocks (8 block slots, 768 thread slots, 8192 registers)

**If a thread needs 10 registers then:**

- A block needs  $10 \times 16 \times 16 = 2560$  registers
- 3 blocks  $\rightarrow$  7680 registers (under the 8192 limit)
- We can't add another block (will make it 10240)
- 3 blocks  $\times$  256 threads/block = 768 (within limit)

By using 1 extra variable the program saw a 1/3 reduction in warp parallelism  
 **$\rightarrow$  performance cliff**

**Assume the programmer declares one more auto var:**

- $11 \times 16 \times 16 = 2816$  registers per block
- 3 blocks  $\rightarrow 3 \times 2816 = 8448$  (above limit)
- SM reduces #blocks by 1  $\rightarrow$  5632 registers required
- This reduces the number of threads in SM to  $2 \times 256 \rightarrow 512$

# Performance Issue: SM Resources

**Example:** Still with G80:

- An instruction takes 4 cycles
- Assume 4 independent instructions between global memory load and its use
- Global memory latency is 200 cycles

To keep execution units fully utilized:

We need to have  $200/(4 \times 4) = 14$  warps

Assume an extra register allows the programmer to use a transformation to increase independent instructions from 4 to 8, then:

- Now we need  $200/(4 \times 8) = 7$  warps
- Blocks reduced from 3 to 2 -> warps reduced from 24 to 16
- Still we can fully utilize execution units

**Trading thread-level  
parallelism with increased  
thread performance**

# Performance Issue: Instruction Mix

```
for (int k = 0; k < BLOCK_SIZE; ++k)
    Pvalue += Ms[ty][k] * Ns[k][tx];
```

**Is the above code efficient?**

- Extra instructions to update loop counter
- Extra instructions for conditional branch at the end of each iteration
- Using k to access matrices incurs address arithmetic instructions.
- All of the above compete with the floating-point calculations for limited instruction processing bandwidth.

**2 FP arithmetic**

**2 address arithmetic instructions**

**1 loop branch instructions**

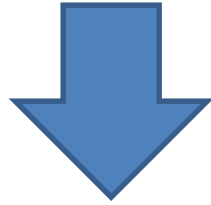
**1 loop increment instructions**



**only 1/3 instructions are FP operations**

# Performance Issue: Instruction Mix

```
for (int k = 0; k < BLOCK_SIZE; ++k)  
    Pvalue += Ms[ty][k] * Ns[k][tx];
```



```
Pvalue += Ms[ty][0] * Ns[0][tx] + ...  
          Ms[ty][15] * Ns[15][tx];
```

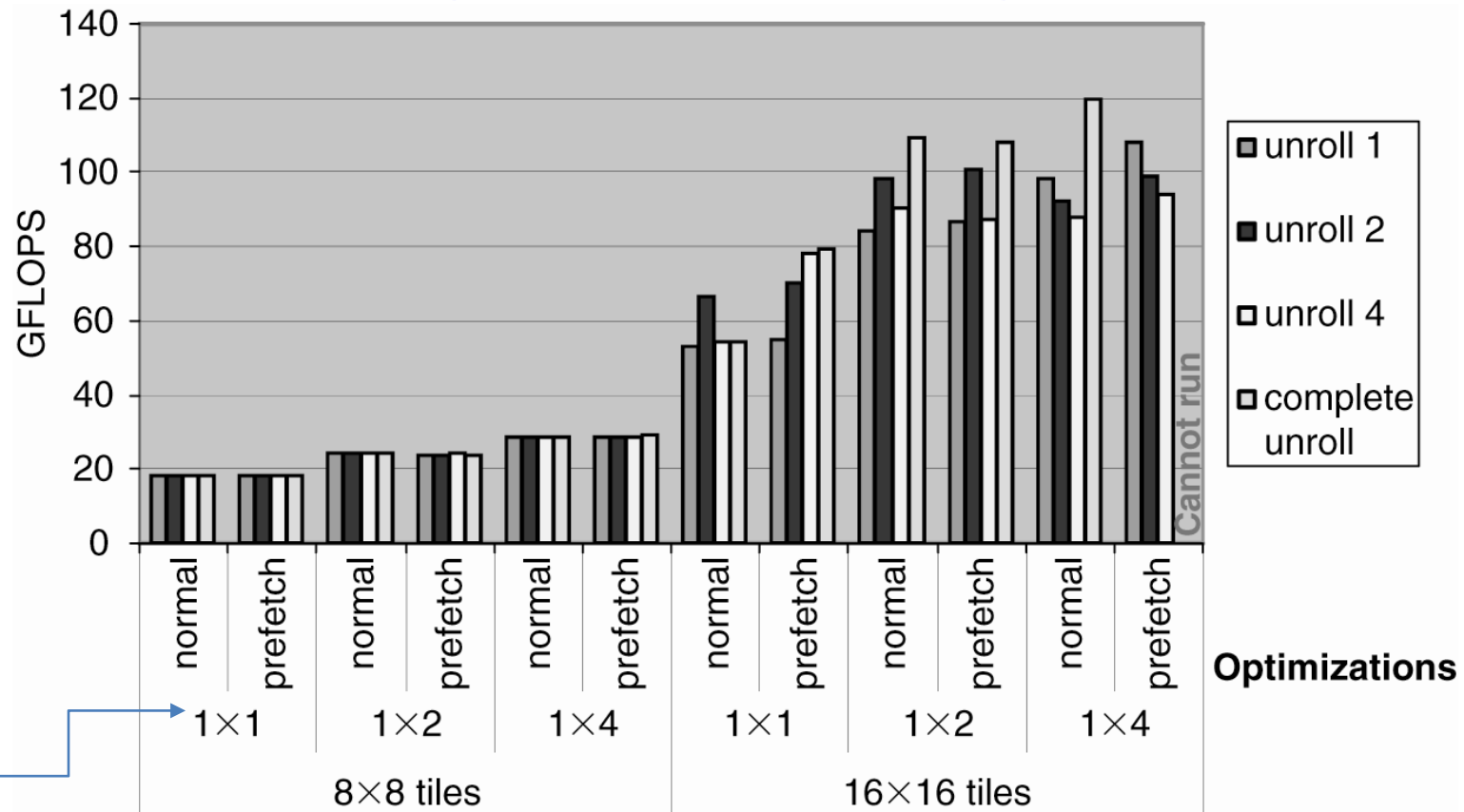
**Loop unrolling**



# Performance Issue: Thread Granularity

- Algorithmic decision
- It is often advantageous to put more work into each thread and use fewer threads
  - When redundant work exists between threads
  - Example: Let a thread compute 2 tiles
- + Less redundant work
- + Potentially more independent instructions
- More resources requirements

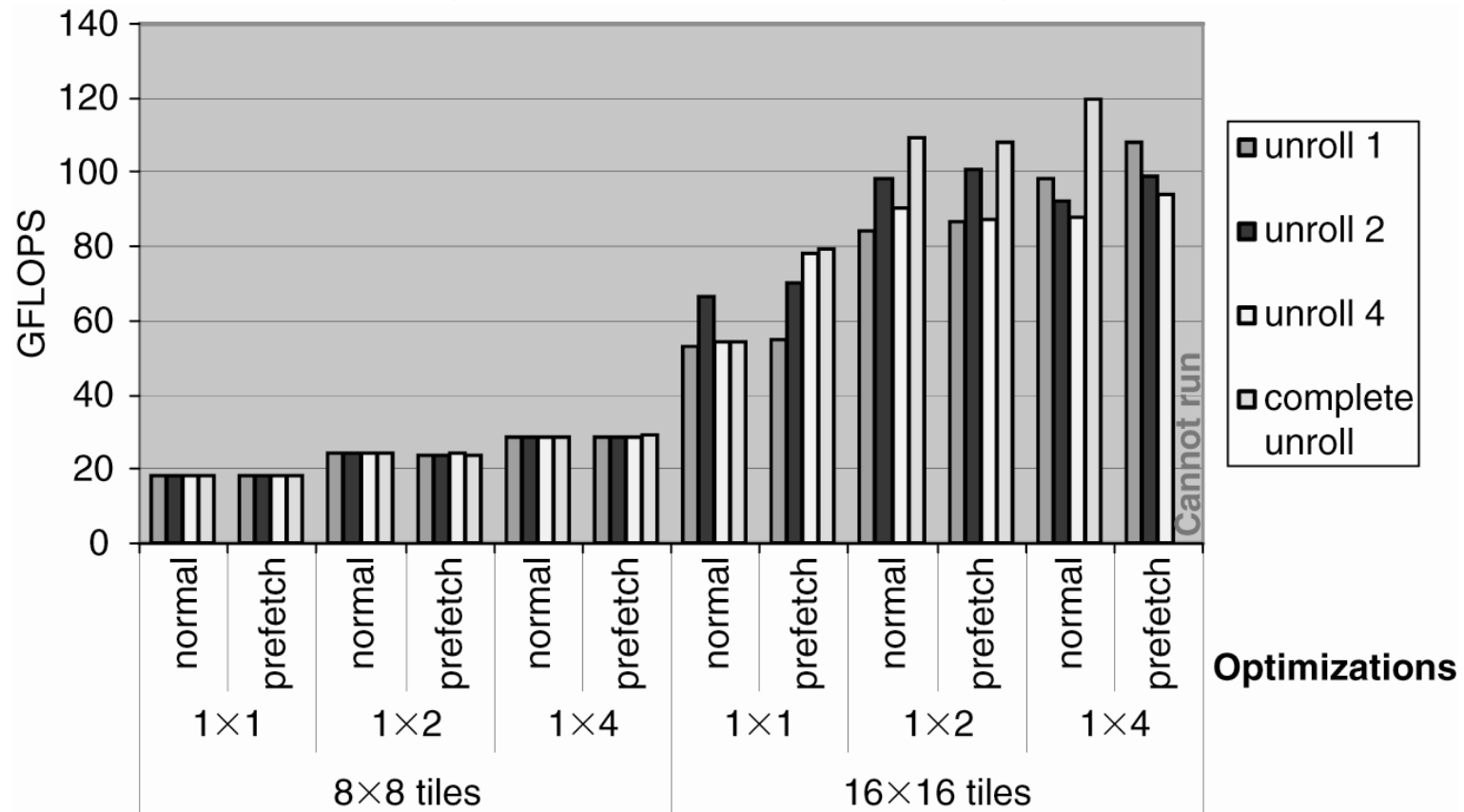
# Putting It All Together



Thread Granularity:

- normal
- merging 2 blocks
- merging 4 blocks

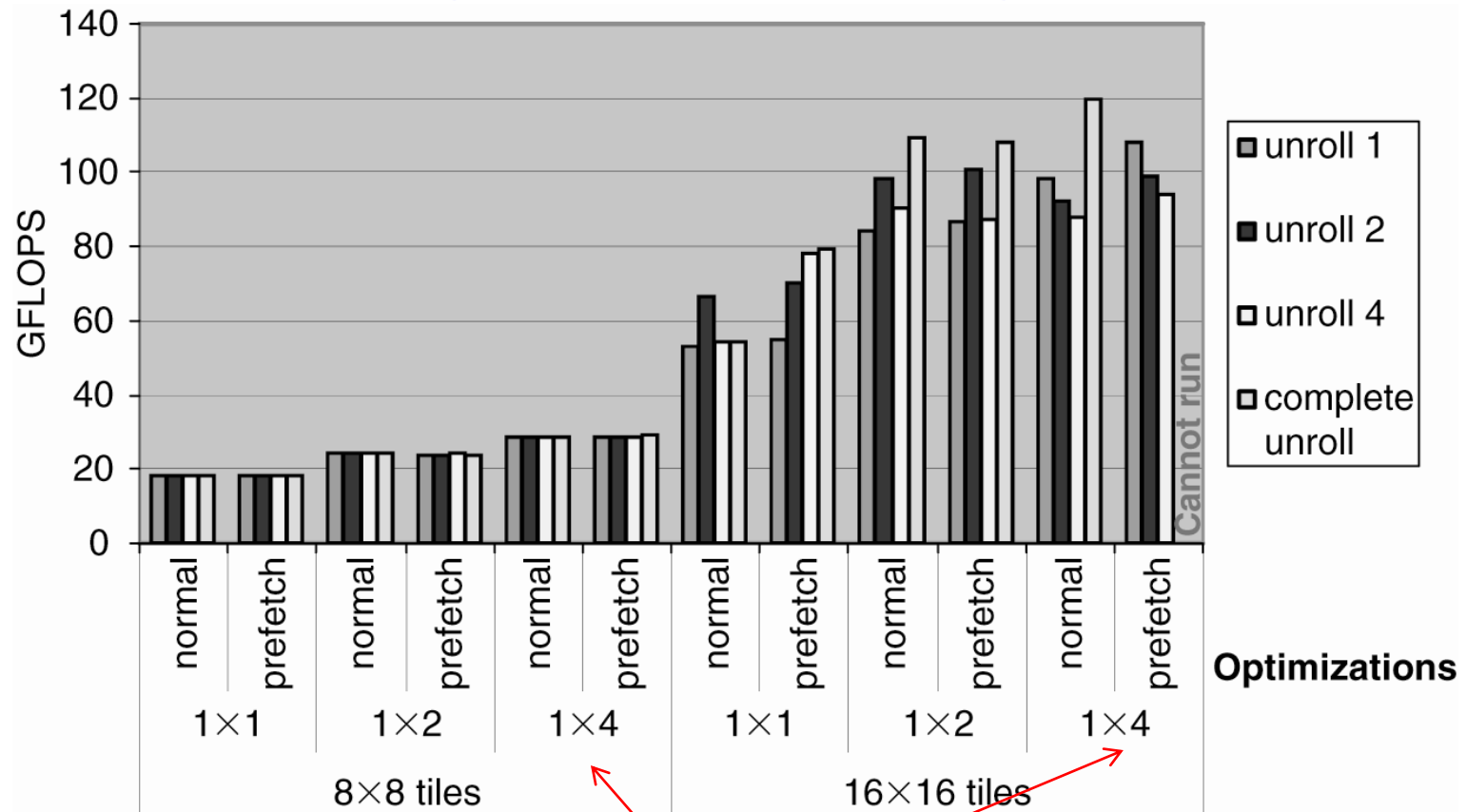
# Putting It All Together



**Until tile reaches 16x16 neither loop unrolling nor data prefetch helps.**

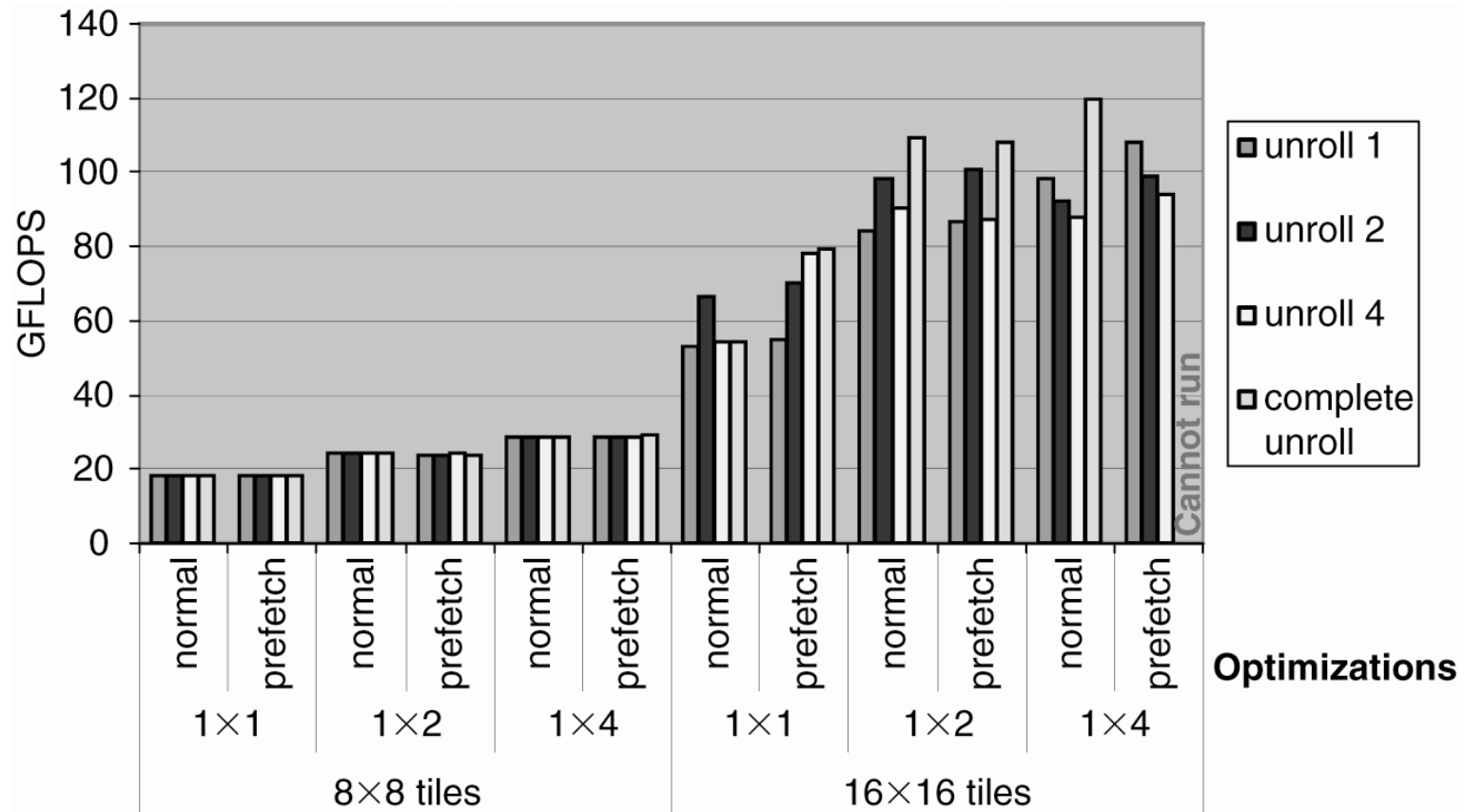
For small tile size, global memory bandwidth severely limits performance.

# Putting It All Together



**Granularity adjustment can reduce global memory access.**

# Putting It All Together



**Data prefetching becomes less beneficial as thread granularity increases.**

# About Algorithms for GPUs

- When designing an algorithm for GPU, the two main characteristics that determine its performance are:
  1. How much parallelism is available
  2. How much data must move through the memory hierarchy
- Then when you move from algorithm to code you need to take hardware constraints into account.

A note about compilation ...  
And some useful tools!

# NVCC device specific switches

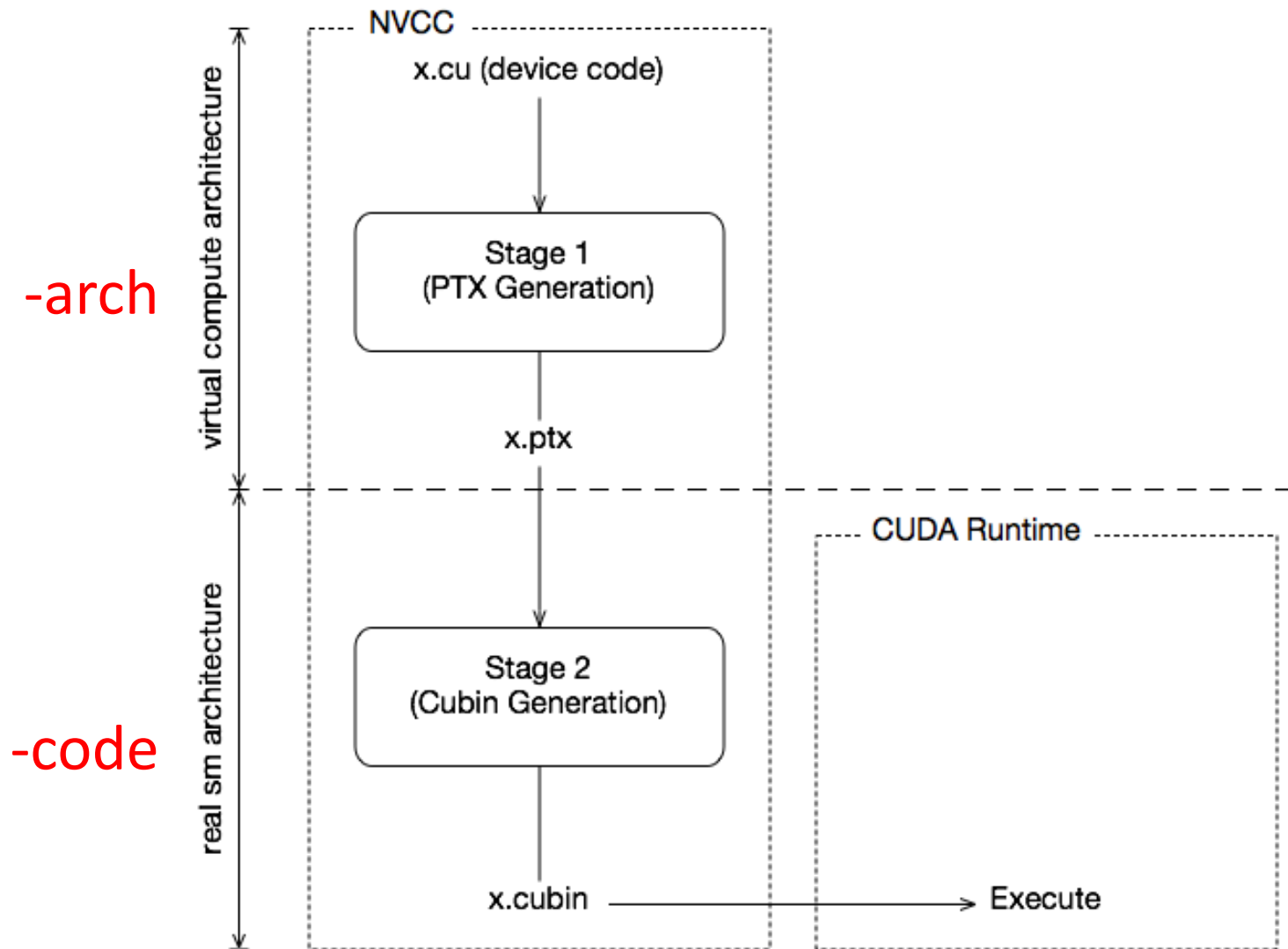
- `-arch` : controls the "virtual" architecture that will be used for the generation of the PTX code.
- `-code` : specifies the actual device that will be targeted by the cubin binary.

Virtual Architecture ( <code>-arch</code> )	Streaming Multipr. code ( <code>-code</code> )	Feature Enabled
compute_10	sm_10	Basic features
compute_11	sm_11	global memory atomic operations
compute_12	sm_12	shared memory atomic operations and vote instructions
compute_13	sm_13	double precision floating point support
compute_20	sm_20	Fermi support
	sm_21	SM structure changes (e.g more cores)
compute_30	sm_30	Kepler support
compute_35	sm_35	Dynamic parallelism (enables recursion)
compute_50	sm_50	Maxwell support



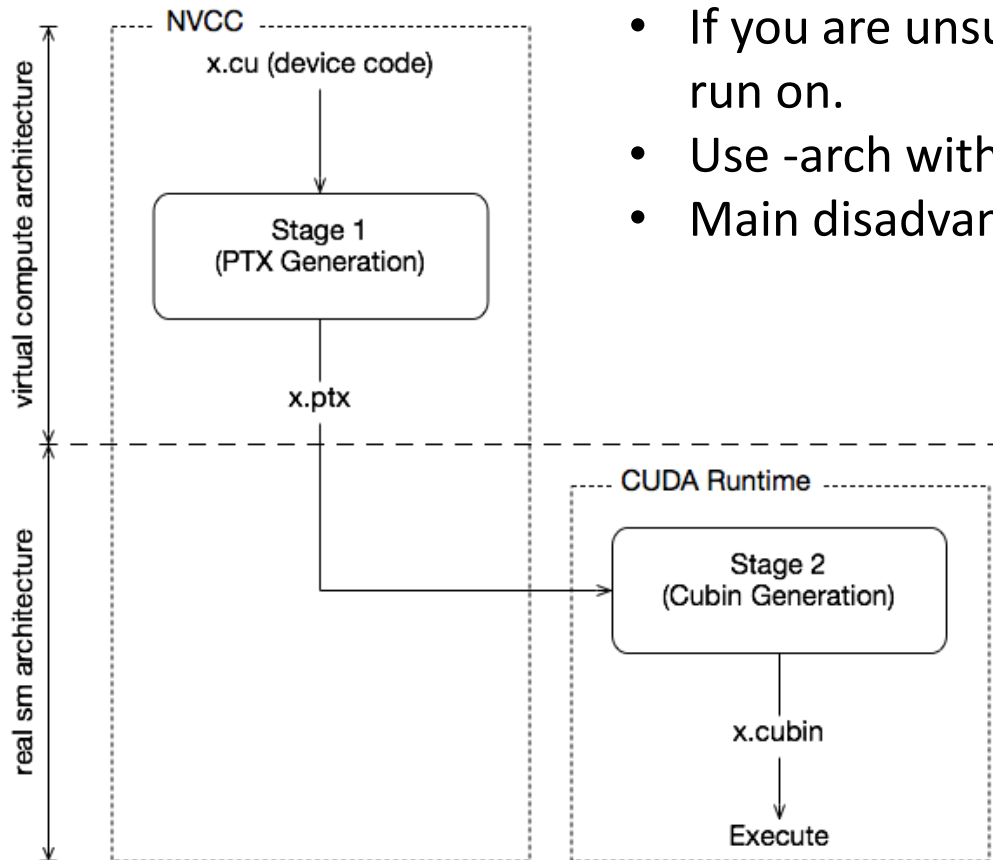
# sm\_xy

- x is the GPU generation number
- y is the version within that generation
- **Binary compatibility** of GPU applications is not guaranteed across different generations.
  - Example: a CUDA application that has been compiled for a Fermi GPU will very likely not run on a Kepler GPU (and vice versa).
- This is why nvcc relies on a two stage compilation model for ensuring **application compatibility** with future GPU generations.



Source: CUDA compiler driver nvcc manual (NVIDIA website)

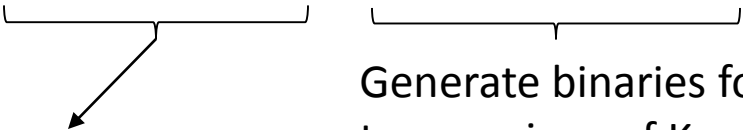
# JIT Compilation



- If you are unsure which exact GPU the code will run on.
- Use `-arch` without `-code`
- Main disadvantage: slower startup

# Fatbinaries

```
nvcc x.cu -arch=compute_30 -code=compute_30,sm_30,sm_35
```



generate PTX  
and keep it in the  
binary generated,  
for JIT on future GPUs

Generate binaries for  
two versions of Kepler

At runtime, the CUDA driver will select the most appropriate translation when the device function is launched.

Till now we have single virtual architecture and several real architectures.  
How about several virtual architectures?

# --generate-code

```
nvcc x.cu \
```

```
--generate-code arch=compute_20,code=sm_20 \
```

```
--generate-code arch=compute_20,code=sm_21 \
```

```
--generate-code arch=compute_30,code=sm_30
```

# The Default

`nvcc x.cu`



is equivalent to

`nvcc x.cu -arch=compute_20 -code=sm_20,compute_20`

# nvcc

- Some nvcc features: `--ptxas-options=-v`
  - Print the smem, register and other resource usages
- Generates CUDA binary file: `nvcc -cubin`
  - cubin file is the cuda executable

# Dealing with binary files

	cuobjdump	nvdiasm
Extract ptx and extract and disassemble cubin from the following input files: <ul style="list-style-type: none"><li>▶ Host binaries<ul style="list-style-type: none"><li>▶ Executables</li><li>▶ Object files</li><li>▶ Static libraries</li></ul></li><li>▶ External fatbinary files</li></ul>	Yes	No
Control flow analysis and output	No	Yes
Advanced display options	No	Yes



# nvprof

- CUDA profiler: profiling data from the command line

```
$ nvprof [nvprof_args] <app> [app_args]
```

- To profile a region of the application:
  1. `#include <cuda_profiler_api.h>`
  2. in the host function surround the region with:
    - `cudaProfilerStart()`
    - `cudaProfilerStop()`
  3. `nvcc myprog.cu`
  4. `nvprof --profile-from-start-off ./a.out`

# nvprof summary mode (default)

```
$ nvprof dct8x8
```

```
===== Profiling result:
```

Time(%)	Time	Calls	Avg	Min	Max	Name
49.52	9.36ms	101	92.68us	92.31us	94.31us	CUDAKernel2DCT(float*, float*, int)
37.47	7.08ms	10	708.31us	707.99us	708.50us	CUDAKernel1DCT(float*,int, int,int)
3.75	708.42us	1	708.42us	708.42us	708.42us	CUDAKernel1IDCT(float*,int,int,int)
1.84	347.99us	2	173.99us	173.59us	174.40us	CUDAKernelQuantizationFloat()
1.75	331.37us	2	165.69us	165.67us	165.70us	[CUDA memcpy DtoH]
1.41	266.70us	2	133.35us	89.70us	177.00us	[CUDA memcpy HtoD]
1.00	189.64us	1	189.64us	189.64us	189.64us	CUDAKernelShortDCT(short*, int)
0.94	176.87us	1	176.87us	176.87us	176.87us	[CUDA memcpy HtoA]
0.92	174.16us	1	174.16us	174.16us	174.16us	CUDAKernelShortIDCT(short*, int)
0.76	143.31us	1	143.31us	143.31us	143.31us	CUDAKernelQuantizationShort(short*)
0.52	97.75us	1	97.75us	97.75us	97.75us	CUDAKernel2IDCT(float*, float*)
0.12	22.59us	1	22.59us	22.59us	22.59us	[CUDA memcpy DtoA]

# nvprof trace mode

```
$ nvprof --print-gpu-trace dct8x8
```

```
==== Profiling result:
```

Start	Duration	Grid Size	Block Size	Regs	SSMem	DSMem	Size	Throughput	Name
167.82ms	176.84us	-	-	-	-	-	1.05MB	5.93GB/s	[CUDA memcpy HtoA]
168.00ms	708.51us	(64 64 1)	(8 8 1)	28	512B	0B	-	-	CUDAKernel1DCT(float*, ...)
168.95ms	708.51us	(64 64 1)	(8 8 1)	28	512B	0B	-	-	CUDAKernel1DCT(float*, ...)
169.74ms	708.26us	(64 64 1)	(8 8 1)	28	512B	0B	-	-	CUDAKernel1DCT(float*, ...)
170.53ms	707.89us	(64 64 1)	(8 8 1)	28	512B	0B	-	-	CUDAKernel1DCT(float*, ...)
171.32ms	708.12us	(64 64 1)	(8 8 1)	28	512B	0B	-	-	CUDAKernel1DCT(float*, ...)
172.11ms	708.05us	(64 64 1)	(8 8 1)	28	512B	0B	-	-	CUDAKernel1DCT(float*, ...)
172.89ms	708.38us	(64 64 1)	(8 8 1)	28	512B	0B	-	-	CUDAKernel1DCT(float*, ...)
173.68ms	708.31us	(64 64 1)	(8 8 1)	28	512B	0B	-	-	CUDAKernel1DCT(float*, ...)
174.47ms	708.15us	(64 64 1)	(8 8 1)	28	512B	0B	-	-	CUDAKernel1DCT(float*, ...)
175.26ms	707.95us	(64 64 1)	(8 8 1)	28	512B	0B	-	-	CUDAKernel1DCT(float*, ...)
176.05ms	173.87us	(64 64 1)	(8 8 1)	27	0B	0B	-	-	CUDAKernelQuantization (...)
176.23ms	22.82us	-	-	-	-	-	1.05MB	45.96GB/s	[CUDA memcpy DtoA]

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GPU-trace mode provides a timeline of all activities taking place on the GPU in chronological order.

Print individual kernel invocations  
and sort them in chronological order

Print CUDA runtime/driver  
API trace

```
$ nvprof --print-gpu-trace --print-api-trace dct8x8
```

=====  
Profiling result:

Start	Duration	Grid Size	Block Size	Regs	SSMem	DSMem	Size	Throughput	Name
167.82ms	176.84us	-	-	-	-	-	1.05MB	5.93GB/s	[CUDA memcpy HtoA]
167.81ms	2.00us	-	-	-	-	-	-	-	cudaSetupArgument
167.81ms	38.00us	-	-	-	-	-	-	-	cudaLaunch
167.85ms	1.00ms	-	-	-	-	-	-	-	cudaDeviceSynchronize
168.00ms	708.51us	(64 64 1)	(8 8 1)	28	512B	0B	-	-	CUDAKernel11DCT(float*, ...)
168.86ms	2.00us	-	-	-	-	-	-	-	cudaConfigureCall
168.86ms	1.00us	-	-	-	-	-	-	-	cudaSetupArgument
168.86ms	1.00us	-	-	-	-	-	-	-	cudaSetupArgument
168.86ms	1.00us	-	-	-	-	-	-	-	cudaSetupArgument
168.87ms	0ns	-	-	-	-	-	-	-	cudaSetupArgument
168.87ms	24.00us	-	-	-	-	-	-	-	cudaLaunch
168.89ms	761.00us	-	-	-	-	-	-	-	cudaDeviceSynchronize
168.95ms	708.51us	(64 64 1)	(8 8 1)	28	512B	0B	-	-	CUDAKernel11DCT(float*, ...)

`nvcc --devices x --events y ./a.out`

- *x*: device number in case of multi-GPU
- *y*: event name
  - Gives very useful information, such as:
    - number of global memory loads, stores, ...
    - number of global memory coalesced
    - branch divergences

```
$ nvprof --devices 0 --events branch,divergent_branch dct8x8
```

# Revisiting Shared Memory

# Till now

- We know that we have to use `__shared__`
- But, you leave it to the runtime system to distribute shared memory among blocks. Can you change that?
- What if you don't know how much shared memory we need before execution?

# This is what we know

```
__global__ void staticShared(int  
*d, int n)  
{  
    __shared__ int array[64];  
    int t = threadIdx.x;  
    s[t] = d[t];  
}
```

What if you don't  
know this size  
beforehand?






# Dynamic Shared Memory

```
__global__ void DynamicShared(int
*d, int n)
{
    extern __shared__ int array[];
    int t = threadIdx.x;
    s[t] = d[t];
}

main()
{
    ...
    DynamicShared<<<100, 100, N*sizeof(int)>>>(int *, int);
}
```

Step 1:  
A different declaration  
with no size in the kernel

Step 2:  
The 3<sup>rd</sup> argument of kernel  
launch is the shared memory  
allocation size per thread block  
in bytes.



# What if you need more than one array in the shared memory?

- Declare a single extern unsized array (as in the previous slide).
- Use pointers into it to divide it into multiple arrays.
- Example:

```
extern __shared__ int s[];
```

```
int *integerData = s;          // nI ints
```

```
float *floatData = (float*)&integerData[nI]; // nF floats
```

```
char *charData = (char*)&floatData[nF];    // nC chars
```

# Potential Performance Loss: Shared Memory Bank Conflicts

- Shared memory is divided into equally sized memory modules called **banks**.
- Banks can be accessed simultaneously.
- Shared memory accesses that span  $b$  distinct banks yield an effective bandwidth that is  $b$  times as high as the bandwidth of when accesses map to the same bank.
- Exceptions are: broadcast and multicast.

# Potential Performance Loss: Shared Memory Bank Conflicts

- Shared memory banks are organized such that successive 32-bit words are assigned to successive banks.
- Bandwidth is 32 bits per bank per clock cycle.
- Warp size is 32

# Potential Performance Loss: Shared Memory Bank Conflicts

- Now, depending on compute capability:
  - 1.x: the number of banks is 16. A shared memory request for a warp is split into one request for the first half of the warp and one request for the second half of the warp.
  - 2.x: the number of banks is 32. A shared memory request for a warp is not split.
  - 3.x: have configurable bank size

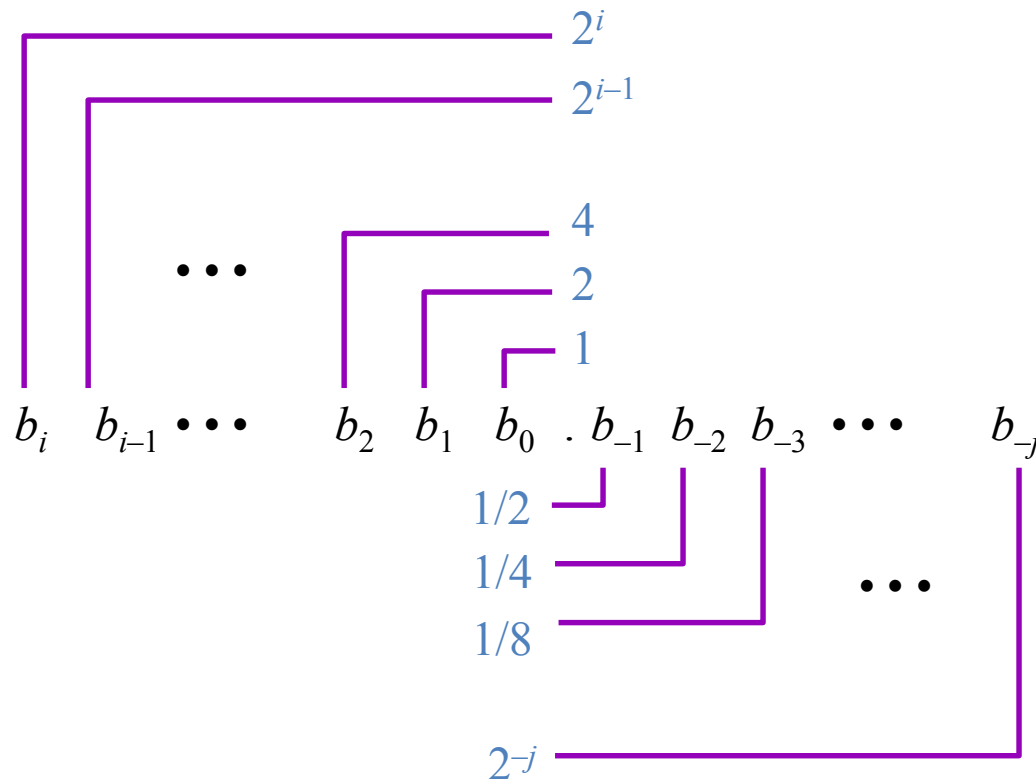
# cudaDeviceSetSharedMemConfig(*i*)

- *i* above can be:
  - cudaSharedMemBankSizeFourByte ← default
  - cudaSharedMemBankSizeEightByte

# Floating Points

# Importance of Floating Points

- Many graphics operations are floating point operations
- GPU performance is measure in GFLOPS







Turing Award 1989 to William Kahan for design of the  
IEEE Floating Point Standards 754 (binary) and 854  
(decimal)

# What is Excel doing?

A1:

1.33333333333333330000	=4/3
------------------------	------

---

# Floating Point

- We need a way to represent
  - numbers with fractions, e.g., 3.1416
  - very small numbers, e.g., .0000000001
  - very large numbers, e.g.,  $3.15576 \times 10^9$
- Representation:
  - sign, exponent, mantissa:  $(-1)^{\text{sign}} \times \text{mantissa} \times 2^{\text{exponent}}$
  - more bits for mantissa gives more accuracy
  - more bits for exponent increases range
- IEEE 754 floating point standard:
  - single precision: 8 bit exponent, 23 bit mantissa
  - double precision: 11 bit exponent, 52 bit mantissa

# IEEE 754 floating-point standard

- Leading "1" bit of significand is implicit (called **hidden 1 technique**, except when  $\text{exp} = -127$ )
- Exponent is "biased" to make sorting easier
  - all 0s is smallest exponent
  - all 1s is largest exponent
  - bias of 127 for single precision and 1023 for double precision
  - summary:  $(-1)^{\text{sign}} \times (1 + \text{significand}) \times 2^{\text{exponent} - \text{bias}}$
- Example:
  - decimal:  $-.75 = -\left(\frac{1}{2} + \frac{1}{4}\right)$
  - binary:  $-.11 = -1.1 \times 2^{-1}$
  - floating point:  $\text{exponent} = 126 = 01111110$
  - IEEE single precision: 10111111010000000000000000000000

# More about IEEE floating Point Standard

Single Precision:

$$(-1)^{\text{sign}} \times (1 + \text{mantissa}) \times 2^{\text{exponent} - 127}$$

The variables shown in red are the numbers stored in the machine

# Floating Point Example

what is the decimal equivalent of

1      01110110      10110000...0

# Based on **exp** we have 3 encoding schemes

- $\text{exp} \neq 0..0 \text{ or } 11...1 \rightarrow$  normalized encoding
- $\text{exp} = 0... 000 \rightarrow$  denormalized encoding
- $\text{exp} = 1111...1 \rightarrow$  special value encoding
  - $\text{frac} = 000...0$
  - $\text{frac} = \text{something else}$

# 1. Normalized Encoding

- Condition:  $\text{exp} \neq 000\dots 0$  and  $\text{exp} \neq 111\dots 1$

referred to as Bias

- Exponent is:  $E = \text{Exp} - (2^{k-1} - 1)$ ,  $k$  is the # of exponent bits

- Single precision:  $E = \text{exp} - 127$
- Double precision:  $E = \text{exp} - 1023$

- Significand is:  $M = 1.\overbrace{\text{xxx}\dots\text{x}_2}^{\text{frac}}$
- $\text{Range}(M) = [1.0, 2.0 - \epsilon)$
- Get extra leading bit for free

$\text{Range}(E) = [-126, 127]$

$\text{Range}(E) = [-1022, 1023]$



# Normalized Encoding Example

- Value: Float  $F = 15213.0$ ;  
–  $15213_{10} = 11101101101101_2$   
–  $= 1.1101101101101_2 \times 2^{13}$

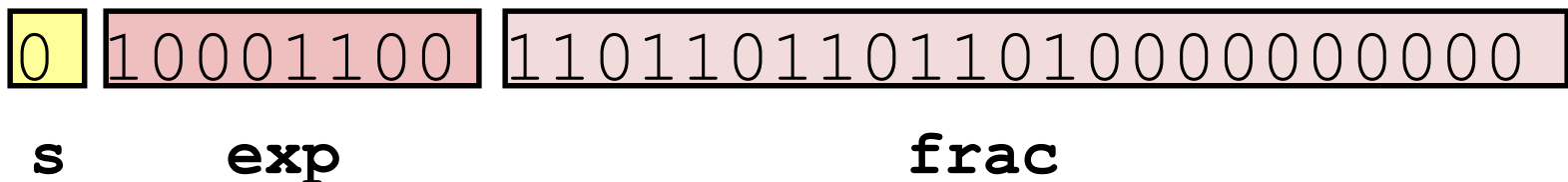
- Significand

$$\begin{array}{ll} M &= 1.\underline{1101101101101}_2 \\ \text{frac} &= \underline{1101101101101}0000000000_2 \end{array}$$

- Exponent

$$\begin{array}{llll} E &= \text{exp} - \text{Bias} &= \text{exp} - 127 &= 13 \\ \rightarrow \text{exp} &= 140 &= 10001100_2 \end{array}$$

- Result:



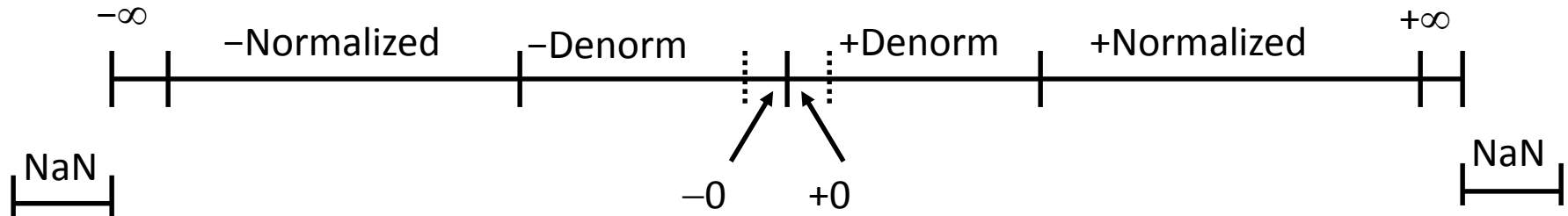
## 2. Denormalized Encoding

- Condition:  $\text{exp} = 000\dots 0$
- Exponent value:  $E = 1 - \text{Bias}$  (instead of  $E = 0 - \text{Bias}$ )
- Significand is:  $M = 0.\underbrace{\text{xxx}\dots\text{x}_2}_{\text{frac}}$  (instead of  $M = 1.\text{xxx}_2$ )
- Cases
  - **exp** = 000...0, **frac** = 000...0
    - Represents zero
    - Note distinct values: +0 and -0
  - **exp** = 000...0, **frac**  $\neq$  000...0
    - Numbers very close to 0.0

# 3. Special Values Encoding

- Condition: **exp** = 111...1
- Case: **exp** = 111...1, **frac** = 000...0
  - Represents value  $\infty$  (infinity)
  - Operation that overflows
  - E.g.,  $1.0/0.0 = -1.0/-0.0 = +\infty$ ,  $1.0/-0.0 = -\infty$
- Case: **exp** = 111...1, **frac**  $\neq$  000...0
  - Not-a-Number (NaN)
  - Represents case when no numeric value can be determined
  - E.g.,  $\text{sqrt}(-1)$ ,  $\infty - \infty$ ,  $\infty \times 0$


# Visualization: Floating Point Encodings



# Floating Point: IEEE 754

## What is the decimal equivalent of:

**10011111101010000000000000000000**



**127**  
So:

11111111

1.1010....0  
= 1.625

- Real exponent =  $127 - 127 = 0$
- There is hidden 1

*Final answer = -1.625*

# In Summary About Floating Points

exponent	mantissa	meaning
11...1	$\neq 0$	NaN
11...1	$=0$	$(-1)^S * \infty$
00...0	$\neq 0$	denormalized
00...0	$=0$	0

# Algorithm Considerations

- Non **representable** numbers are rounded
- This rounding *error* leads to different results depending on the order of operations
  - Non-repeatability makes debugging harder
- A common technique to maximize floating point arithmetic accuracy is to presort data before a reduction computation.

So..

When doing floating-point operations in parallel you have to decide:

- How much accuracy is good enough?
- Do you need single-precision or double precision?
- Can you tolerate presorting overhead, if you care about rounding errors?



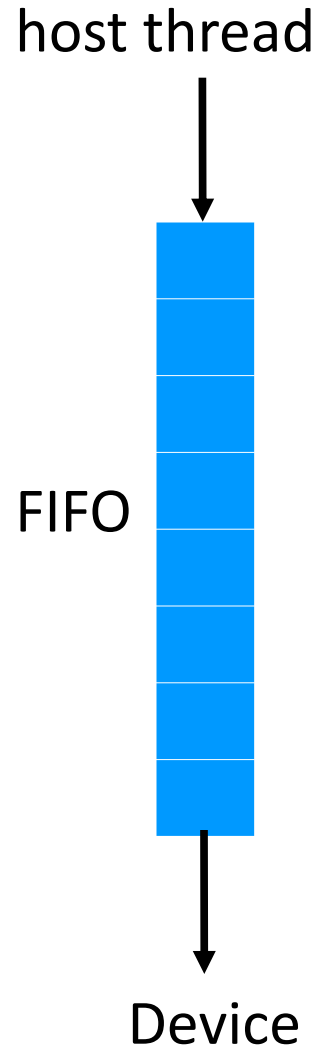
Streams

# Streams

- A sequence of operations that execute on the device in the order in which they are issued by the host code
- Operations in different streams can be interleaved and, when possible, they can even run concurrently.
- A stream can be sequence of kernel launches and host-device memory copies
- Can have several open streams to the same device at once
- Need GPUs with concurrent transfer/execution capability
- Potential performance improvement: can overlap transfer and computation

# Streams

- By default all transfers and kernel launches are assigned to stream 0
  - This means they are executed in order



# Example: Default Stream

```
cudaMemcpy(d_a, a, numBytes, cudaMemcpyHostToDevice);  
increment<<<1,N>>>(d_a);  
cudaMemcpy(a, d_a, numBytes, cudaMemcpyDeviceToHost);
```

- In the code above, from the perspective of the device, all three operations are issued to the same (default) stream and will execute in the order that they were issued.
- From the perspective of the host:
  - data transfers are blocking or synchronous transfers
  - kernel launch is asynchronous.



```
cudaMemcpy(d_a, a, numBytes, cudaMemcpyHostToDevice);  
increment<<<1,N>>>(d_a);  
anyCPUfunction();  
cudaMemcpy(a, d_a, numBytes, cudaMemcpyDeviceToHost);
```

# Example: Non-Default Stream

Non-default streams in CUDA C/C++ are **declared**, **created**, and **destroyed** in host code as follows:

```
cudaStream_t stream1;  
cudaError_t result;  
result = cudaStreamCreate(&stream1);  
result = cudaStreamDestroy(stream1);
```

To issue data transfer to non-default stream (non-blocking):

```
result = cudaMemcpyAsync(d_a, a, N, cudaMemcpyHostToDevice, stream1);
```

To launch a kernel to non-default stream:

```
increment<<<1,N,0,stream1>>>(d_a);
```

# Important

- All operations to non-default streams are non-blocking with respect to the host code.
- Sometimes you need to synchronize the host code with operations in a stream.
- You have several options:
  - `cudaDeviceSynchronize()` → blocks host
  - `cudaStreamSynchronize(stream)` → blocks host
  - `cudaStreamQuery(stream)` → does not block host

# Streams

- The amount of overlap execution between two streams depends on:
  - Device supports overlap transfer and kernel execution ( compute capability 1.1 and higher)
  - Devices supports concurrent kernel execution (compute capability 2.x and higher)
  - Device supports concurrent data transfer (compute capability 2.x and higher)
  - The order on which commands are issued to each stream

# Using streams to overlap device execution with data transfer

- Conditions to be satisfied first:
  - The device must be capable of *concurrent copy and execution*.
  - The kernel execution and the data transfer to be overlapped must both occur in *different, non-default streams*.
  - The host memory involved in the data transfer must be *pinned memory*.



# Using streams to overlap device execution with data transfer

```
for (int i = 0; i < nStreams; ++i) {  
  
    int offset = i * streamSize;  
  
    cudaMemcpyAsync(&d_a[offset], &a[offset],  
                    streamBytes,  
                    cudaMemcpyHostToDevice,  
                    stream[i]);  
  
    kernel<<.....>>(d_a, offset);  
  
    cudaMemcpyAsync(&a[offset], &d_a[offset],  
                    streamBytes,  
                    cudaMemcpyDeviceToHost,  
                    stream[i]);  
  
}
```

**So..**

- **Streams are a good way to overlap execution and transfer, hardware permits.**
- **Don't confuse kernels, threads, and streams.**

# Pinned Pages

- Allocate page(s) from system RAM (`cudaMallocHost()` or `cudaHostAlloc()`)
  - Accessible by device (but wait till next slide)
  - Cannot be paged out
  - Enables highest memory copy performance (`cudaMemcpyAsync()`)
  - Don't forget `cudaFreeHost()`;
- If too much pinned pages, overall system performance may greatly suffer.

# Host page accessible by the device??

- The pointer to the host memory is not directly transferable to device, except with:
  - `cudaHostGetDevicePointer`(void \*\* pDevice,  
void \* pHost,  
unsigned int flags)
  - flags are 0 for now
- Accessing host memory from device without explicit copy is called "zero-copy" mechanism.

# Steps for Zero-Copy

1. `cudaHostAlloc` (`void ** ptr, size_t size, unsigned int flags`)
  - flag here: `cudaHostAllocMapped`
2. `cudaHostGetDevicePointer()`
3. Then use the pointer in your kernel on device as if it is in the GPU memory

```

#include <stdio.h>
#include <cuda.h>
#include <stdlib.h>
#define N 32      // size of vectors

__global__ void add(int *a,int *b, int *c) {
    int tid = blockIdx.x * blockDim.x + threadIdx.x;
    if(tid < N) c[tid] = a[tid]+b[tid];
}

int main(int argc, char *argv[]) {
    int T = 32, B = 1;           // threads per block and blocks per grid
    int *a,*b,*c;               // host pointers
    int *dev_a, *dev_b, *dev_c; // device pointers to host memory

    cudaHostAlloc( (void**)&a, size, cudaHostAllocMapped);
    cudaHostAlloc( (void**)&b, size, cudaHostAllocMapped);
    cudaHostAlloc( (void**)&c, size, cudaHostAllocMapped );

    ... // load arrays with some numbers

    cudaHostGetDevicePointer(&dev_a, a, 0);
    cudaHostGetDevicePointer(&dev_b, b, 0);
    cudaHostGetDevicePointer(&dev_c, c, 0);

    add<<<B,T>>>>(dev_a,dev_b,dev_c);

    cudaFreeHost(a);
    cudaFreeHost(b);
    cudaFreeHost(c);

    return 0;
}

```

So..

- If the CPU program requires a lot of memory, then pinned pages is not a good idea.

# Asynchronous Execution

- Asynchronous = returns to host right-away and does not wait for device
- This includes (but not limited to):
  - Kernel launches;
  - Memory copies between two addresses to the same device memory;
  - Memory copies from host to device of a memory block of 64 KB or less;
  - Memory copies performed by functions that are suffixed with Async;



# Asynchronous Execution

- Some CUDA API calls and all kernel launches are asynchronous with respect to the host code.
- This means error-reporting is also asynchronous.

```
cudaMemcpyAsync(a_d, a_h, size, cudaMemcpyHostToDevice, 0);  
kernel<<<grid, block>>>(a_d);  
cpuFunction();
```

# Other Sources of Concurrency

- Some devices of compute capability 2.x and higher can **execute multiple kernels** concurrently.
- The maximum number of kernel launches that a device can execute concurrently is 32 on devices of compute capability 3.5 and 16 on devices of lower compute capability.
- A kernel from one CUDA context cannot execute concurrently with a kernel from another CUDA context.
- Kernels that use many textures or a large amount of local memory are less likely to execute concurrently with other kernels.
- Some devices of compute capability 2.x and higher can perform **a copy from page-locked host memory to device memory concurrently with a copy from device memory to page locked host memory.**

# Conclusions

- As we program GPUs we need to pay attention to several performance bottlenecks:
  - Branch diversion
  - Global memory latency
  - Global memory bandwidth
  - Shared memory bank conflicts
  - Communication
  - Limited resources
- We have several techniques in our arsenal to enhance performance
  - Try to make threads in the same warp follow the **same control flow**
  - **Tiling**
  - **Coalescing**
  - **Loop unrolling**
  - **Increase thread granularity**
  - **Trade one resource for another**
  - **Memory access pattern**
  - **Streams**
- Pay attention to interaction among techniques