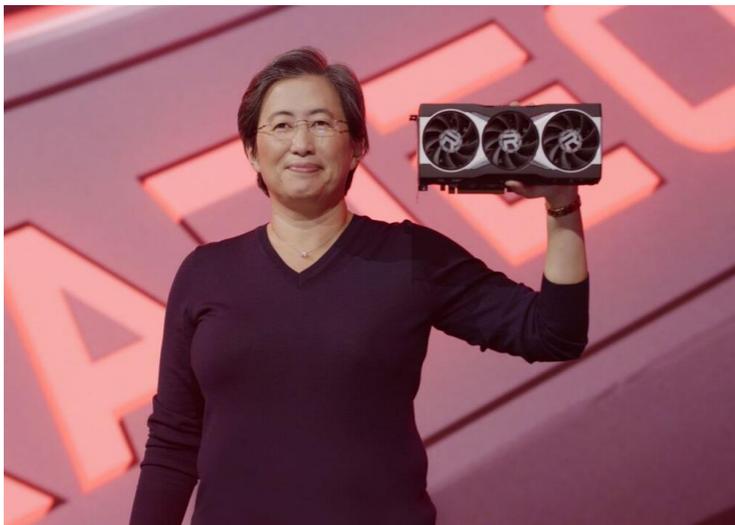


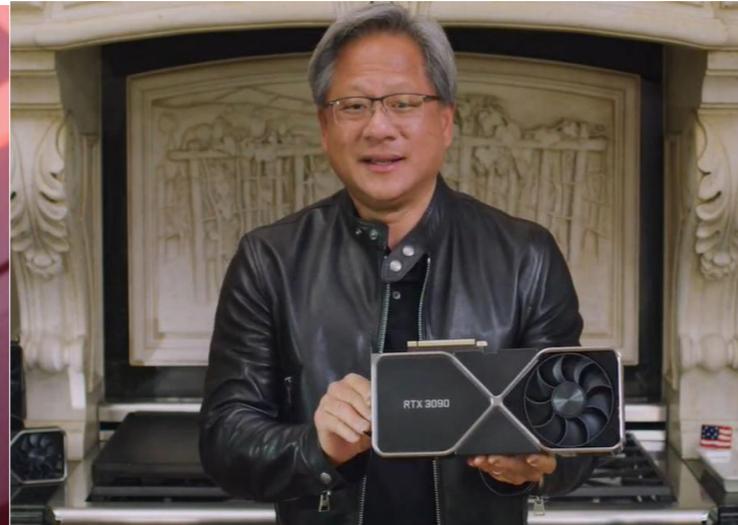
Advanced Computer Architecture

Chapter 9

Data-Level Parallel Architectures: GPUs



Lisa Su, CEO of AMD, launching the rx6000 series



Jensen Huang, CEO of NVIDIA, launching the RTX 30 Series GPUs

November 2023
Paul Kelly

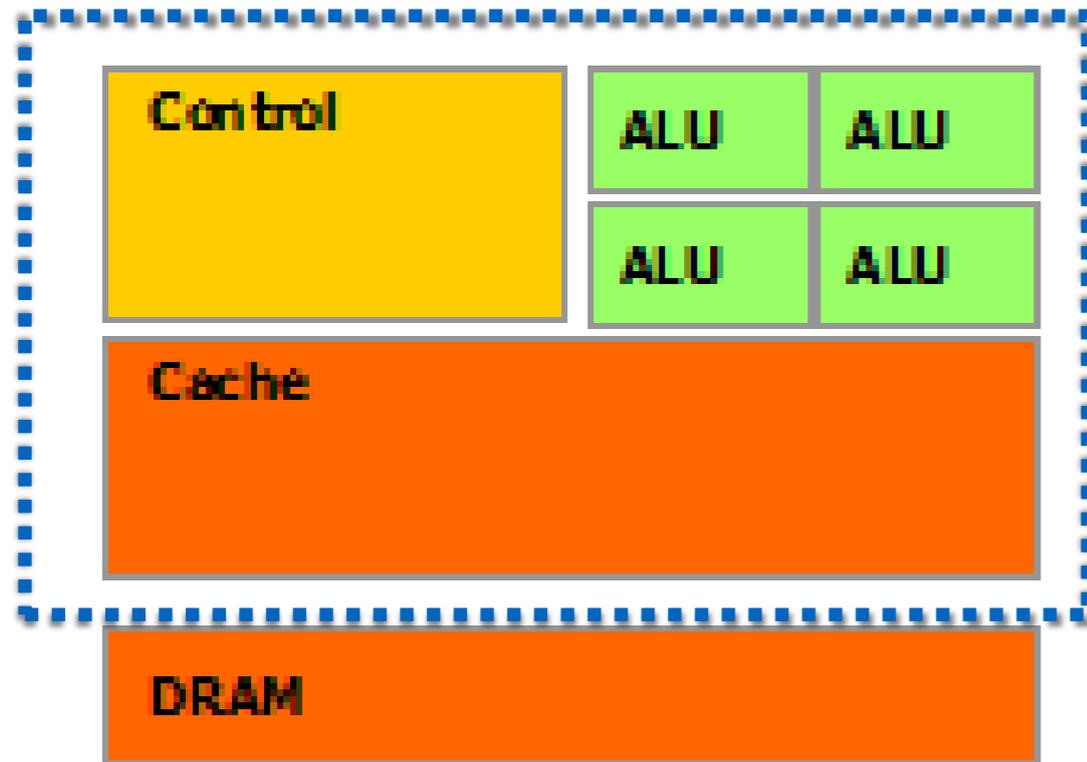
These lecture notes are partly based on:

- **Contributions to the lecture slides from Luigi Nardi (postdoc at Imperial and Stanford, now academic at Lund, Sweden), Fabio Luporini (Imperial PhD, postdoc, now CTO, DevitoCodes), and Nicolai Stawinoga (Imperial PhD, postdoc, now researcher at TU Berlin)**
- **the course text, Hennessy and Patterson's Computer Architecture (5th ed.)**

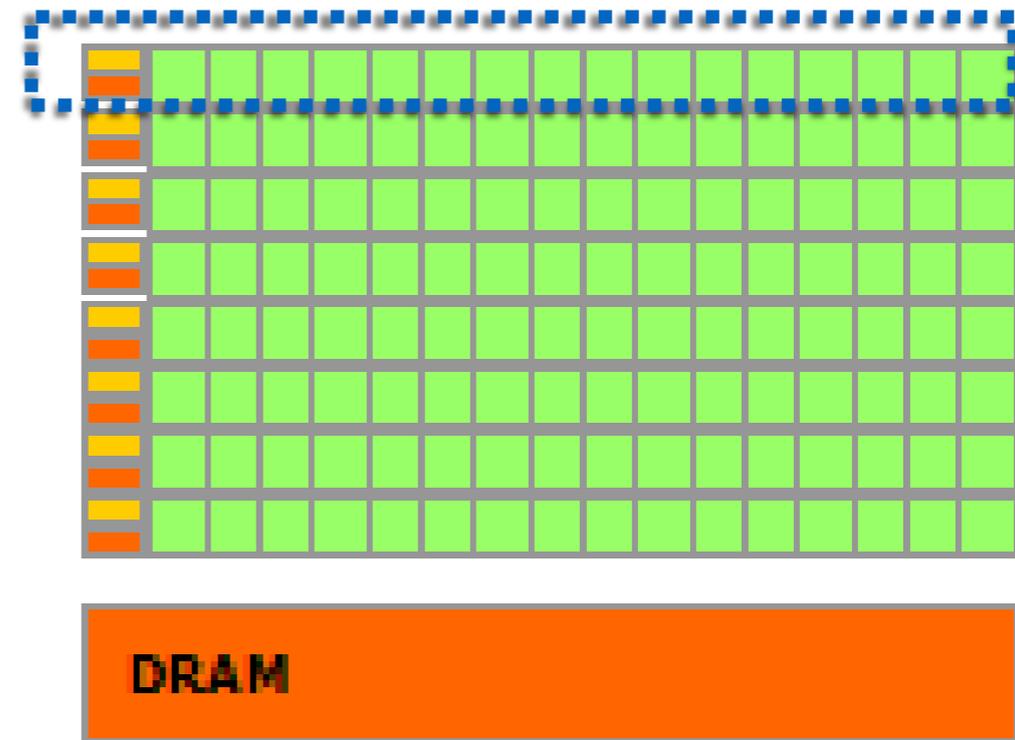
Graphics Processors (GPUs)

- Much of our attention so far has been devoted to making a single core run a single thread faster
- If your workload consists of thousands of threads, *everything* looks different:
 - *Never* speculate: there is always another thread waiting with work you *know* you have to do
 - No speculative branch execution, perhaps even no branch prediction
 - Can use FGMT or SMT to hide cache access latency, and maybe even main memory latency
 - Control is at a premium (Turing tax avoidance):
 - How to launch >10,000 threads?
 - What if they branch in different directions?
 - What if they access random memory blocks/banks?
- This is the “manycore” world
- Initially driven by the gaming market – but with many other applications

A first comparison with CPUs



CPU



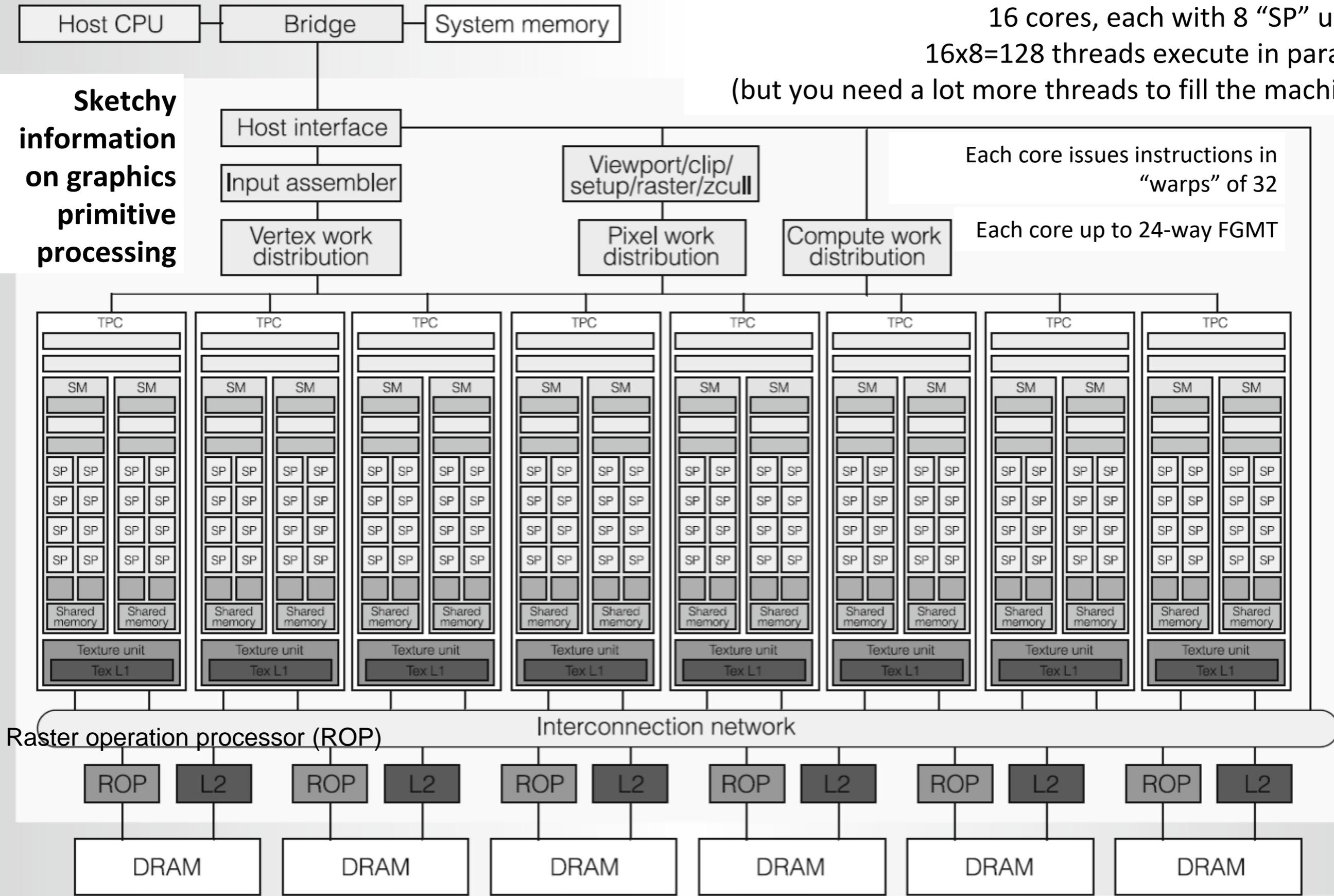
GPU

- “Simpler” cores
- Many functional units (FUs) (implementing the SIMD model)
- Much less cache per core; just thousands of threads and super-fast context switch
- Drop sophisticated branch prediction mechanisms

NVIDIA G80 (2006)

16 cores, each with 8 "SP" units
 16x8=128 threads execute in parallel
 (but you need a lot more threads to fill the machine)

**Sketchy
 information
 on graphics
 primitive
 processing**

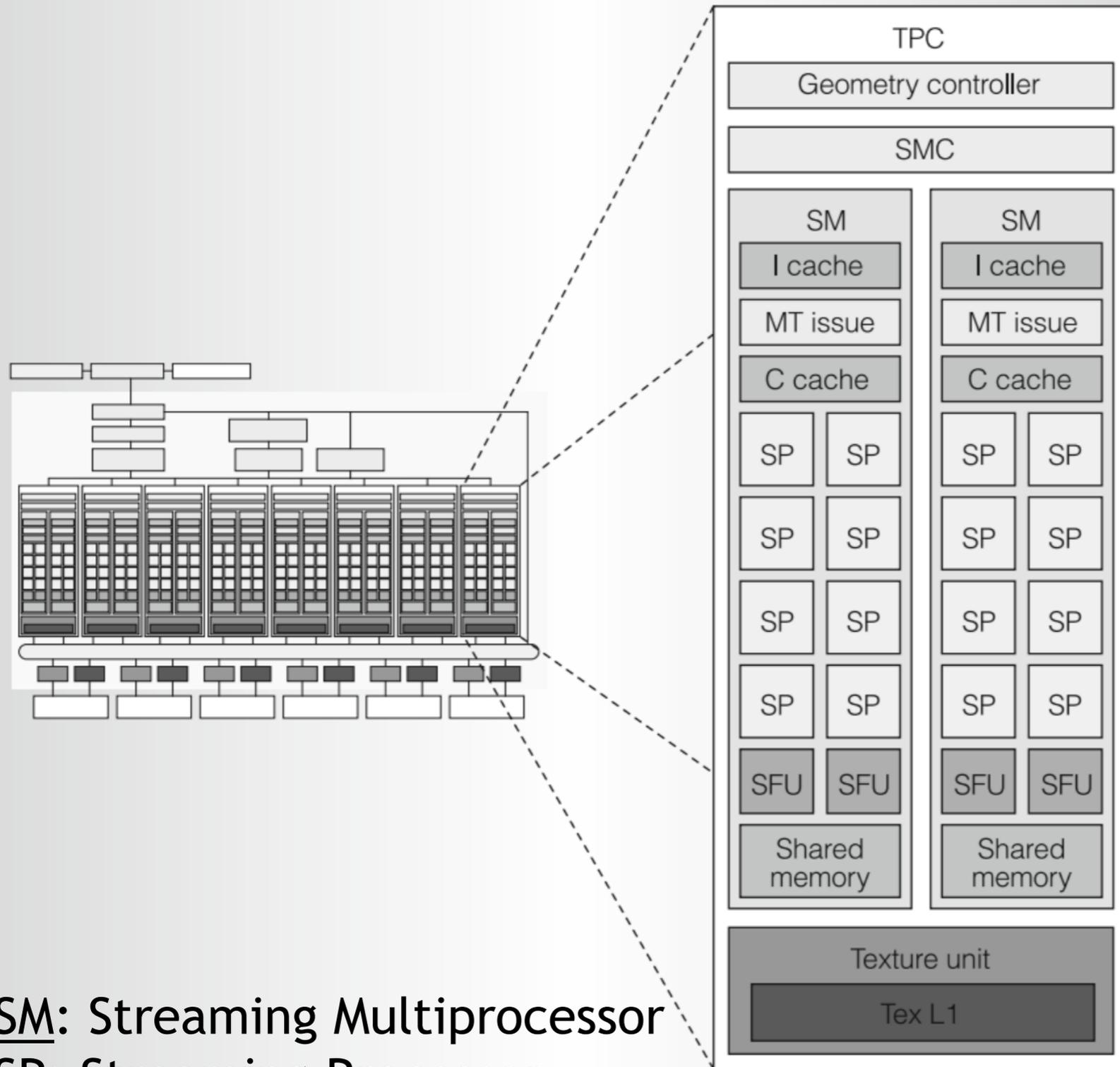


Streaming Processor Array (SPA)

ROP performs colour and depth frame buffer operations directly on memory

No L2 cache coherency problem, data can be in only one cache. Caches are small

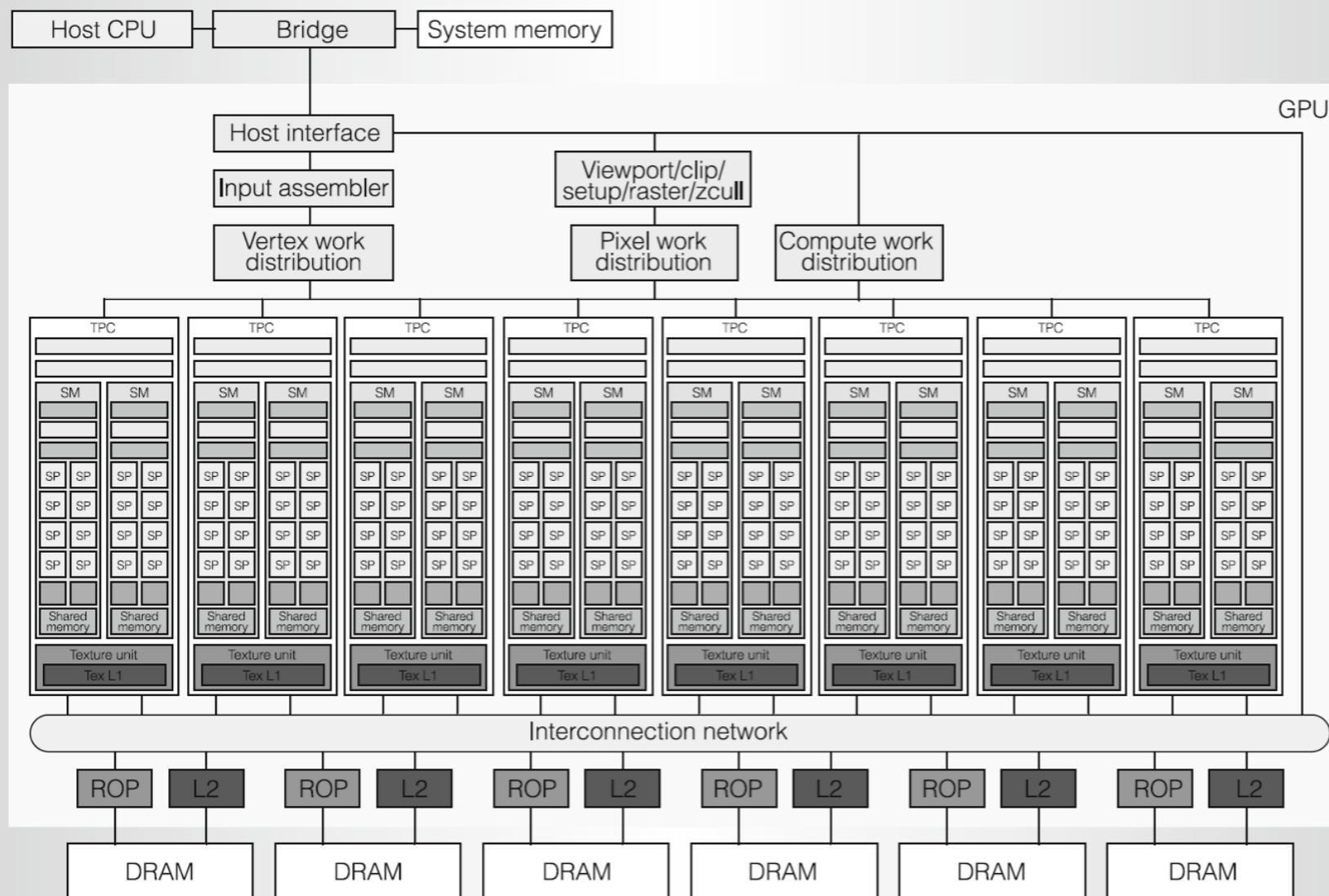
Texture/Processor Cluster (TPC)



- SM: Streaming Multiprocessor
- SP: Streaming Processor

- SMC: Streaming Multiprocessor controller
- MT issue: multithreaded instruction fetch and issue unit
- C cache: constant read-only cache
- I cache: instruction cache
- Geometry controller: directs all primitive and vertex attribute and topology flow in the TPC
- SFU: Special-Function Unit, compute transcendental functions (sin, cos, log x, 1/x)
- Shared memory: scratchpad memory, i.e. user managed cache
- Texture cache does interpolation

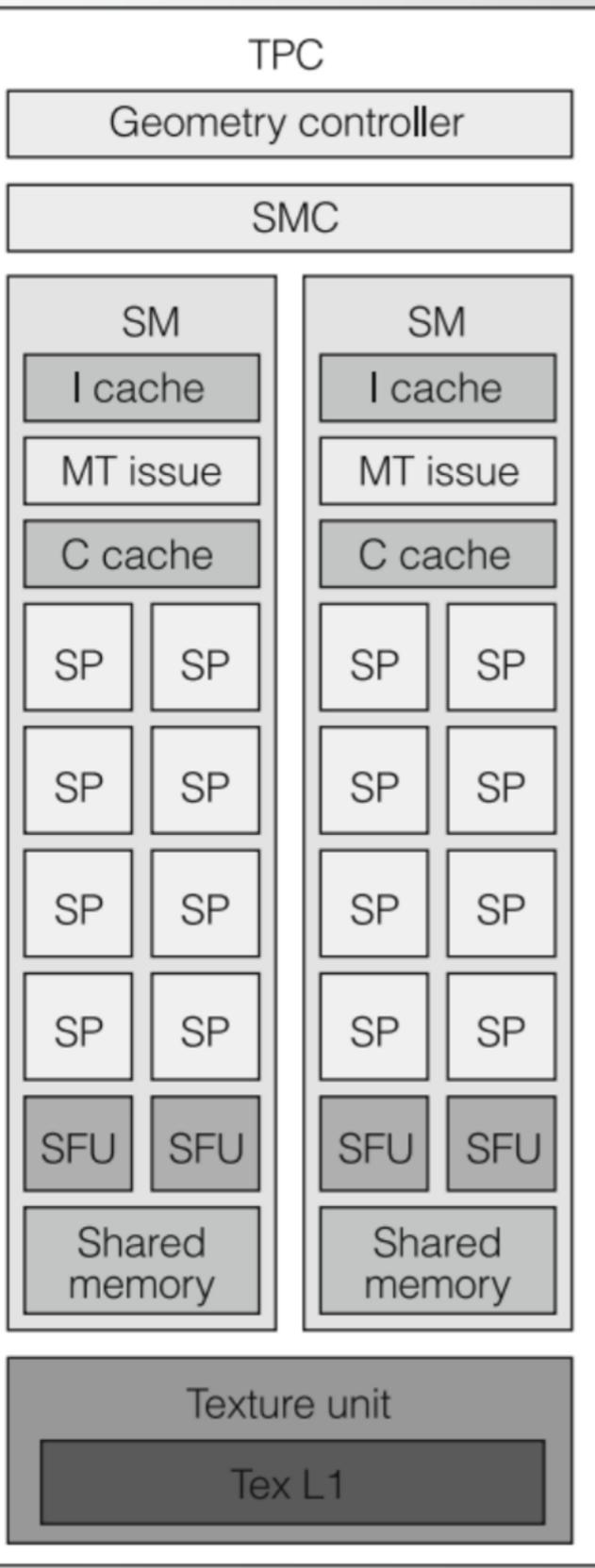
Tesla memory, interconnect, control



- SM's also have an L1 data cache (but no cache-coherency protocol – flushed on kernel launch)
- The chip has multiple DRAM channels, each of which includes an L2 cache
- but each data value can only be in one L2 location, so there's no cache coherency issue at the L2 level
- Tesla has more features specific to graphics, which are not our focus here:
 - Work distribution, load distribution
 - Texture cache, pixel interpolation
 - Z-buffering and alpha-blending (the ROP units, see diagram)

CUDA: using NVIDIA GPUs for general computation

- Designed to do rendering
- Evolved to do general-purpose computing (GPGPU)
 - But to manage thousands of threads, a new programming model is needed, called **CUDA** (Compute Unified Device Architecture)
 - CUDA is proprietary, but the same model lies behind **OpenCL**, an open standard with implementations for multiple vendors' GPUs
- GPU evolved from hardware designed specifically around the OpenGL/DirectX rendering pipeline, with separate vertex- and pixel-shader stages
- “Unified” architecture arose from increased sophistication of shader programs



We focus initially on NVIDIA architecture and terminology. AMD GPUs are quite similar, and the OpenCL programming model is similar to CUDA. Mobile GPUs are somewhat different

CUDA Execution Model

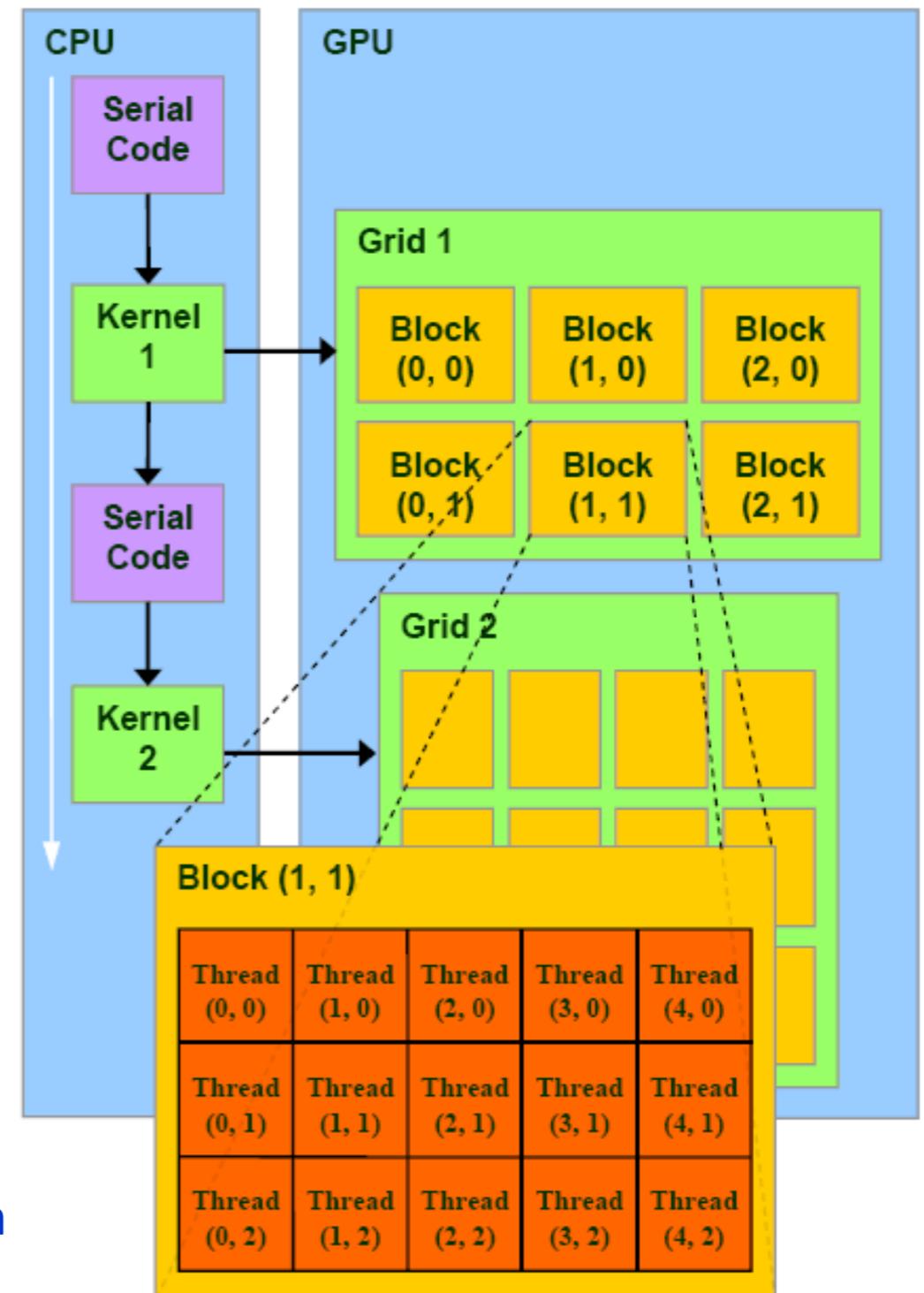
- CUDA is a C extension
 - Serial CPU code
 - Parallel GPU code (*kernels*)
- GPU kernel is a C function
 - Each *thread* executes kernel code
 - A group of threads form a *thread block* (1D, 2D or 3D)
 - Thread blocks are organised into a *grid* (1D, 2D or 3D)
 - Threads within the same thread block can synchronise execution, and share access to local *scratchpad memory*

Key idea: **hierarchy of parallelism**, to handle *thousands* of threads

Thread blocks are allocated (dynamically) to SMs, and run to completion

Threads (warps) within a block **run on the same SM**, so can share data and synchronise

Different blocks in a grid can't interact with each other



Source: CUDA programming guide

```

__global__ void daxpy(int N,
                    double a,
                    double* x,
                    double* y) {
    int i = blockIdx.x *
           blockDim.x +
           threadIdx.x;
    if (i < N)
        y[i] = a*x[i] + y[i];
}

```

CUDA kernel

```

// Invoke DAXPY
daxpy(n, 2.0, x, y);
// DAXPY in C
void daxpy(int n,
           double a,
           double* x,
           double* y) {
    for(int i=0; i < n; ++i)
        y[i] = a*x[i] + y[i];
}

```

C version for comparison

fully parallel loop

CPU code to launch kernel on GPU

```

int main() {
    // Kernel setup
    int N = 1024;
    int blockDim = 256; // These are the threads per block
    int gridDim = N / blockDim; // These are the number of blocks
    daxpy<<<gridDim, blockDim>>>(N, 2.0, x, y); // Kernel invocation
}

```

CUDA example: DAXPY

- ▶ Kernel invocation (“<<<...>>>”) corresponds to enclosing loop nest, managed by hardware
- ▶ Explicitly split into 2-level hierarchy: blocks (256 threads that can share “shared” memory), and grid (N/256 blocks)
- ▶ Kernel commonly consists of just one iteration but could be a loop
- ▶ Multiple tuning parameters trade off register pressure, shared-memory capacity and parallelism

PTX Example (SAXPY code)



```
cvt.u32.u16    $blockid, %ctaid.x;    // Calculate i from thread/block IDs
cvt.u32.u16    $blocksize, %ntid.x;
cvt.u32.u16    $tid, %tid.x;
mad24.lo.u32   $i, $blockid, $blocksize, $tid;
ld.param.u32   $n, [N];              // Nothing to do if n ≤ i
setp.le.u32    $p1, $n, $i;
@$p1 bra      $L_finish;

mul.lo.u32     $offset, $i, 4;        // Load y[i]
ld.param.u32   $yaddr, [Y];
add.u32        $yaddr, $yaddr, $offset;
ld.global.f32  $y_i, [$yaddr+0];
ld.param.u32   $xaddr, [X];          // Load x[i]
add.u32        $xaddr, $xaddr, $offset;
ld.global.f32  $x_i, [$xaddr+0];

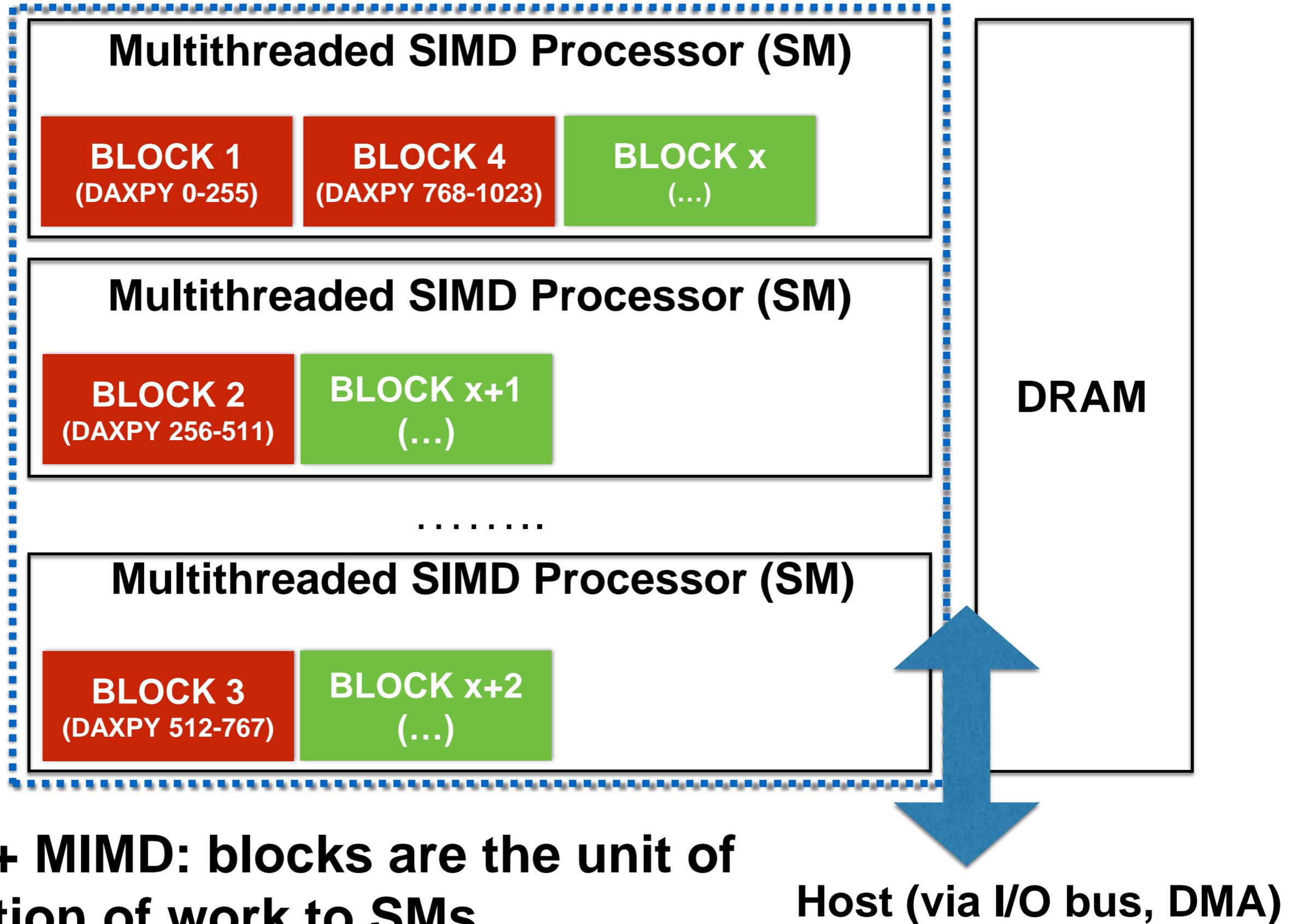
ld.param.f32   $alpha, [ALPHA];      // Compute and store alpha*x[i] + y[i]
mad.f32        $y_i, $alpha, $x_i, $y_i;
st.global.f32  [$yaddr+0], $y_i;

$L_finish:    exit;
```

```
__global__ void daxpy(int N,
                      double a,
                      double* x,
                      double* y) {
    int i = blockIdx.x *
           blockDim.x +
           threadIdx.x;
    if (i < N)
        y[i] = a*x[i] + y[i];
}
```

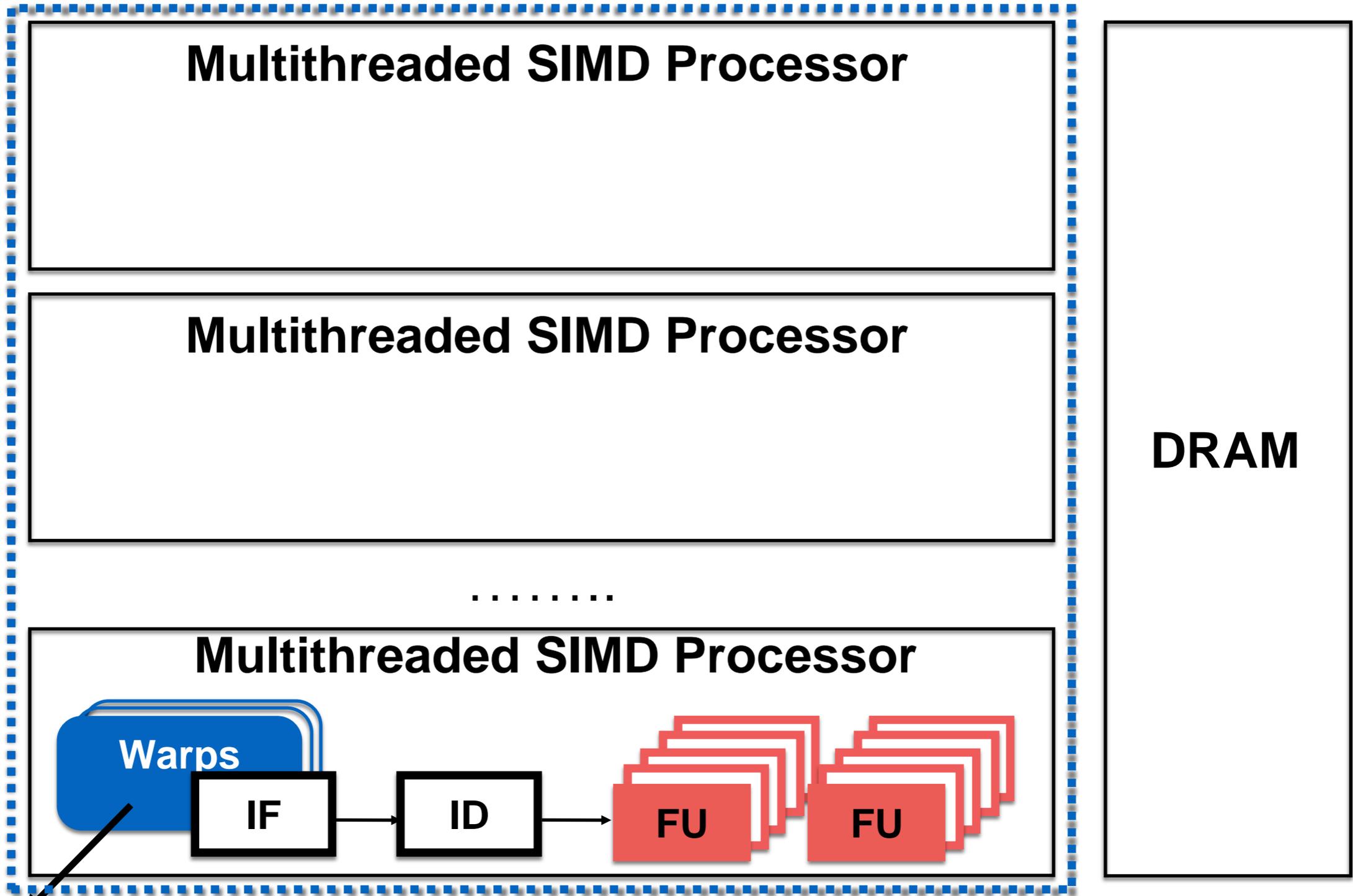
- This is PTX: a pseudo-assembly code that is translated to proprietary ISA
- Threads are scheduled in hardware
- Each thread is provided with its position in the Grid through registers %ctaid, %ntid, %tid
- p1 is a predicate register to determine the outcome of the “if”
- The conditional branch “@\$p1 bra \$L_finish” may be (probably is) translated to predication in the target ISA

Running DAXPY (N=1024) on a GPU



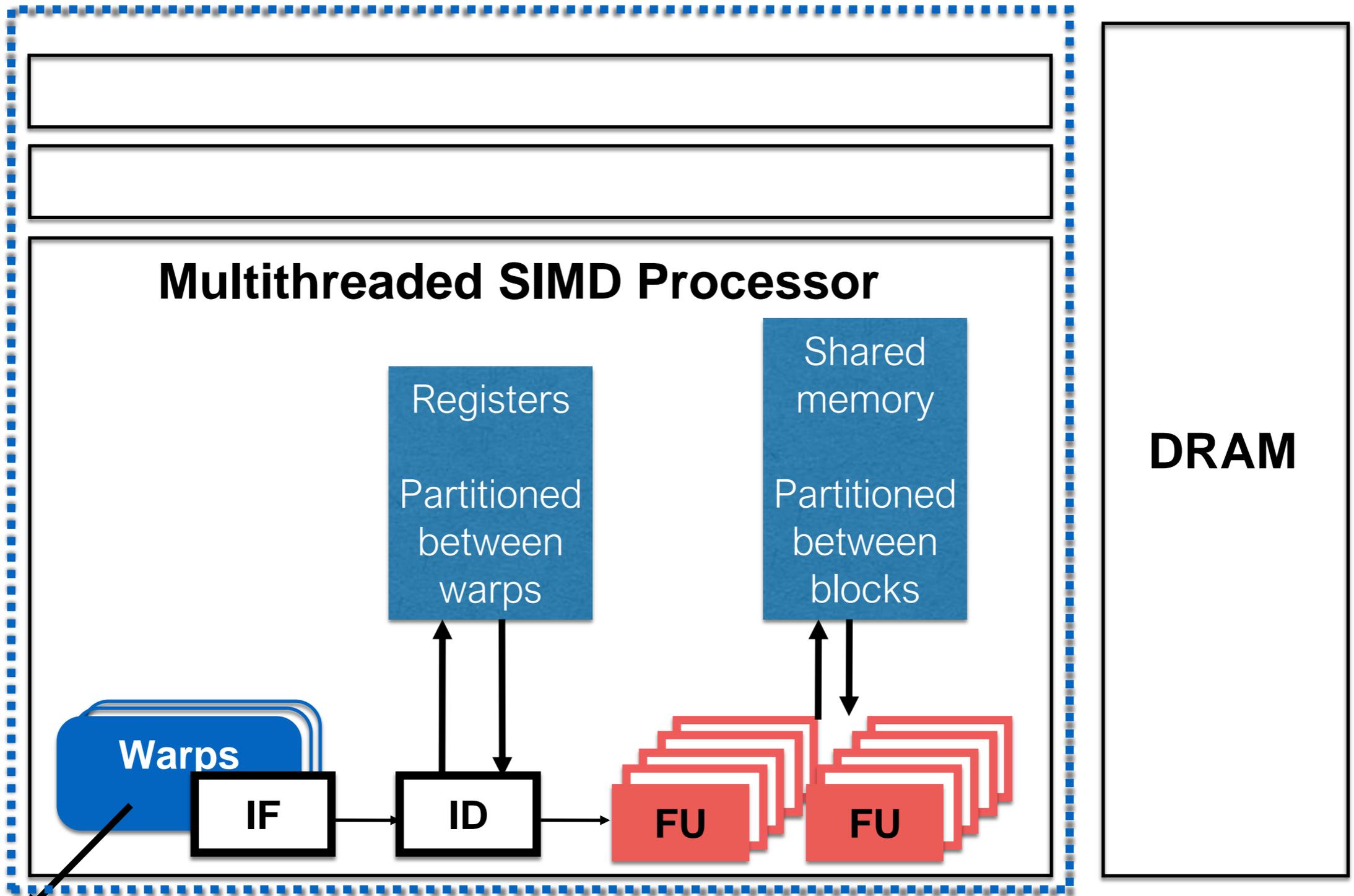
SIMD + MIMD: blocks are the unit of allocation of work to SMs

Running DAXPY on a GPU



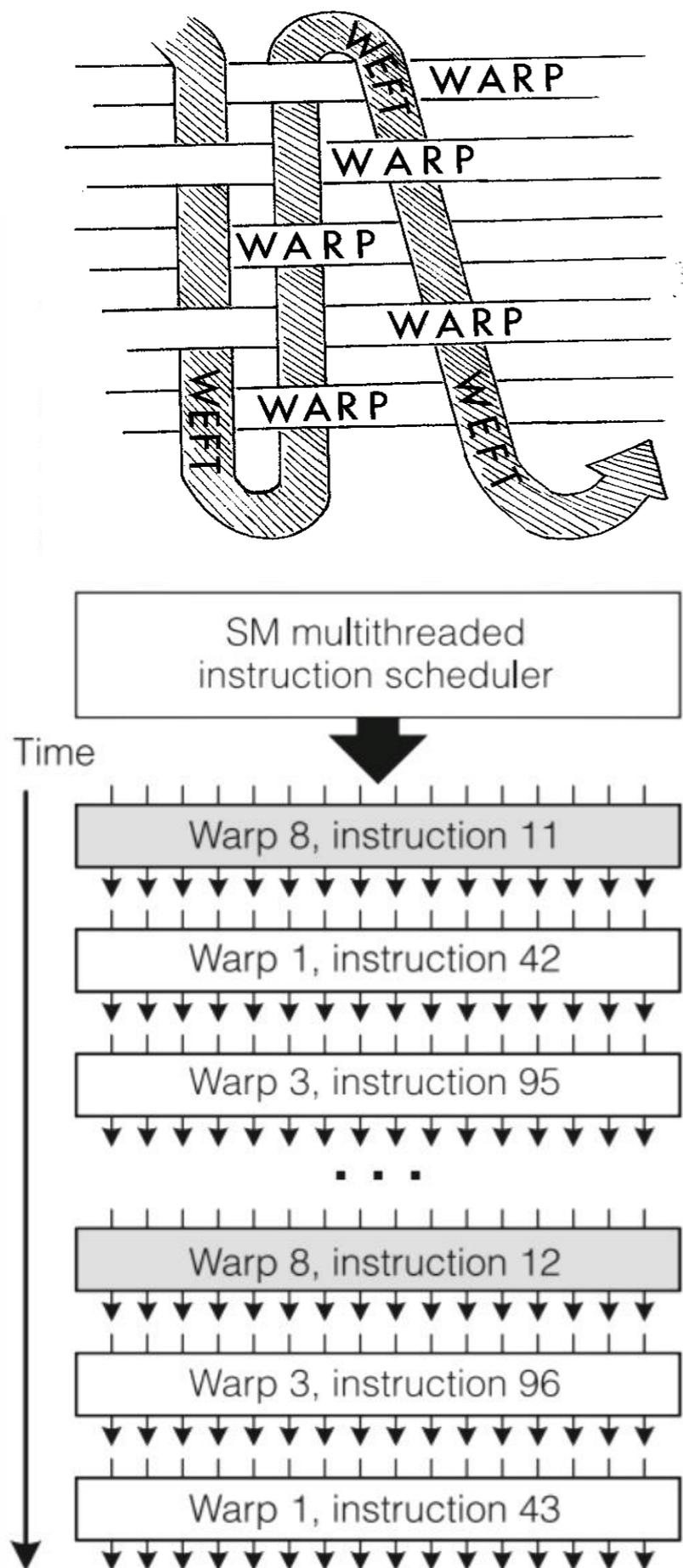
- Each **warp** executes 32 CUDA threads in SIMD lock-step
- Each CUDA thread executes one instance of the kernel
- Each SM is shared by many warps (possibly from the same or different blocks)

Running DAXPY on a GPU



- Each **warp** executes 32 CUDA threads in SIMD lock-step
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- Each SM is shared by many warps (possibly from the same or different blocks)

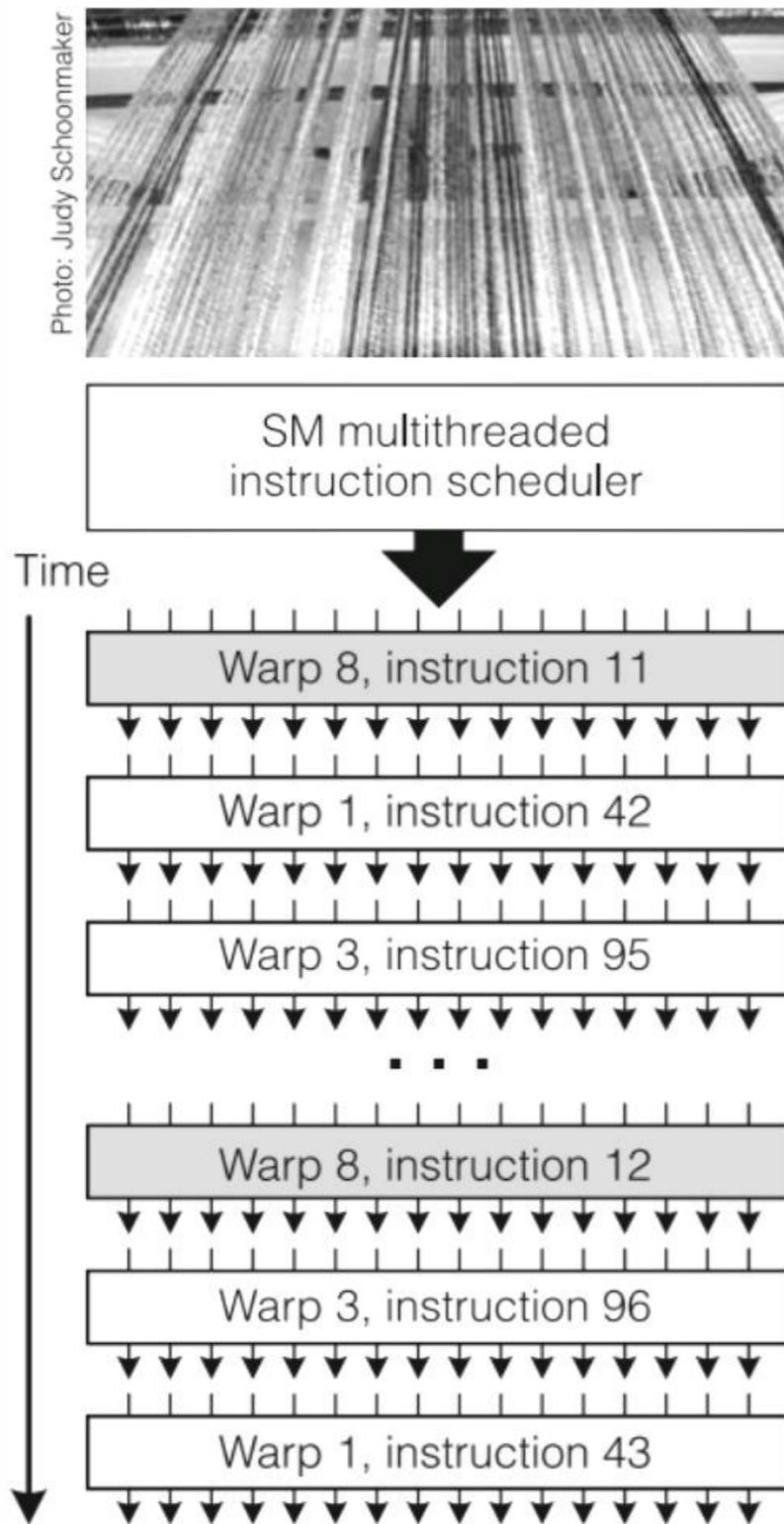
Single-instruction, multiple-thread (SIMT)



- A new parallel programming model: **SIMT**
- The SM's SIMT multithreaded instruction unit creates, manages, schedules, and executes threads in groups of **warps**
- The term warp originates from weaving
- Each SM manages a pool of 24 warps, **24 ways FGMT** (more on later devices)
- Individual threads composing a SIMT warp start together at the same program address, but they are otherwise **free to branch** and execute independently
- At instruction issue time, select **ready-to-run warp** and issue the next instruction to that warp's active threads

Reflecting on SIMT

- SIMT architecture is **similar to SIMD** design, which applies one instruction to multiple data lanes
- The **difference**: SIMT applies one instruction to multiple independent threads in parallel, not just multiple data lanes. A SIMT instruction controls the execution and branching behaviour of one thread
- For **program correctness**, programmers can ignore SIMT executions; but, they can achieve performance improvements if threads in a warp don't diverge
- Correctness/performance analogous to the role of cache lines in traditional architectures
- The SIMT design shares the SM instruction fetch and issue unit efficiently across 32 threads but requires a full warp of active threads for full performance efficiency



Branch divergence

- In a warp, threads all take the same path (good!) or **diverge!**
- A warp serially executes each path, disabling some of the threads
- When all paths complete, the threads reconverge
- **Divergence only occurs within a warp** - different warps execute independently
- **Control-flow coherence**: when all the threads in a warp goes the same way we get good utilisation (a form of locality – spatial branch locality)

```

:
:
if (x == 10)
    c = c + 1;
:

```



Predicate bits: enable/disable each lane

```

:
LDR r5, X
p1 <- r5 eq 10
<p1> LDR r1 <- C
<p1> ADD r1, r1, 1
<p1> STR r1 -> C
:

```

SIMT vs SIMD – GPUs without the hype

- GPUs combine many architectural techniques:
 - Multicore
 - Simultaneous multithreading (SMT)
 - Vector instructions
 - Predication
- So basically a GPU core is a lot like the processor architectures we have studied!
- But the SIMT programming model makes it look different

- ▶ **Overloading the same architectural concept doesn't help GPU beginners**
- ▶ **GPU learning curve is steep in part because of using terms such as “Streaming Multiprocessor” for the SIMD Processor, “Thread Processor” for the SIMD Lane, and “Shared Memory” for Local Memory - especially since Local Memory is not shared between SIMD Processor**

SIMT vs SIMD – GPUs without the hype

SIMT:

- One thread per lane
- Adjacent threads (“warp” / “wavefront”) execute in lockstep
- SMT: multiple “warps” run on the same core, to hide memory latency

SIMD:

- Each thread may include SIMD vector instructions
- SMT: a small number of threads run on the same core to hide memory latency

Which one is easier for the programmer?

SIMT vs SIMD – spatial locality & coalescing

SIMT:

- Spatial locality = adjacent threads access adjacent data
- A load instruction can result in a completely different address being accessed by each lane
- “Coalesced” loads, where accesses are (almost) adjacent, run *much* faster

SIMD:

- Spatial locality = adjacent loop iterations access adjacent data
- A SIMD vector load usually *has* to access adjacent locations
- Some recent processors have “gather” instructions which can fetch from a different address per lane
- But performance is often serialised

SIMT vs SIMD – spatial locality & coalescing

SIMD (on CPU):

```
void add (float *c, float *a, float *b)
{
    for (int i=0; i <= N; i++)
        #pragma omp simd
        for (int j=0; j <= N; j++)
            c[i][j]=a[i][j]+b[i][j];
}
```

Using OpenMP

This example has good spatial locality because it traverses the data in layout order:

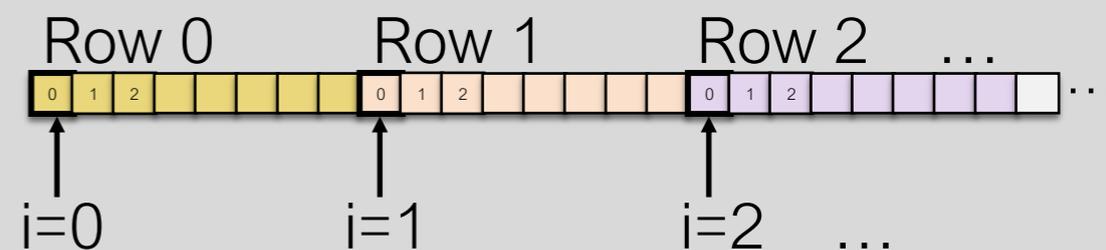
```
void add (float *c, float *a, float *b)
{
    for (int i=0; i <= N; i++) {
        __m128* pa = (__m128*) &a[i][0];
        __m128* pb = (__m128*) &b[i][0];
        __m128* pc = (__m128*) &c[i][0];
        for (int i=0; i <= N/4; i++)
            *pc++ = _mm_add_ps(*pa++, *pb++);
    }
}
```

Using intrinsics

SIMT (on GPU):

```
__global__ void add(int N,
                    double* a,
                    double* b,
                    double* c) {
    int i = blockIdx.x *
            blockDim.x +
            threadIdx.x;
    for (int j=0; j <= N; j++)
        c[i][j] = a[i][j] + b[i][j];
}
```

This example has terrible spatial locality because adjacent threads access different columns



Threads with adjacent thread ids access data in different cache lines

SIMT vs SIMD – spatial *control* locality

SIMT:

- Branch coherence = adjacent threads in a warp all usually branch the same way (*spatial* locality for branches, across threads)

SIMD:

- Branch predictability = each *individual* branch is mostly taken or not-taken (or is well-predicted by global history)

NVIDIA Volta GPU (2017)



GV100 with 84 SMs



GV100's SM includes 8 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

Tensor core computes matrix-matrix multiply on FP16s with FP32 accumulation

Pre-Volta



Volta



32 thread warp with independent scheduling

Each CUDA thread has its own PC and stack, enabling dynamic scheduling in hardware to heuristically enhance branch convergence

It is a heterogeneous world



TITAN
4998 GFLOPS
< 400 W



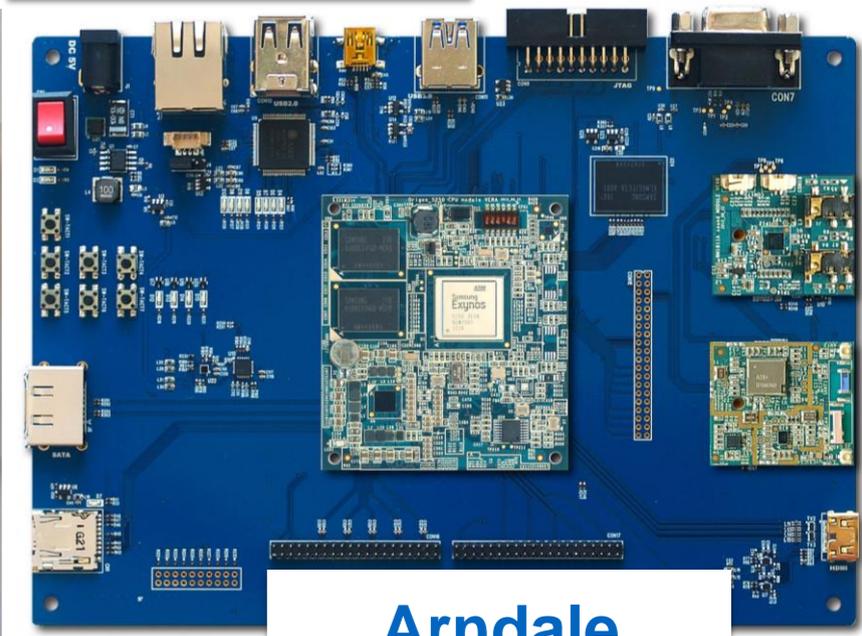
GTX 870M
2827 GFLOPS
< 100 W



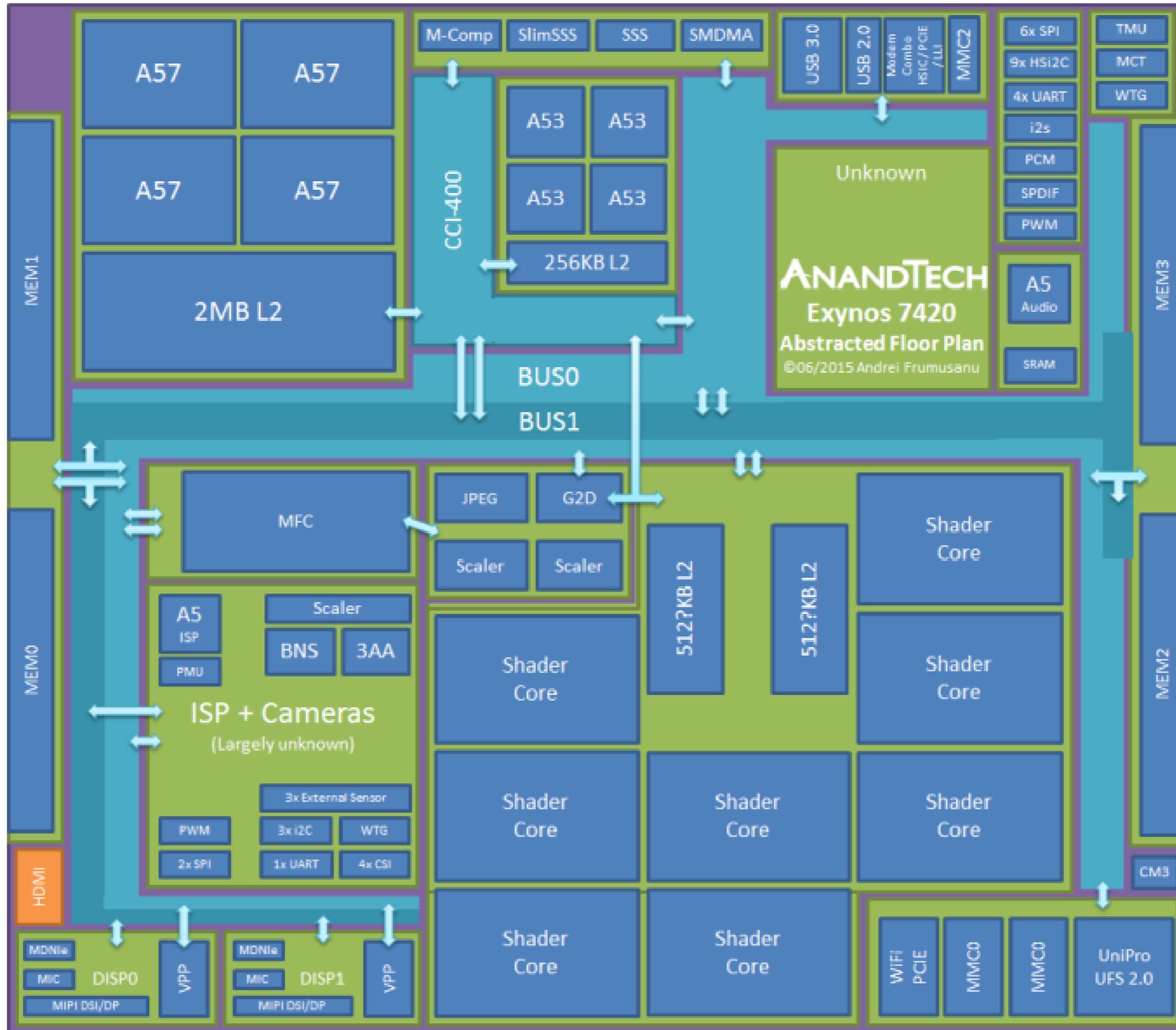
TK1
404 GFLOPS
< 20 W



ODROID
170 GFLOPS
< 10 W



Arndale
87 GFLOPS
< 5 W

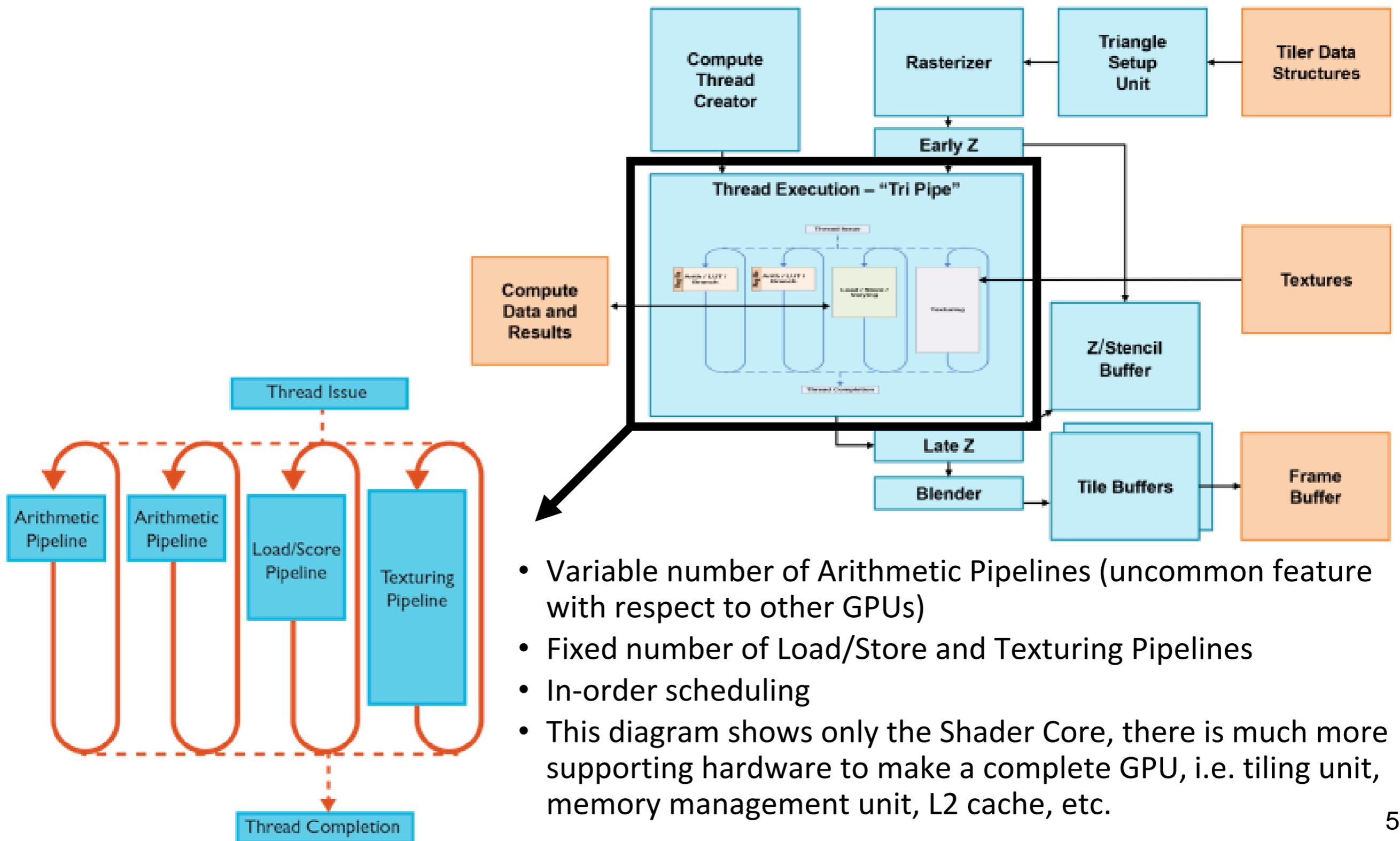


ARM-based Samsung Exynos 7420 SoC
Reverse engineered

spare slides for interest

ARM MALI GPU: Midgard microarchitecture

Shader Core Architecture



- Variable number of Arithmetic Pipelines (uncommon feature with respect to other GPUs)
- Fixed number of Load/Store and Texturing Pipelines
- In-order scheduling
- This diagram shows only the Shader Core, there is much more supporting hardware to make a complete GPU, i.e. tiling unit, memory management unit, L2 cache, etc.

Midgard arithmetic Pipe

ARM Mali Midgard Arithmetic Pipe

V_MUL



V_ADD



V_SFU



- Very flexible SIMD
- Simply fill the SIMD with as many (identical) operations as will fit, and the SIMD will handle it

- ARM Midgard is a VLIW design with SIMD characteristics (power efficient)
- So, at a high level ARM is feeding multiple ALUs, including SIMD units, with a single long word of instructions (ILP)
- Support a wide range of data types, integer and FP: I8, I16, I32, I64, FP16, FP32, FP64
- 17 SP GFLOPS per core at 500 MHz (if you count also the SFUs)

Optimising for MALI GPUs

How to run optimally OpenCL code on Mali GPUs means mainly to locate and remove optimisations for alternative compute devices:

- Use of local or private memory: Mali GPUs use caches instead of local memories. There is therefore no performance advantage using these memories on a Mali
- Barriers: data transfers to or from local or private memories are typically synchronised with barriers. If you remove copy operations to or from these memories, also remove the associated barriers
- Use of scalars: some GPUs work with scalars whereas Mali GPUs can also use vectors. Do vectorise your code
- Optimisations for divergent threads: threads on a Mali are independent and can diverge without any performance impact. If your code contains optimisations for divergent threads in warps, remove them
- Modifications for memory bank conflicts: some GPUs include per-warp memory banks. If the code includes optimisations to avoid conflicts in these memory banks, remove them
- No host-device copies: Mali shares the same memory with the CPU

Source: http://infocenter.arm.com/help/topic/com.arm.doc.dui0538f/DUI0538F_mali_t600_opencldg.pdf

Texture cache

- GPUs were built for rendering
- Critical element:
 - Mapping from a stored texture onto a triangular mesh
- To render each triangle:
 - enumerate the pixels,
 - map each pixel to the texture and interpolate
- Texture cache
 - Can be accessed with 2d float index
 - Cache includes dedicated hardware to implement bilinear interpolation
 - Can be configured to clamp, border, wrap or mirror at texture boundary
 - Hardware support to decompress compressed textures on cache miss
 - Custom hardware-specific storage layout (blocked/Morton) to exploit 2d locality
 - Triangle/pixel enumeration is tiled for locality

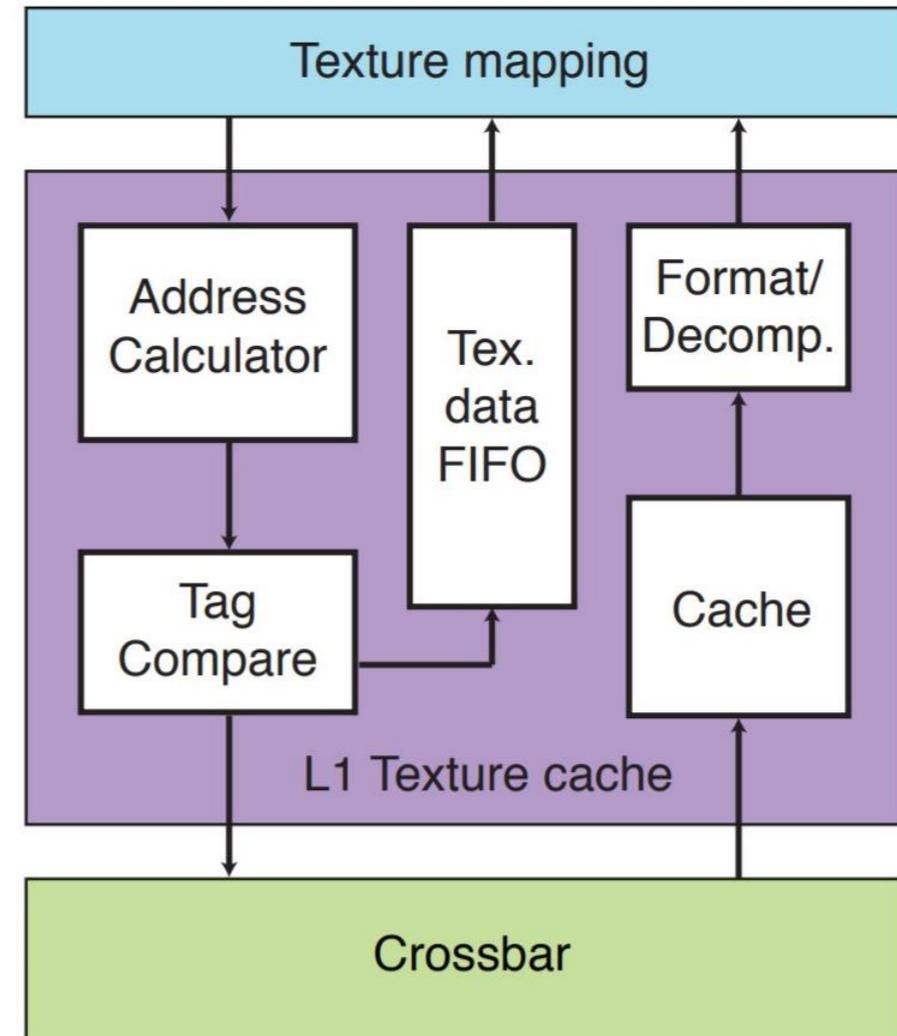
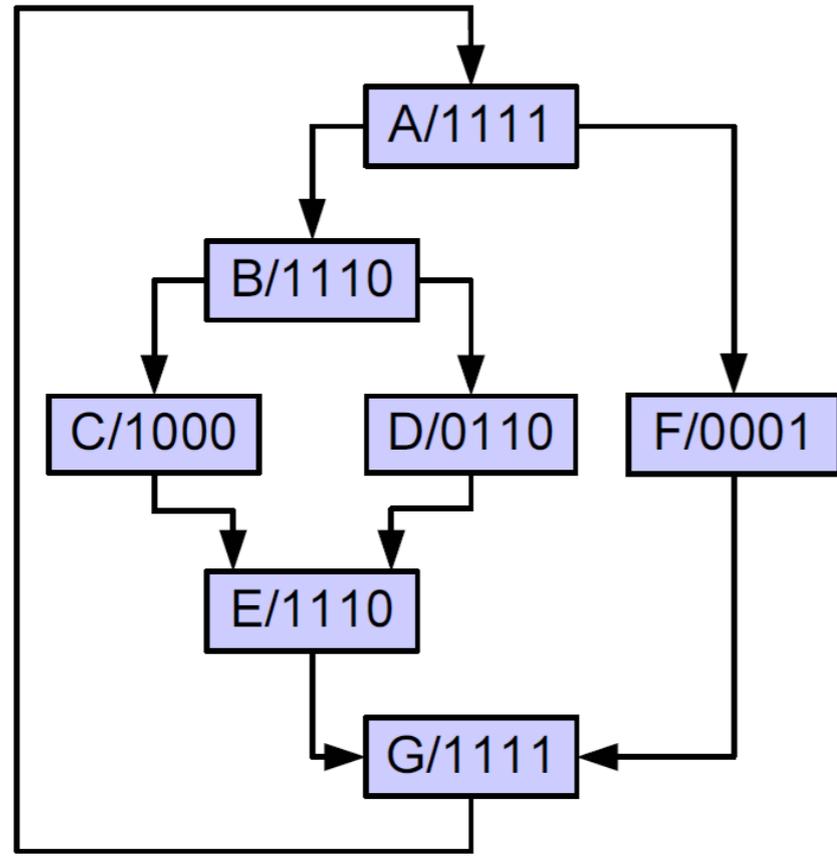


Fig. 5. An overview of a texture cache architecture. The texture mapping unit provides texture coordinates for which a memory address is calculated. The address is sent to the tag compare to determine if the data is in the cache. If the data isn't in the cache, a request is sent via the crossbar to the L2 cache. Any state associated with the original request is sent into a FIFO to return to the texture mapping unit with the texel data. Once the data arrives in the cache, or is already available in the cache, it is returned to the texture mapping unit. If the data is compressed, it is decompressed and any formatting that is required is done.

Nested if-then-else execution



(a) Example Program

	Ret./Reconv. PC	Next PC	Active Mask
	-	G	1111
	G	F	0001
TOS →	G	B	1110

(c) Initial State

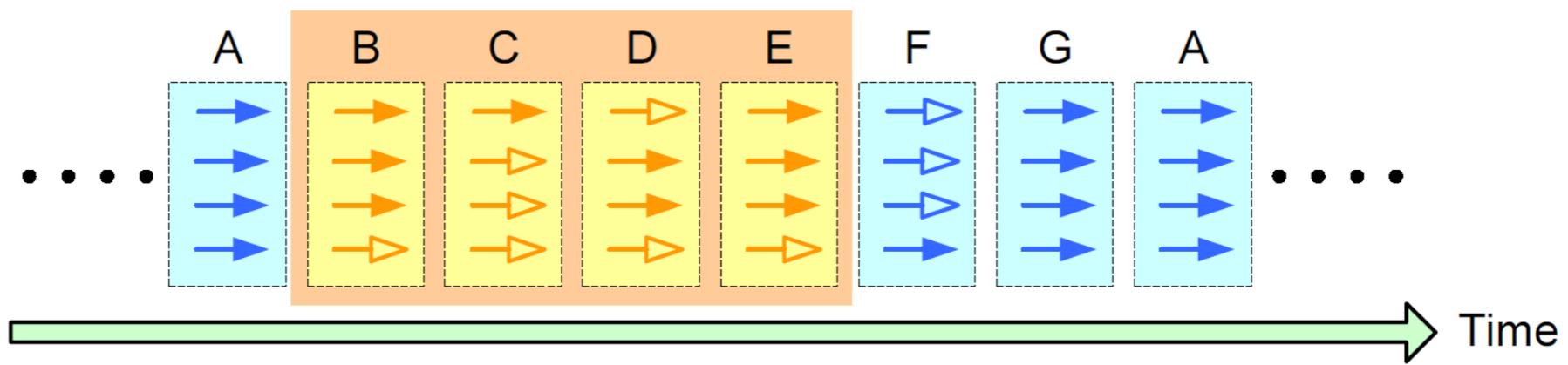
	Ret./Reconv. PC	Next PC	Active Mask
	-	G	1111
	G	F	0001
	G	E	1110
	E	D	0110
TOS →	E	C	1000

(i)
(ii)
(iii)

(d) After Divergent Branch

	Ret./Reconv. PC	Next PC	Active Mask
	-	G	1111
	G	F	0001
TOS →	G	E	1110

(e) After Reconvergence



(b) Re-convergence at Immediate Post-Dominator of B