CS 380 - GPU and GPGPU Programming
Lecture 4: GPU Architecture 3

Markus Hadwiger, KAUST

Reading Assignment #2 (until Feb. 9)

Read (required):

- GLSL book, chapter 4 (*The OpenGL Programmable Pipeline*)
- GPU Gems 2 book, chapter 30 (*The GeForce 6 Series GPU Architecture*) available online:

`http://download.nvidia.com/developer/GPU_Gems_2/GPU_Gems2_ch30.pdf`

From Shader Code to a Teraflop: How Shader Cores Work

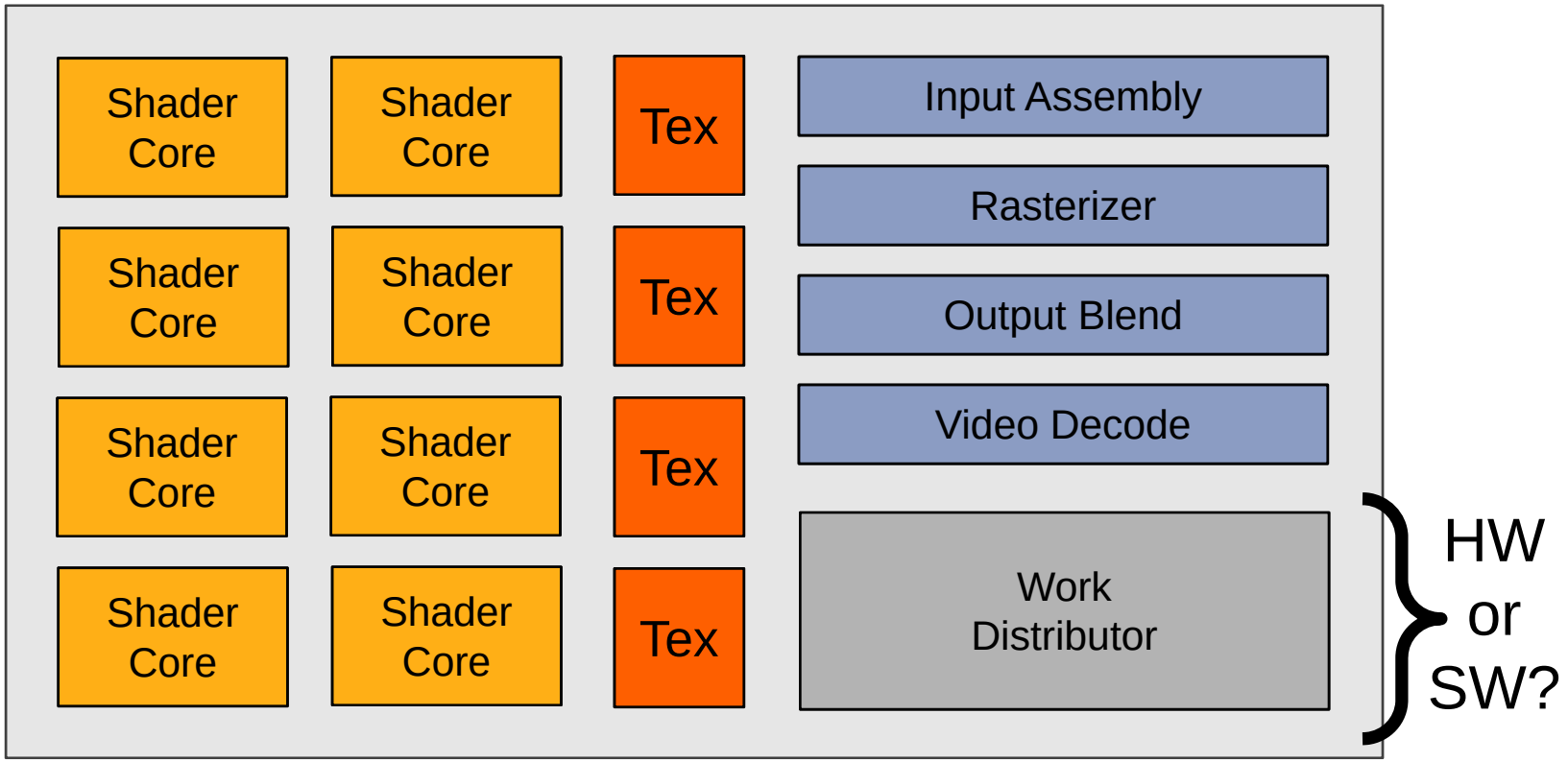
Kayvon Fatahalian
Stanford University

Part 1: throughput processing

- Three key concepts behind how modern GPU processing cores run code

- Knowing these concepts will help you:
  1. Understand space of GPU core (and throughput CPU processing core) designs
  2. Optimize shaders/compute kernels
  3. Establish intuition: what workloads might benefit from the design of these architectures?

What's in a GPU?



Heterogeneous chip multi-processor (highly tuned for graphics)

A diffuse reflectance shader

```
sampler mySamp;
Texture2D<float3> myTex;
float3 lightDir;

float4 diffuseShader(float3 norm, float2 uv)
{
  float3 kd;
  kd = myTex.Sample(mySamp, uv);
  kd *= clamp( dot(lightDir, norm), 0.0, 1.0);
  return float4(kd, 1.0);
}
```

Independent, but no explicit parallelism

Compile shader

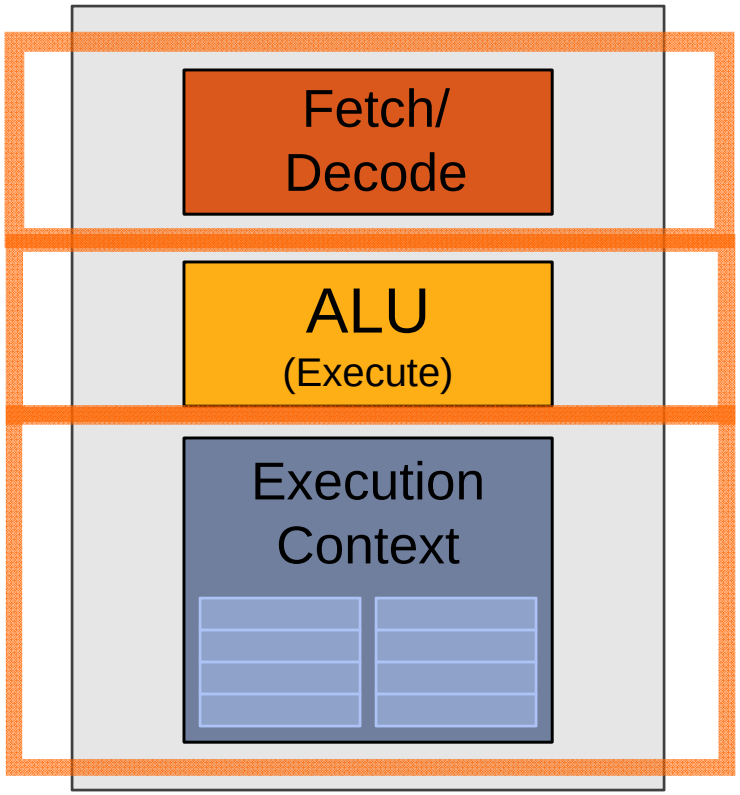
1 unshaded fragment input record

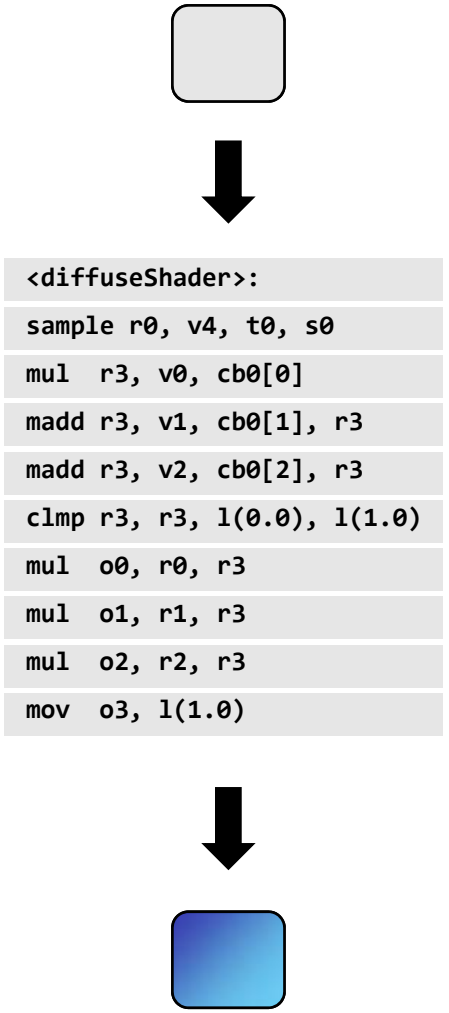
```
sampler mySamp;
Texture2D<float3> myTex;
float3 lightDir;

float4 diffuseShader(float3 norm, float2 uv)
{
  float3 kd;
  kd = myTex.Sample(mySamp, uv);
  kd *= clamp ( dot(lightDir, norm), 0.0, 1.0);
  return float4(kd, 1.0);
}
```

```
<diffuseShader>:
sample r0, v4, t0, s0
mul  r3, v0, cb0[0]
madd r3, v1, cb0[1], r3
madd r3, v2, cb0[2], r3
clmp r3, r3, l(0.0), l(1.0)
mul  o0, r0, r3
mul  o1, r1, r3
mul  o2, r2, r3
mov  o3, l(1.0)
```

1 shaded fragment output record

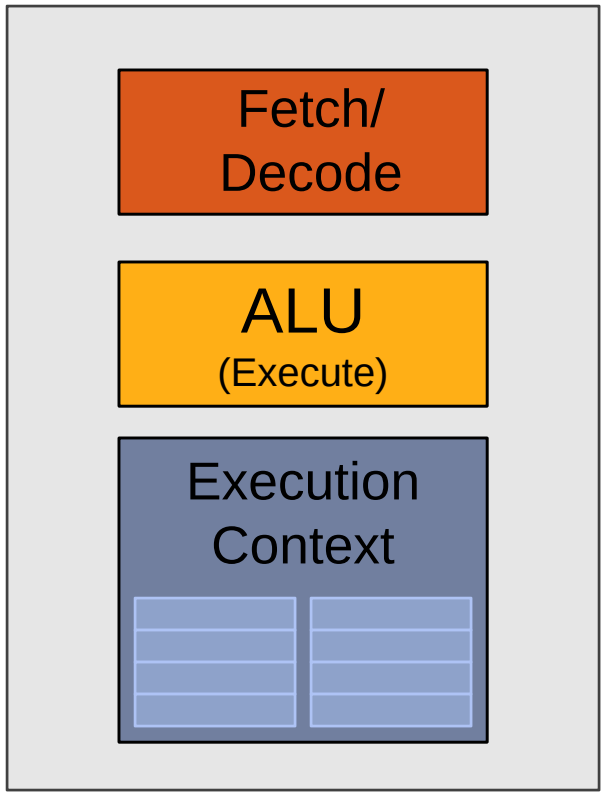
Execute shader

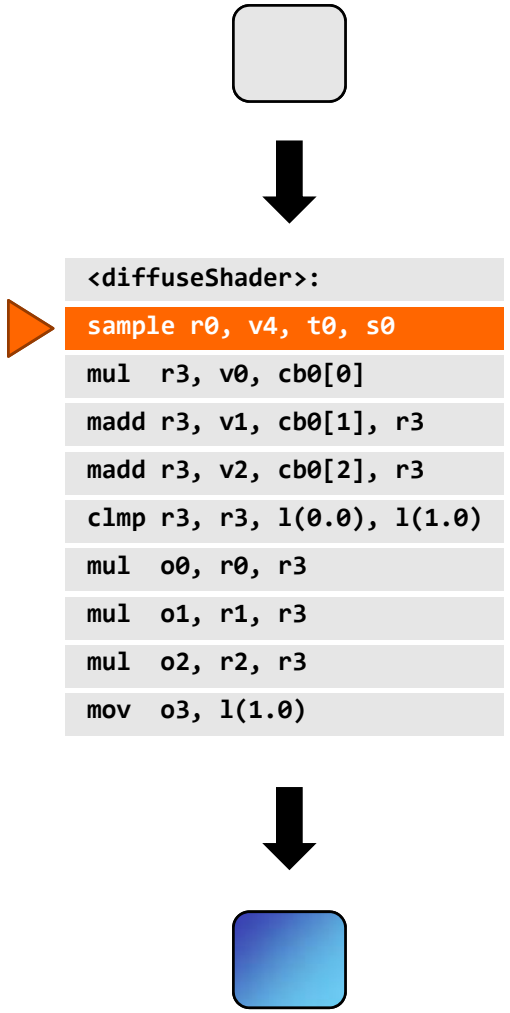




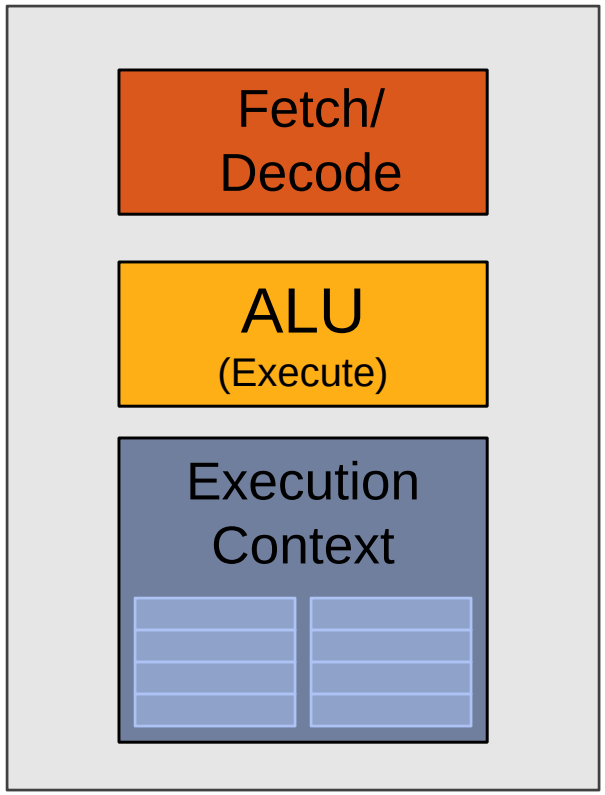
```
<diffuseShader>:
sample r0, v4, t0, s0
mul  r3, v0, cb0[0]
madd r3, v1, cb0[1], r3
madd r3, v2, cb0[2], r3
clmp r3, r3, l(0.0), l(1.0)
mul  o0, r0, r3
mul  o1, r1, r3
mul  o2, r2, r3
mov  o3, l(1.0)
```

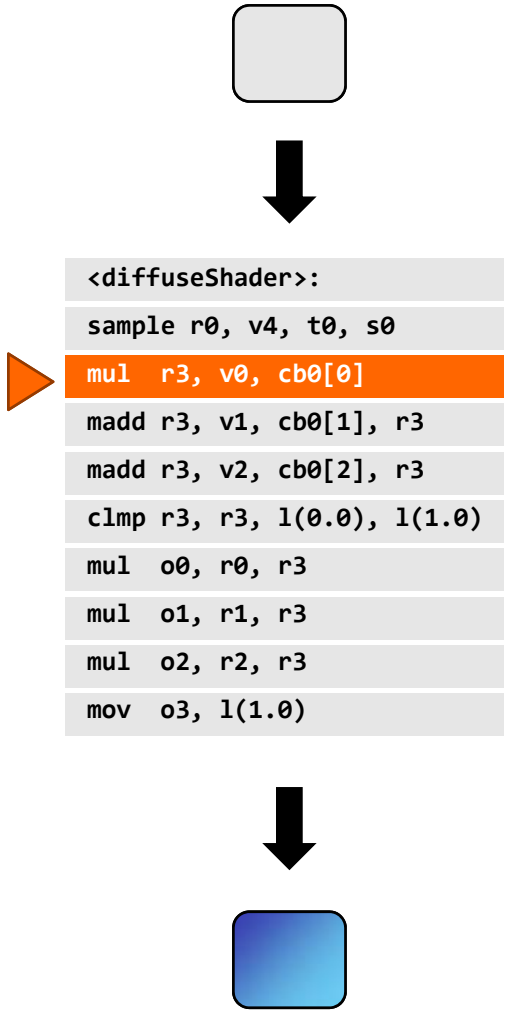
Execute shader



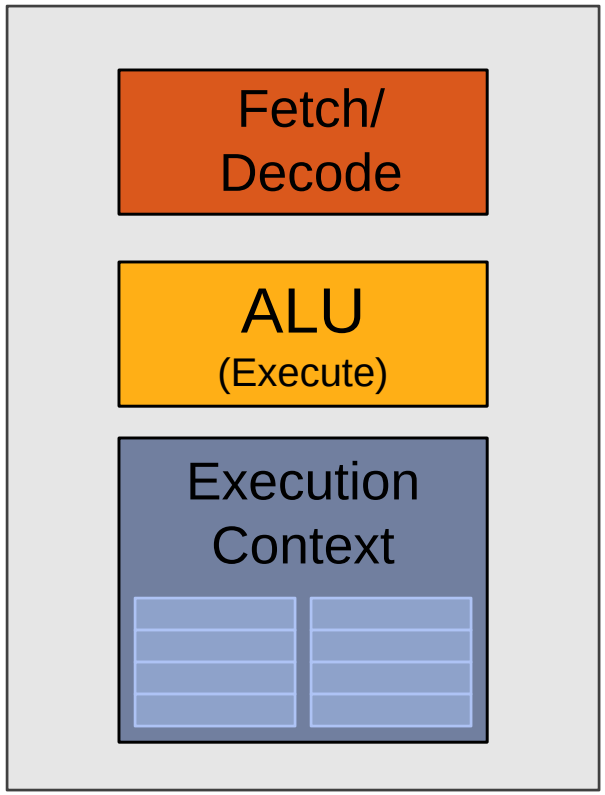


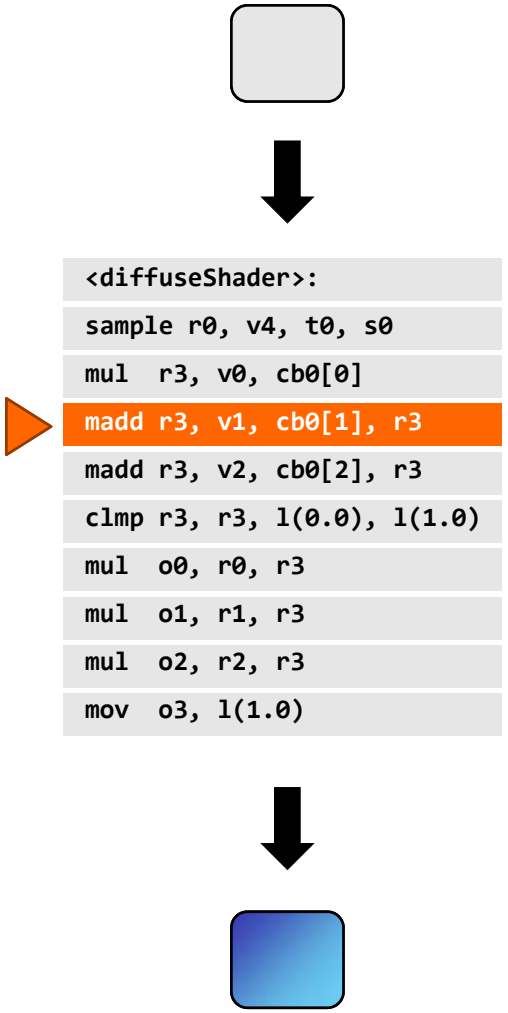
Execute shader



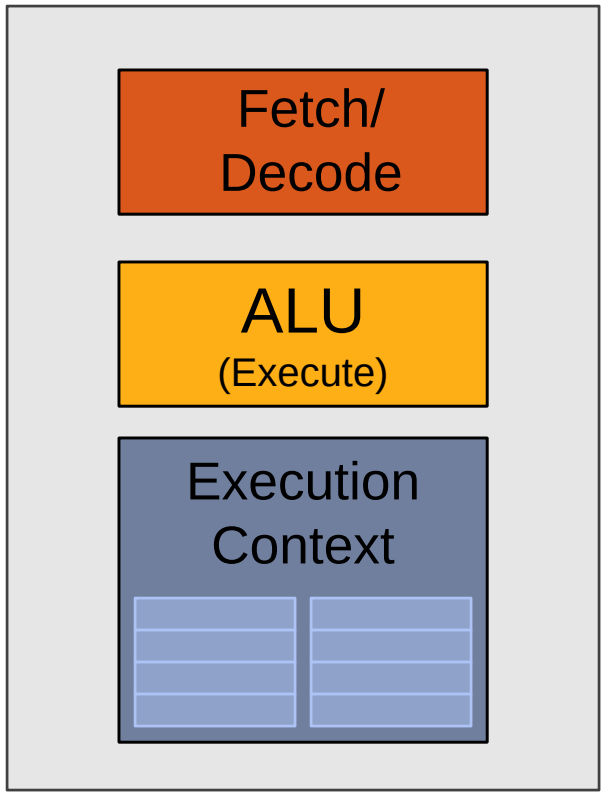


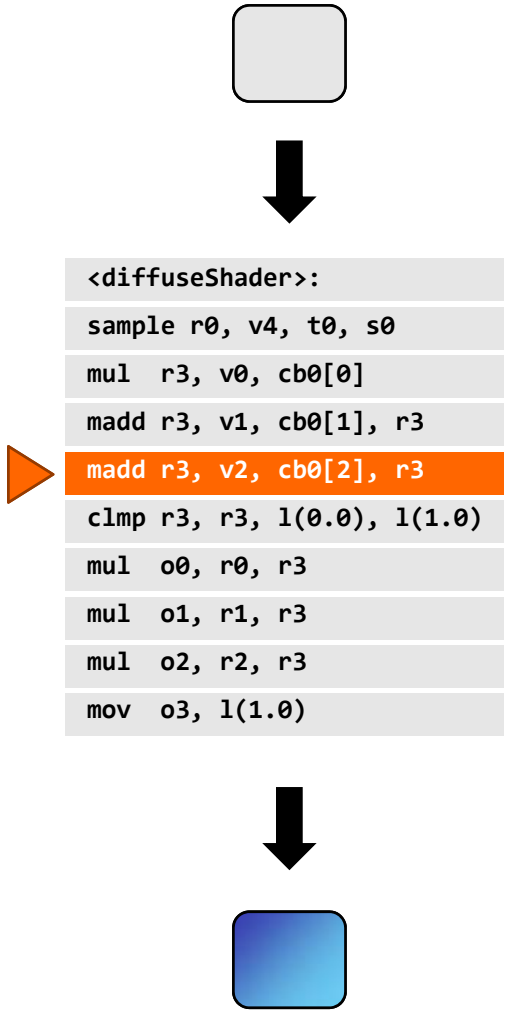
Execute shader



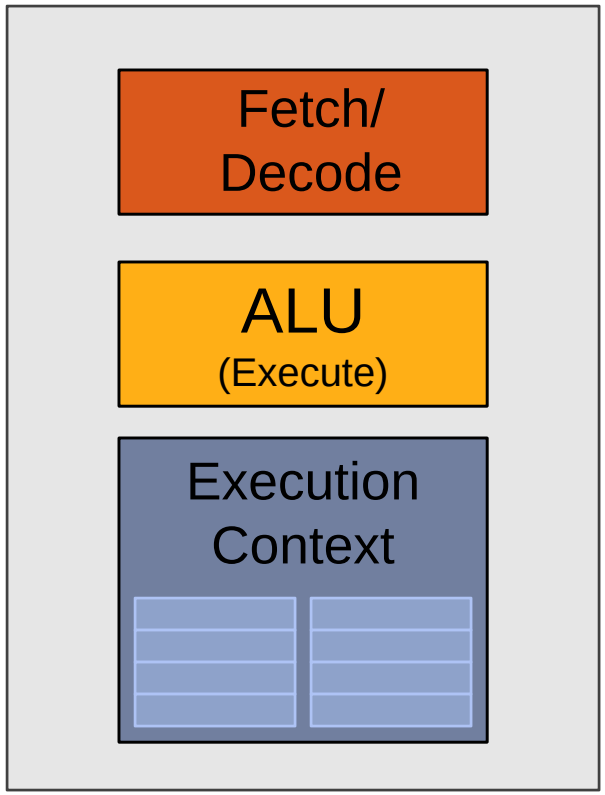


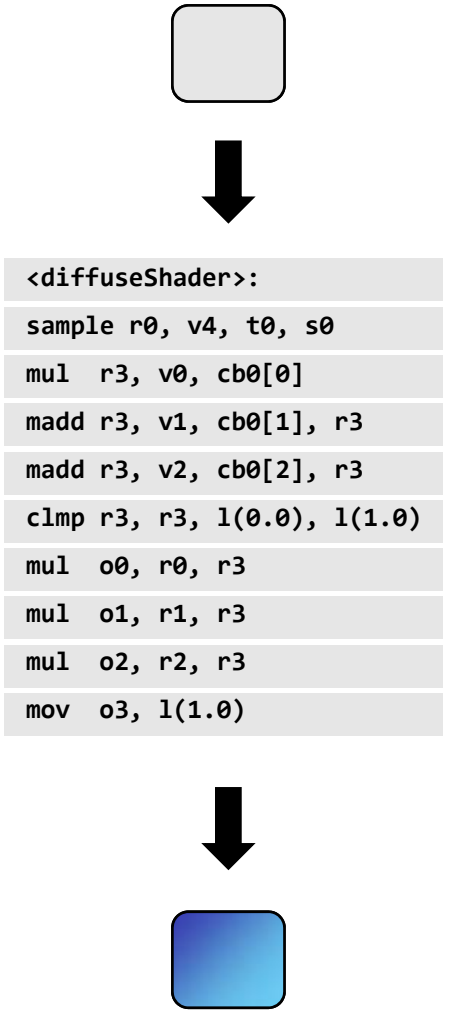
Execute shader





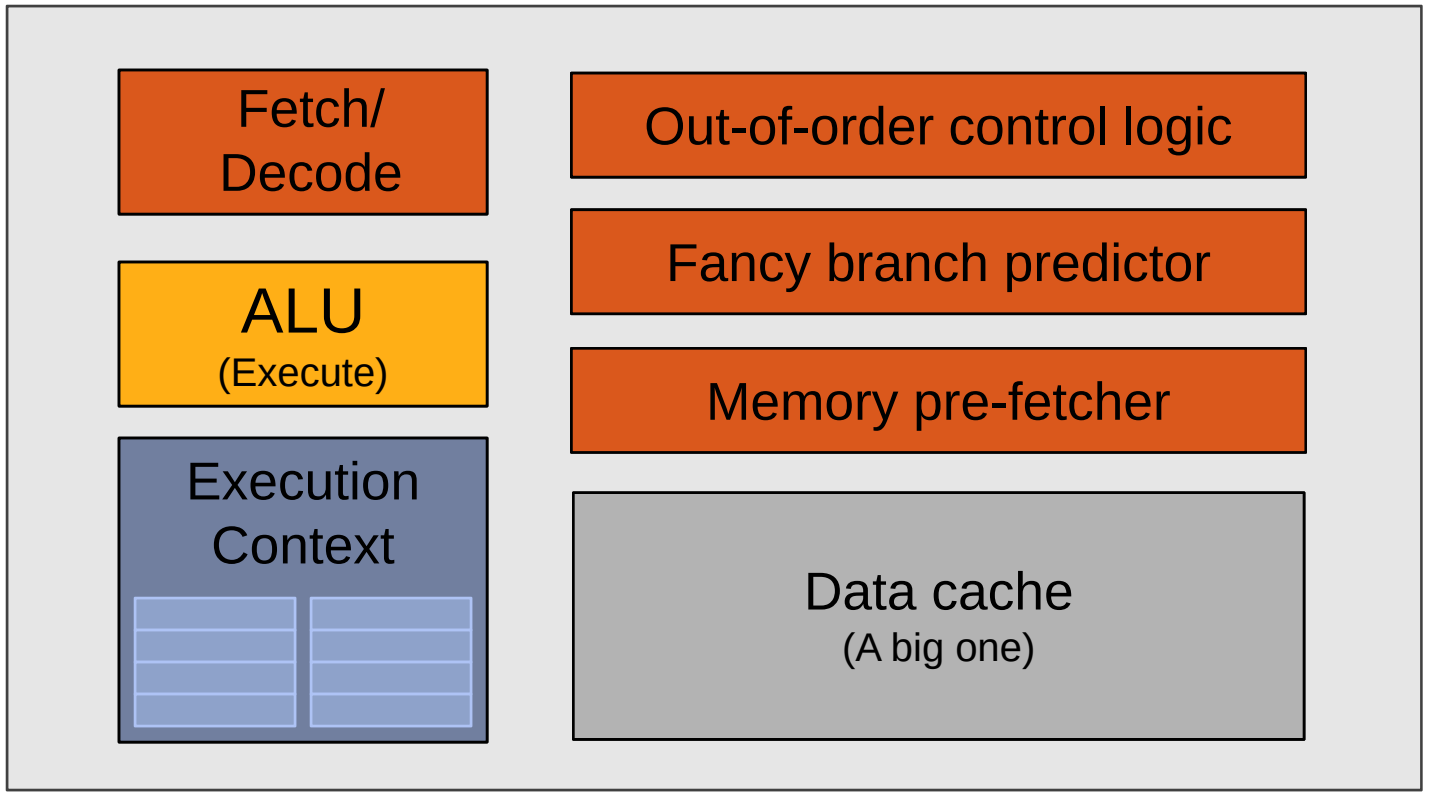
Execute shader



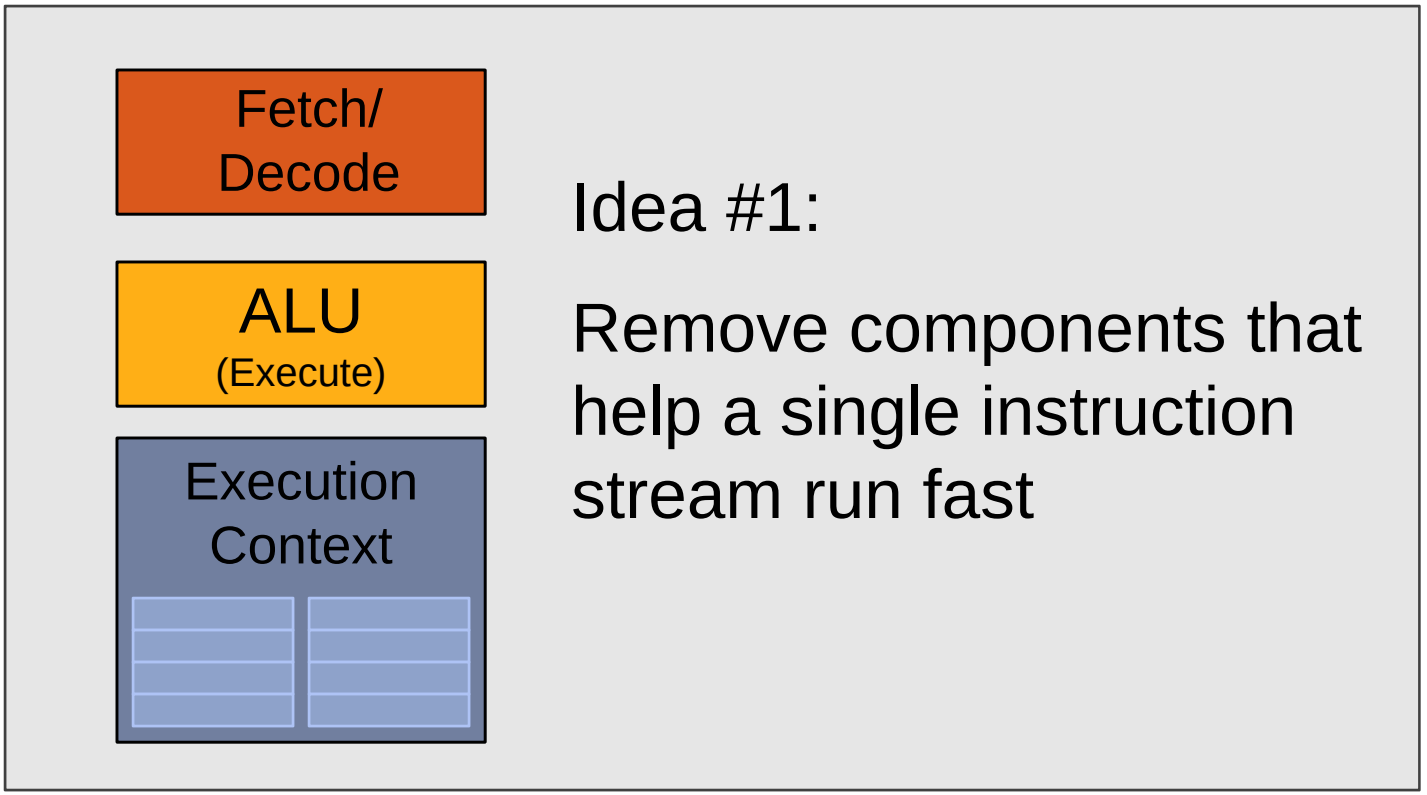


```
<diffuseShader>:
sample r0, v4, t0, s0
mul  r3, v0, cb0[0]
madd r3, v1, cb0[1], r3
madd r3, v2, cb0[2], r3
clmp r3, r3, l(0.0), l(1.0)
mul  o0, r0, r3
mul  o1, r1, r3
mul  o2, r2, r3
mov  o3, l(1.0)
```

CPU-“style” cores



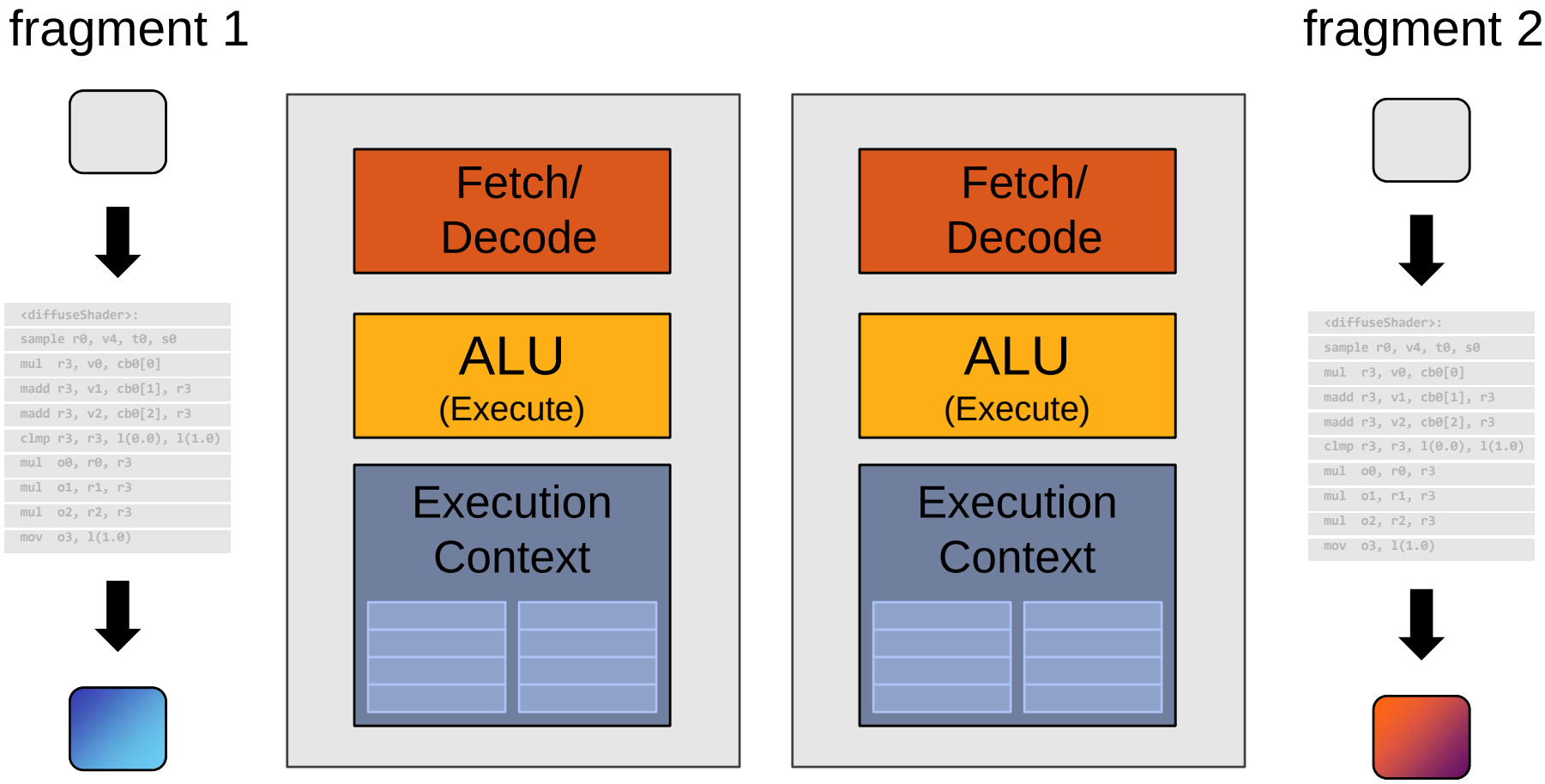
Slimming down



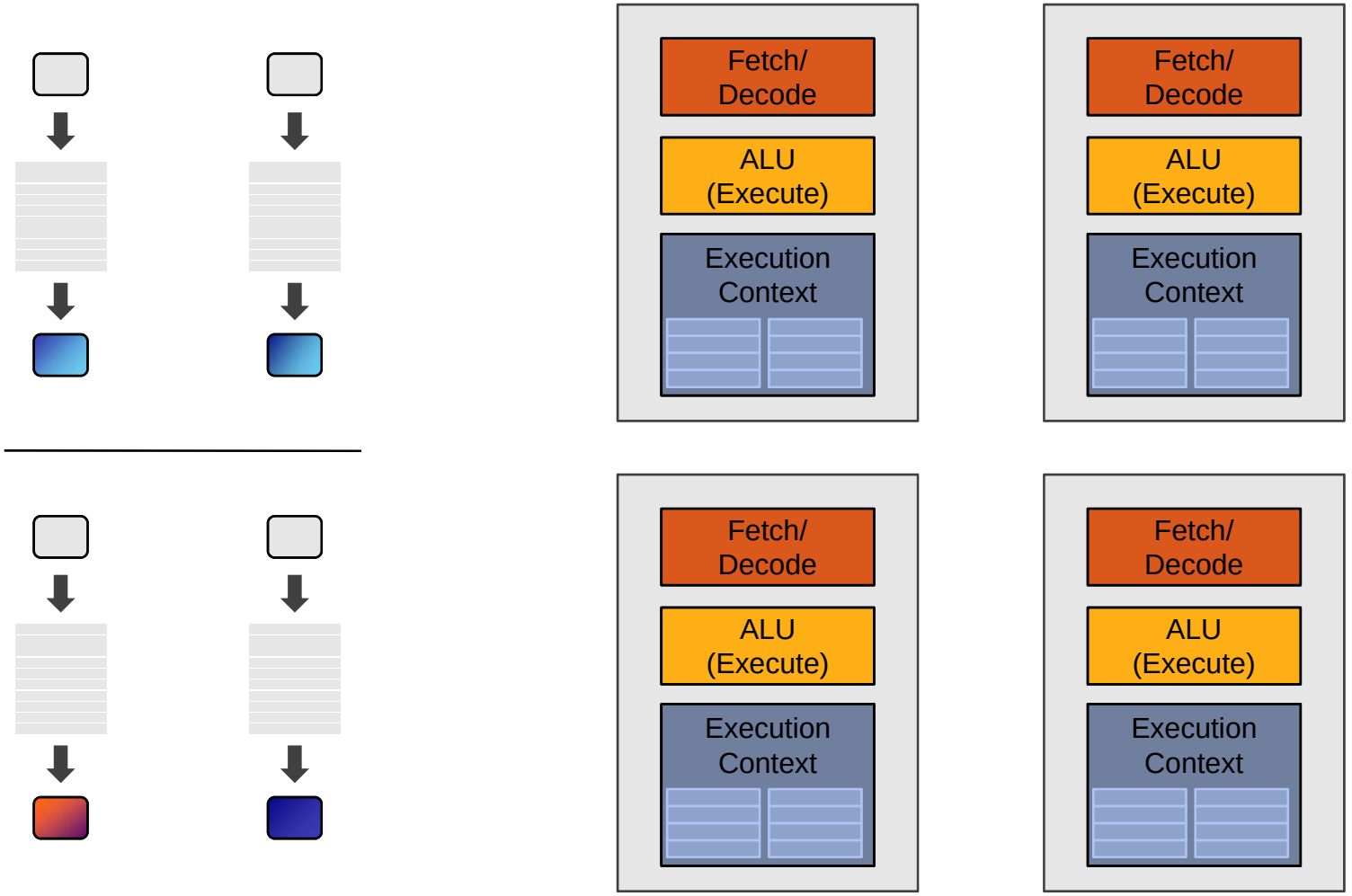
Idea #1:

Remove components that help a single instruction stream run fast

Two cores (two fragments in parallel)



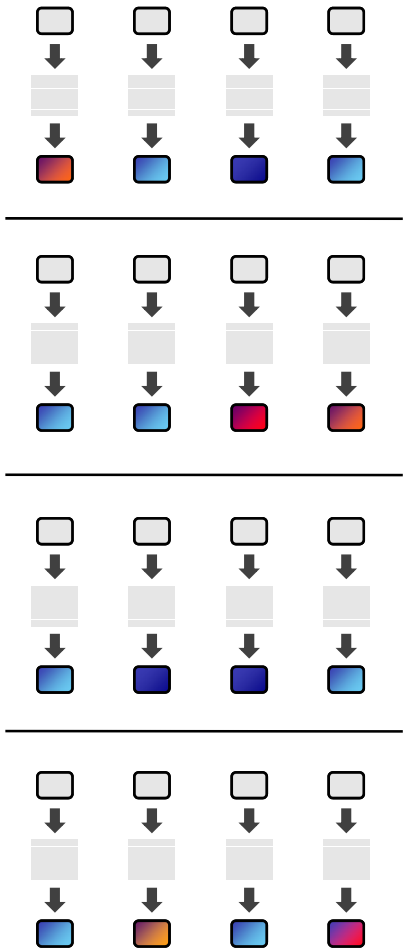
Four cores (four fragments in parallel)



Sixteen cores (sixteen fragments in parallel)

16 cores = 16 simultaneous instruction streams

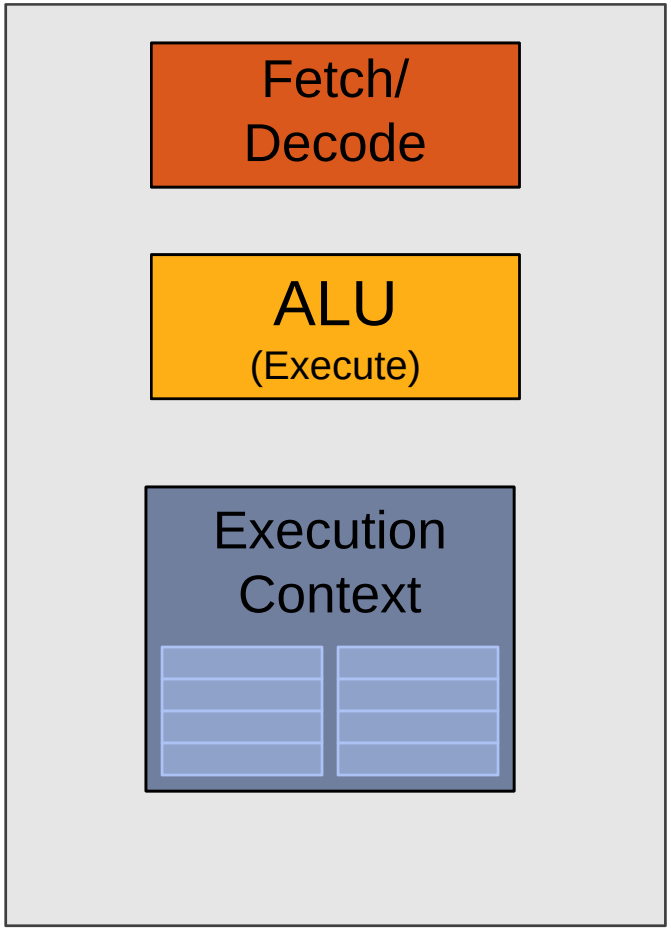
Instruction stream sharing



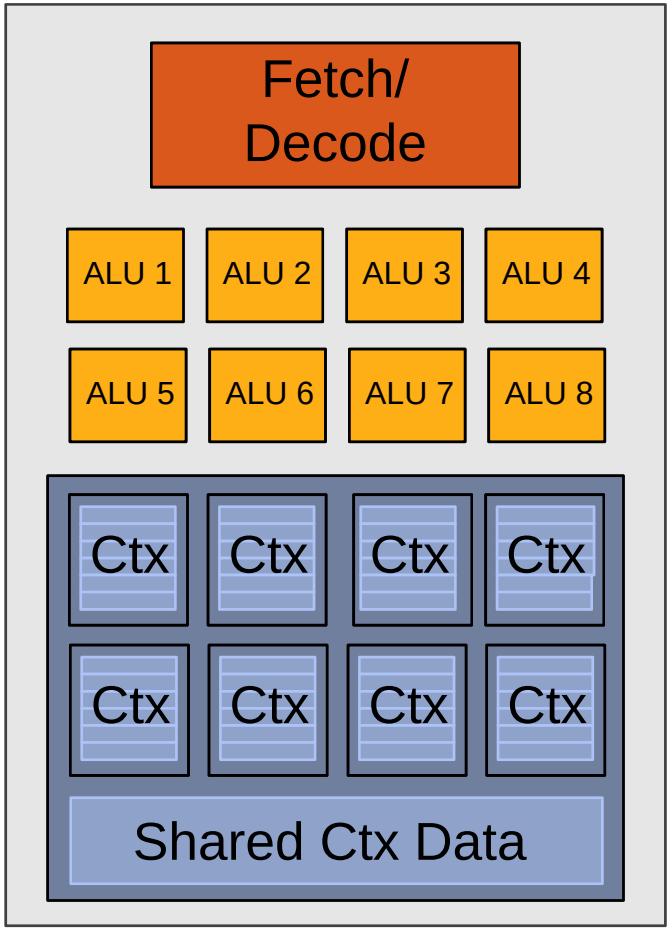
But… many fragments should be able to share an instruction stream!

```
<diffuseShader>:
sample r0, v4, t0, s0
mul  r3, v0, cb0[0]
madd r3, v1, cb0[1], r3
madd r3, v2, cb0[2], r3
clmp r3, r3, l(0.0), l(1.0)
mul  o0, r0, r3
mul  o1, r1, r3
mul  o2, r2, r3
mov  o3, l(1.0)
```

Recall: simple processing core



Add ALUs



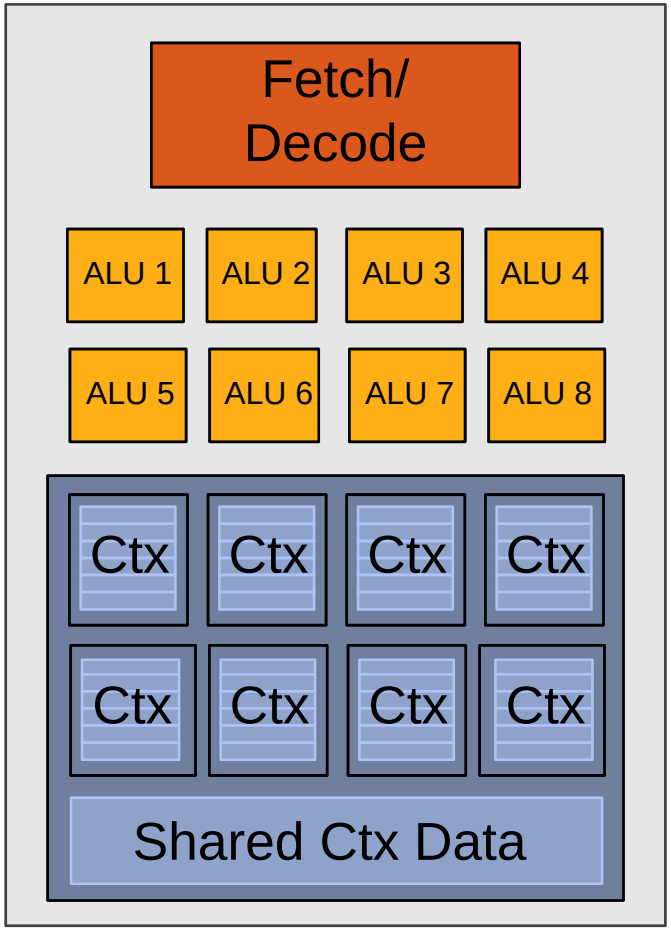
Idea #2:

Amortize cost/complexity of managing an instruction stream across many ALUs

SIMD processing

(or SIMT, SPMD)

Modifying the shader

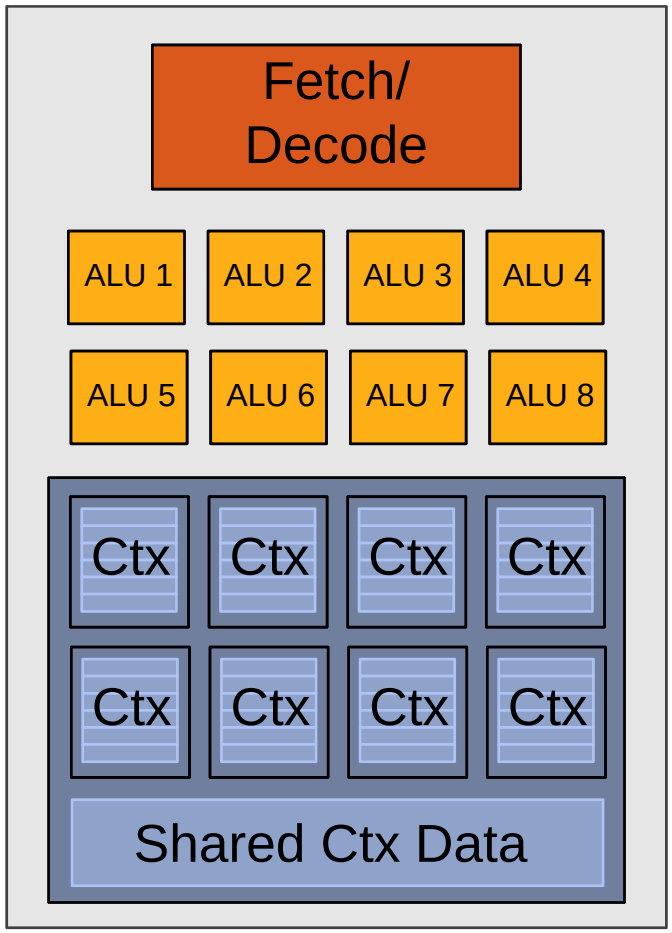



```
<diffuseShader>:
sample r0, v4, t0, s0
mul  r3, v0, cb0[0]
madd r3, v1, cb0[1], r3
madd r3, v2, cb0[2], r3
clmp r3, r3, l(0.0), l(1.0)
mul  o0, r0, r3
mul  o1, r1, r3
mul  o2, r2, r3
mov  o3, l(1.0)
```

Original compiled shader:

Processes one fragment using scalar ops on scalar registers

Modifying the shader

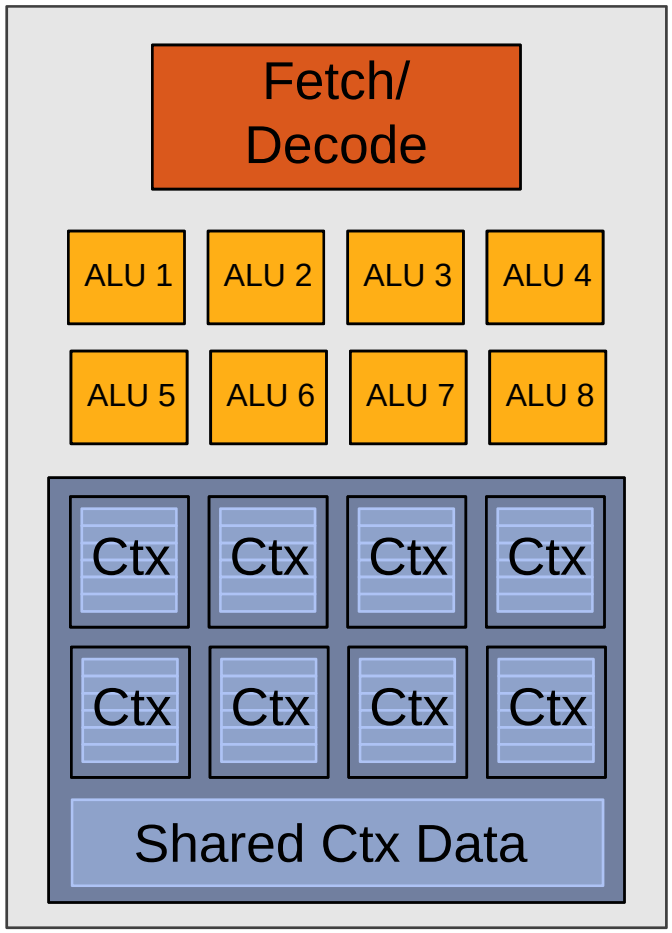


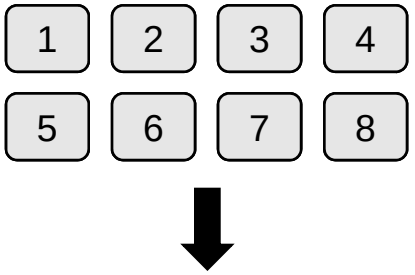

```
<VEC8_diffuseShader>:
VEC8_sample vec_r0, vec_v4, t0, vec_s0
VEC8_mul  vec_r3, vec_v0, cb0[0]
VEC8_madd vec_r3, vec_v1, cb0[1], vec_r3
VEC8_madd vec_r3, vec_v2, cb0[2], vec_r3
VEC8_clmp vec_r3, vec_r3, l(0.0), l(1.0)
VEC8_mul  vec_o0, vec_r0, vec_r3
VEC8_mul  vec_o1, vec_r1, vec_r3
VEC8_mul  vec_o2, vec_r2, vec_r3
VEC8_mov  vec_o3, l(1.0)
```

New compiled shader:

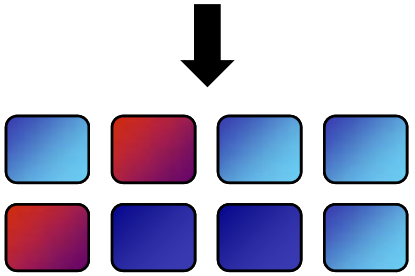
Processes 8 fragments using vector ops on vector registers

Modifying the shader

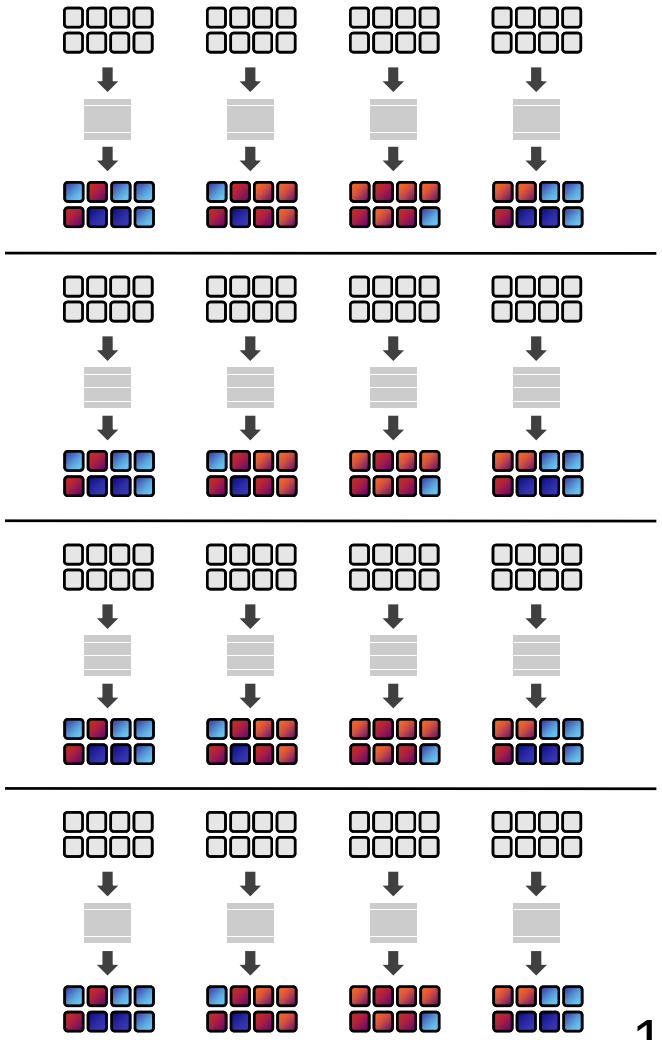


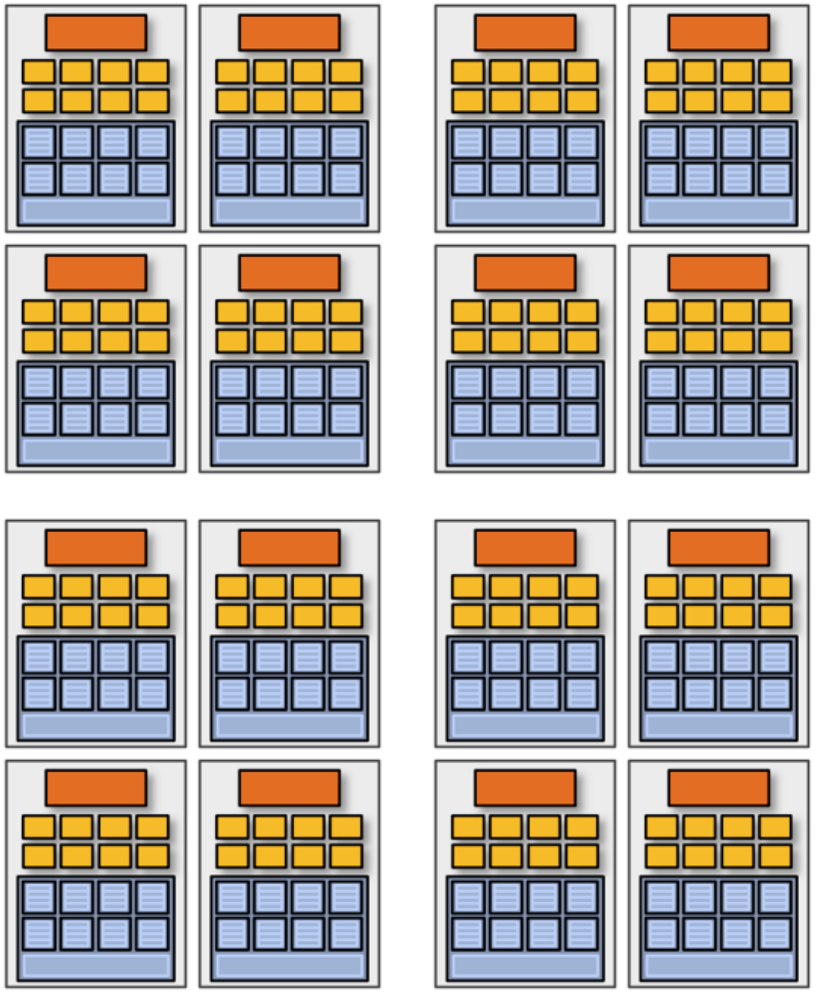



```
<VEC8_diffuseShader>:
VEC8_sample vec_r0, vec_v4, t0, vec_s0
VEC8_mul   vec_r3, vec_v0, cb0[0]
VEC8_madd vec_r3, vec_v1, cb0[1], vec_r3
VEC8_madd vec_r3, vec_v2, cb0[2], vec_r3
VEC8_clmp vec_r3, vec_r3, l(0.0), l(1.0)
VEC8_mul   vec_o0, vec_r0, vec_r3
VEC8_mul   vec_o1, vec_r1, vec_r3
VEC8_mul   vec_o2, vec_r2, vec_r3
VEC8_mov   vec_o3, l(1.0)
```



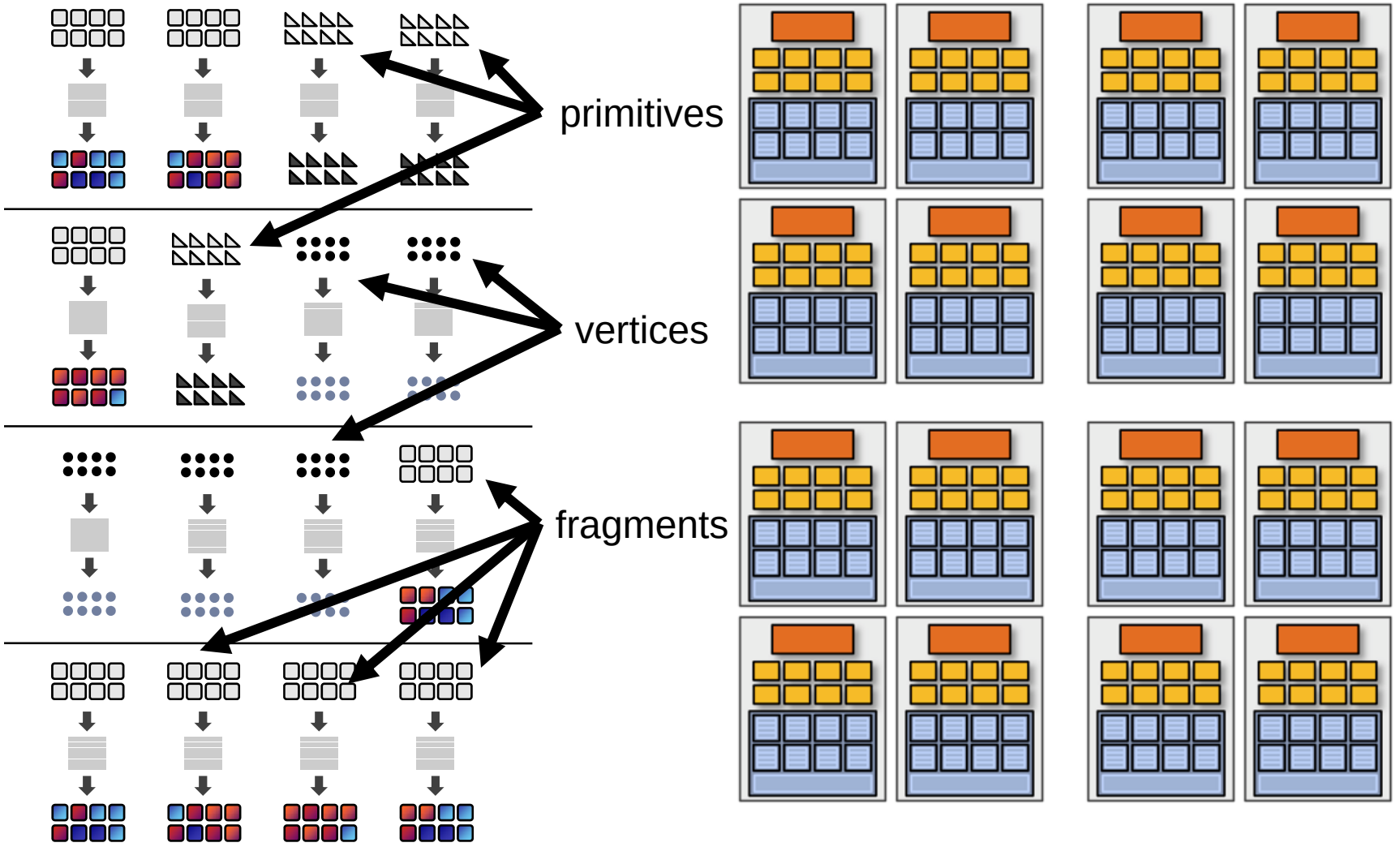
128 fragments in parallel





16 cores = 128 ALUs
= 16 simultaneous instruction streams

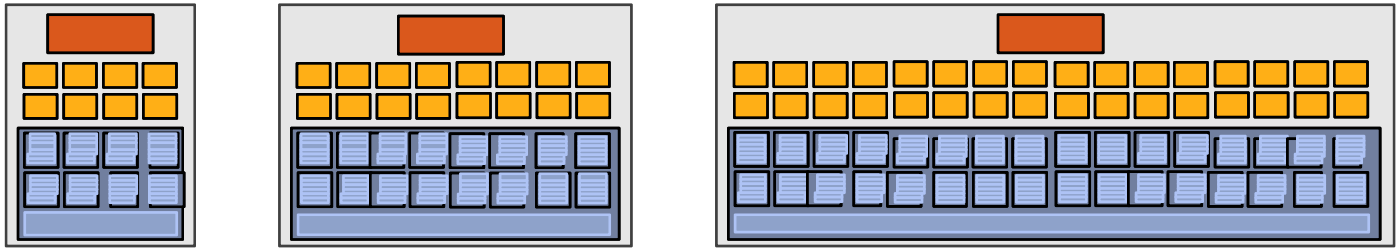
128 [vertices / fragments
primitives
CUDA threads
OpenCL work items
compute shader threads] in parallel



Clarification

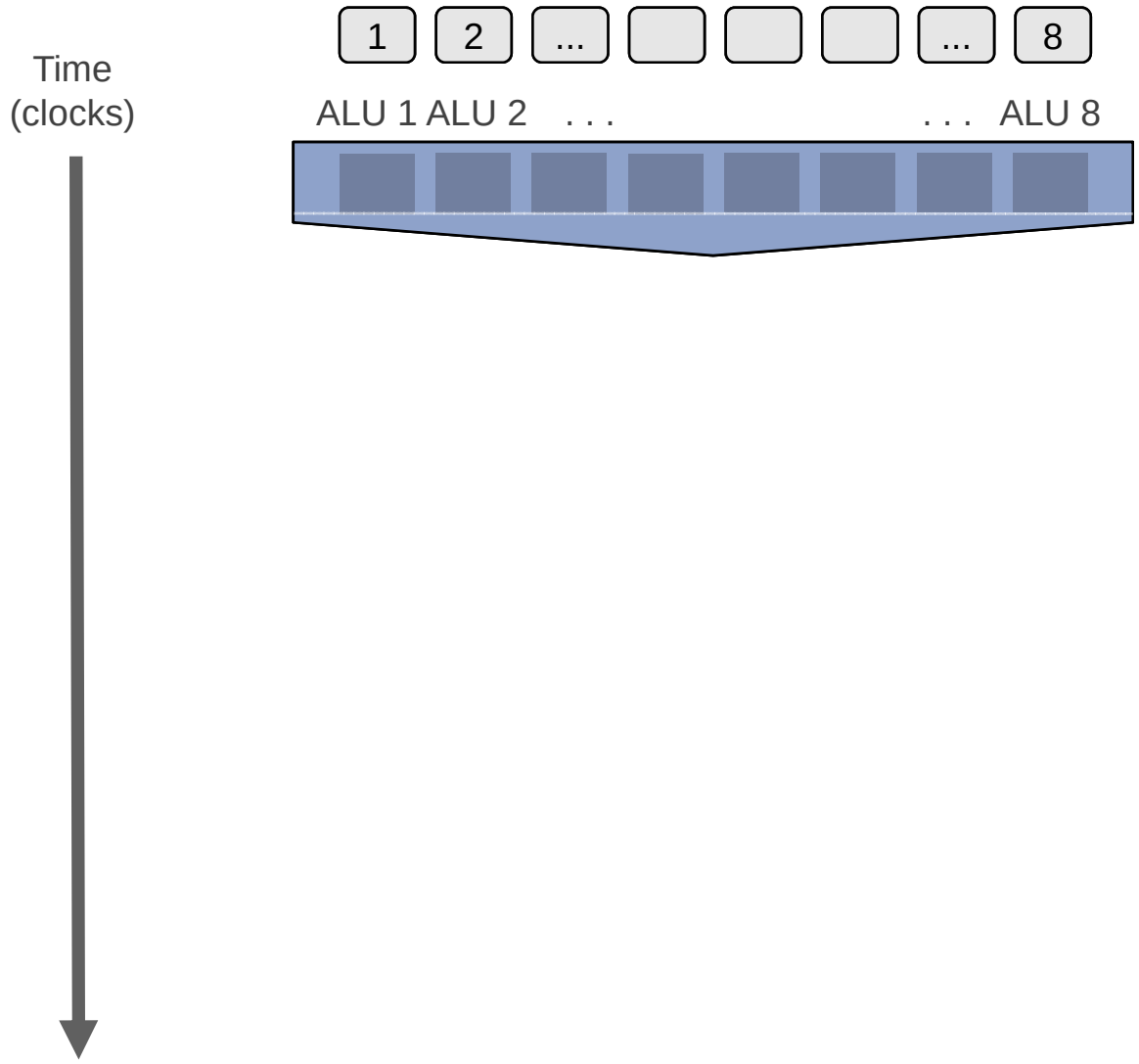
SIMD processing does not imply SIMD instructions

- Option 1: Explicit vector instructions
  - Intel/AMD x86 SSE, Intel Larrabee
- Option 2: Scalar instructions, implicit HW vectorization
  - HW determines instruction stream sharing across ALUs (amount of sharing hidden from software)
  - NVIDIA GeForce (“SIMT” warps), AMD Radeon architectures



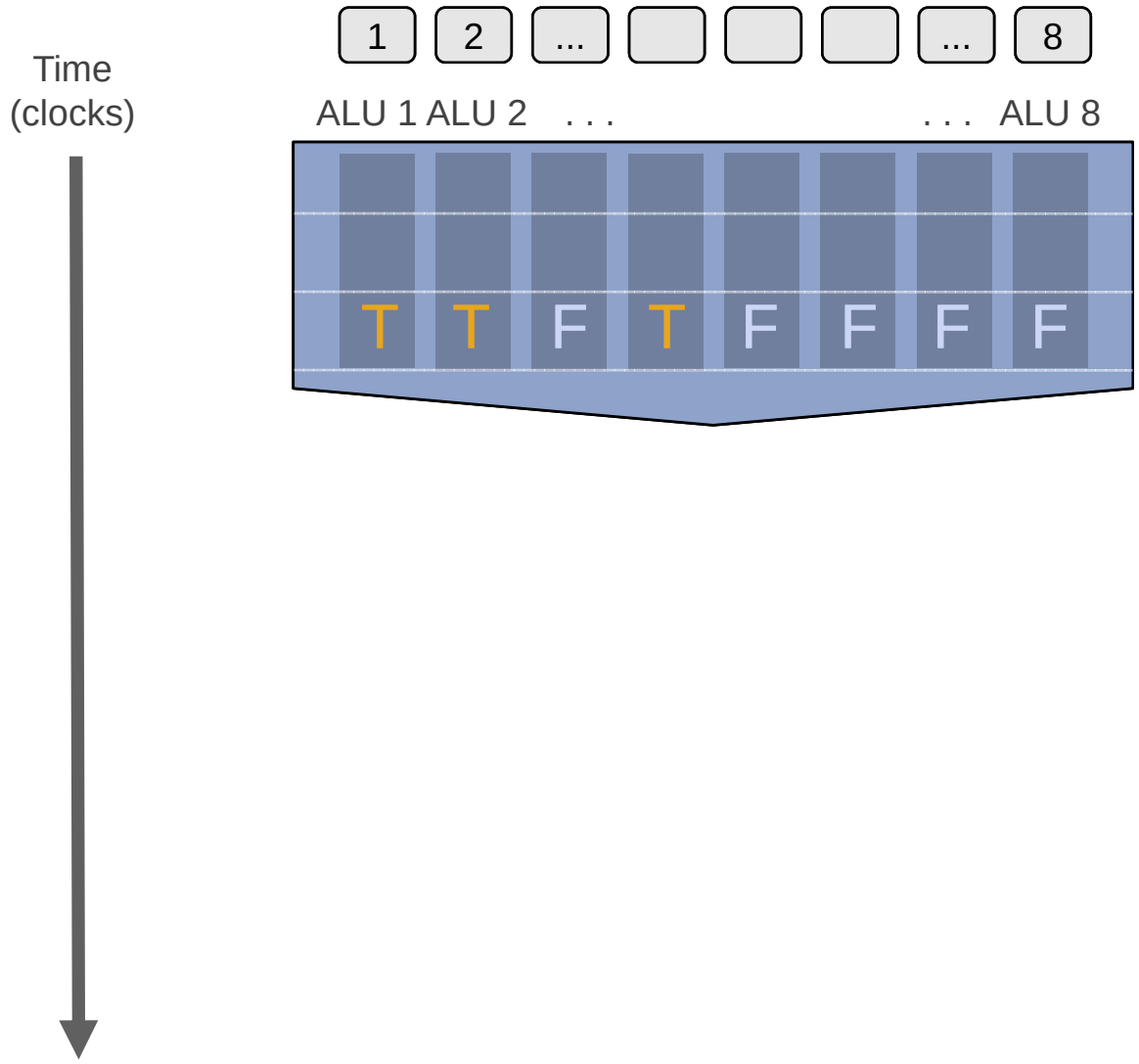
In practice: 16 to 64 fragments share an instruction stream

But what about branches?



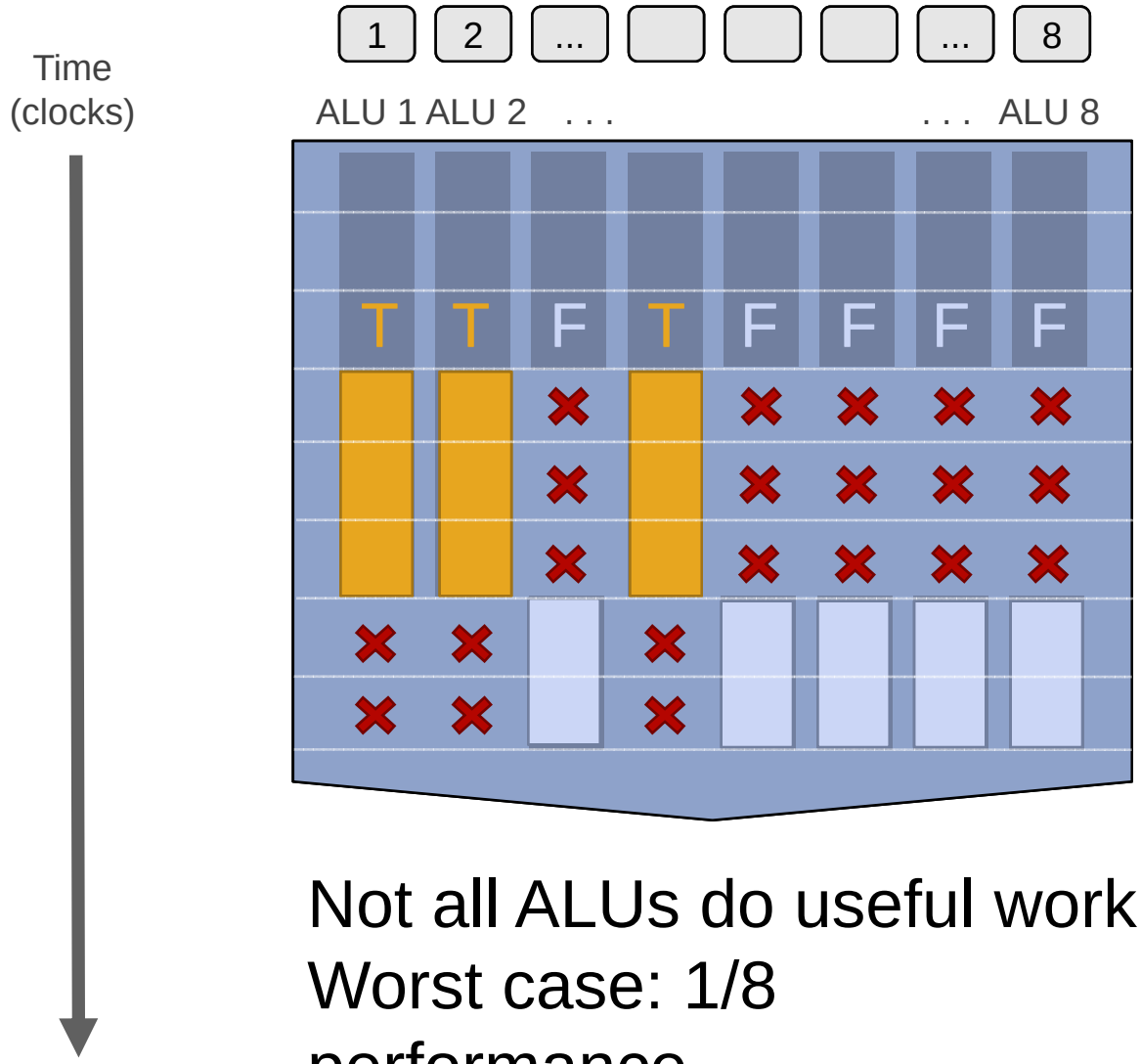

```
<unconditional
shader code>

if (x > 0) {
    y = pow(x, exp);
    y *= Ks;
    refl = y + Ka;
} else {
    x = 0;
    refl = Ka;
}

<resume unconditional
shader code>
```

But what about branches?




```
<unconditional
shader code>

if (x > 0) {
    y = pow(x, exp);
    y *= Ks;
    refl = y + Ka;
} else {
    x = 0;
    refl = Ka;
}

<resume unconditional
shader code>
```

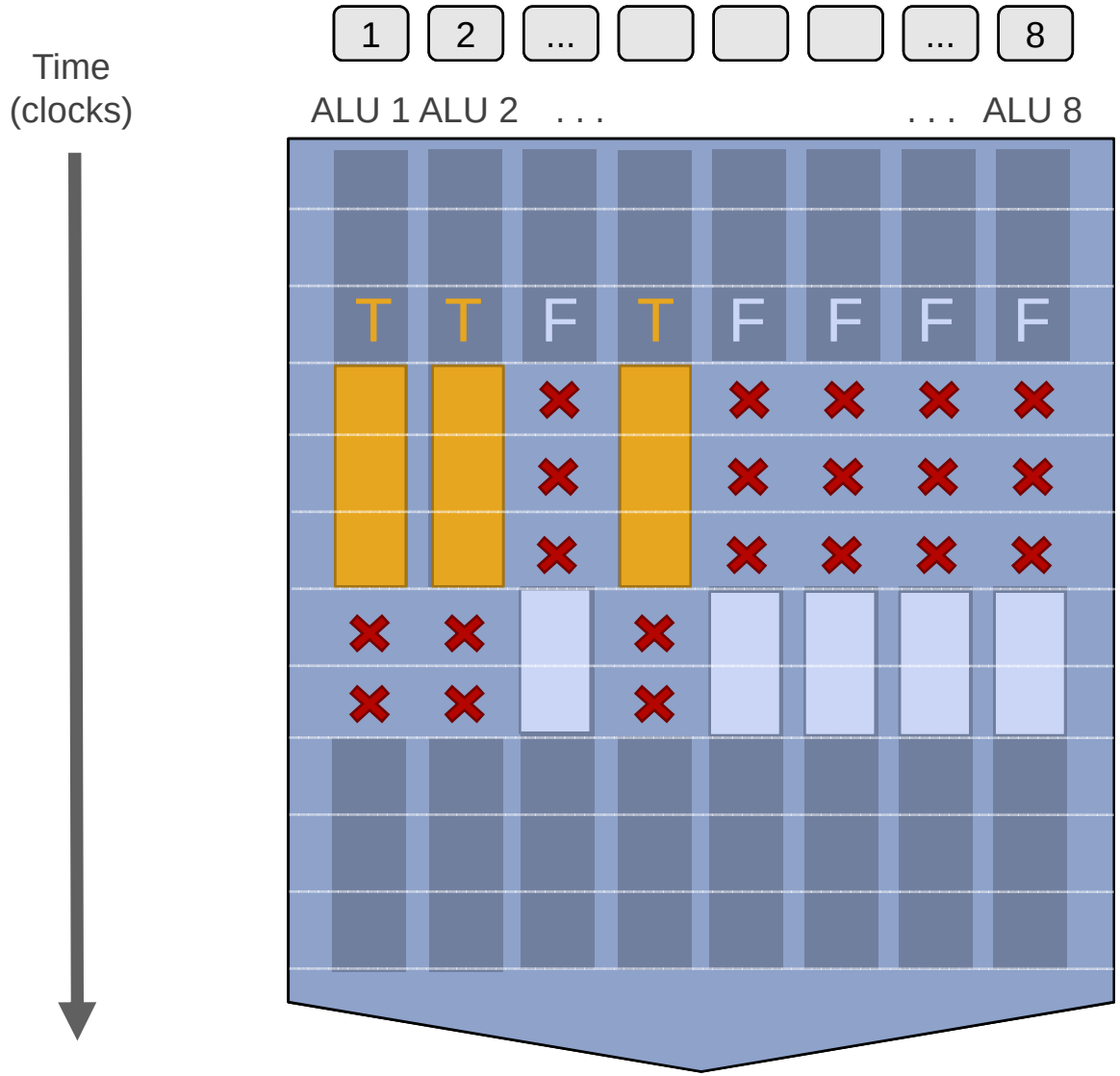
But what about branches?




```
<unconditional
shader code>

if (x > 0) {
    y = pow(x, exp);
    y *= Ks;
    refl = y + Ka;
} else {
    x = 0;
    refl = Ka;
}

<resume unconditional
shader code>
```

Not all ALUs do useful work!
Worst case: 1/8
performance

But what about branches?




```
<unconditional
shader code>

if (x > 0) {
    y = pow(x, exp);
    y *= Ks;
    refl = y + Ka;
} else {
    x = 0;
    refl = Ka;
}

<resume unconditional
shader code>
```

Stalls!

Stalls occur when a core cannot run the next instruction because of a dependency on a previous operation.

Texture access latency = 100’s to 1000’s of cycles

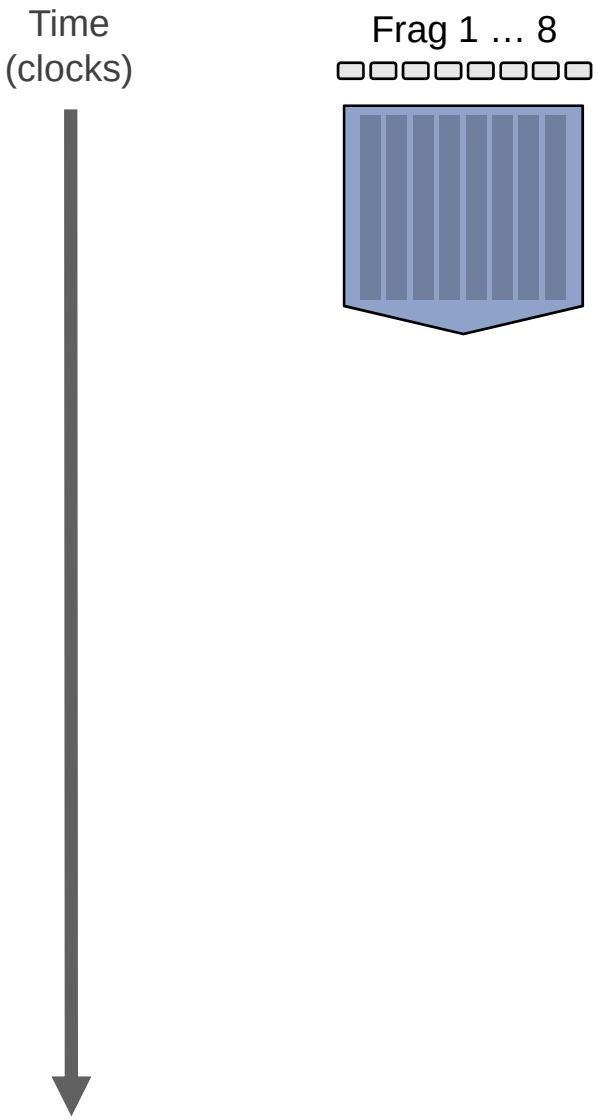
We’ve removed the fancy caches and logic that helps avoid stalls.

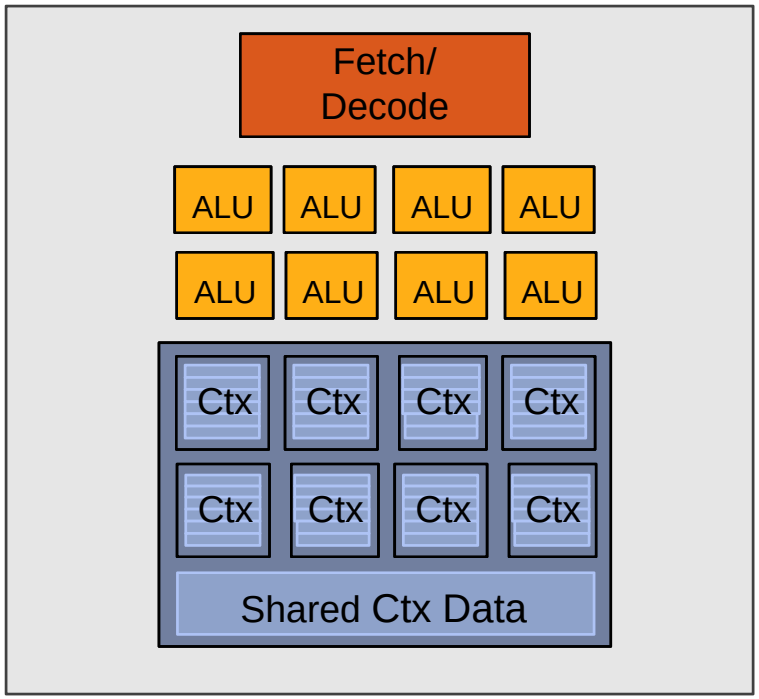
But we have LOTS of independent fragments.

Idea #3:

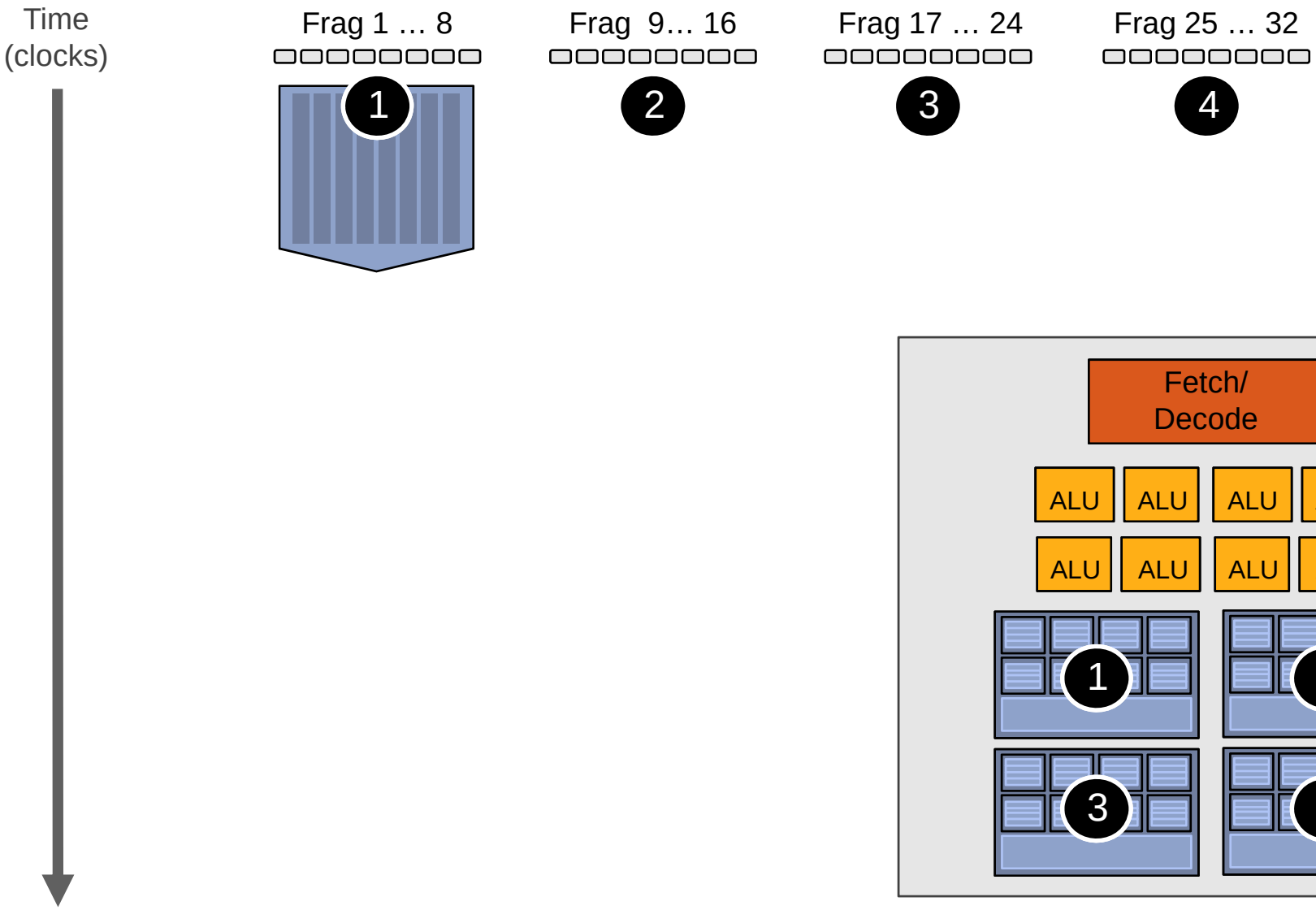
Interleave processing of many fragments on a single core to avoid stalls caused by high latency operations.

Hiding shader stalls

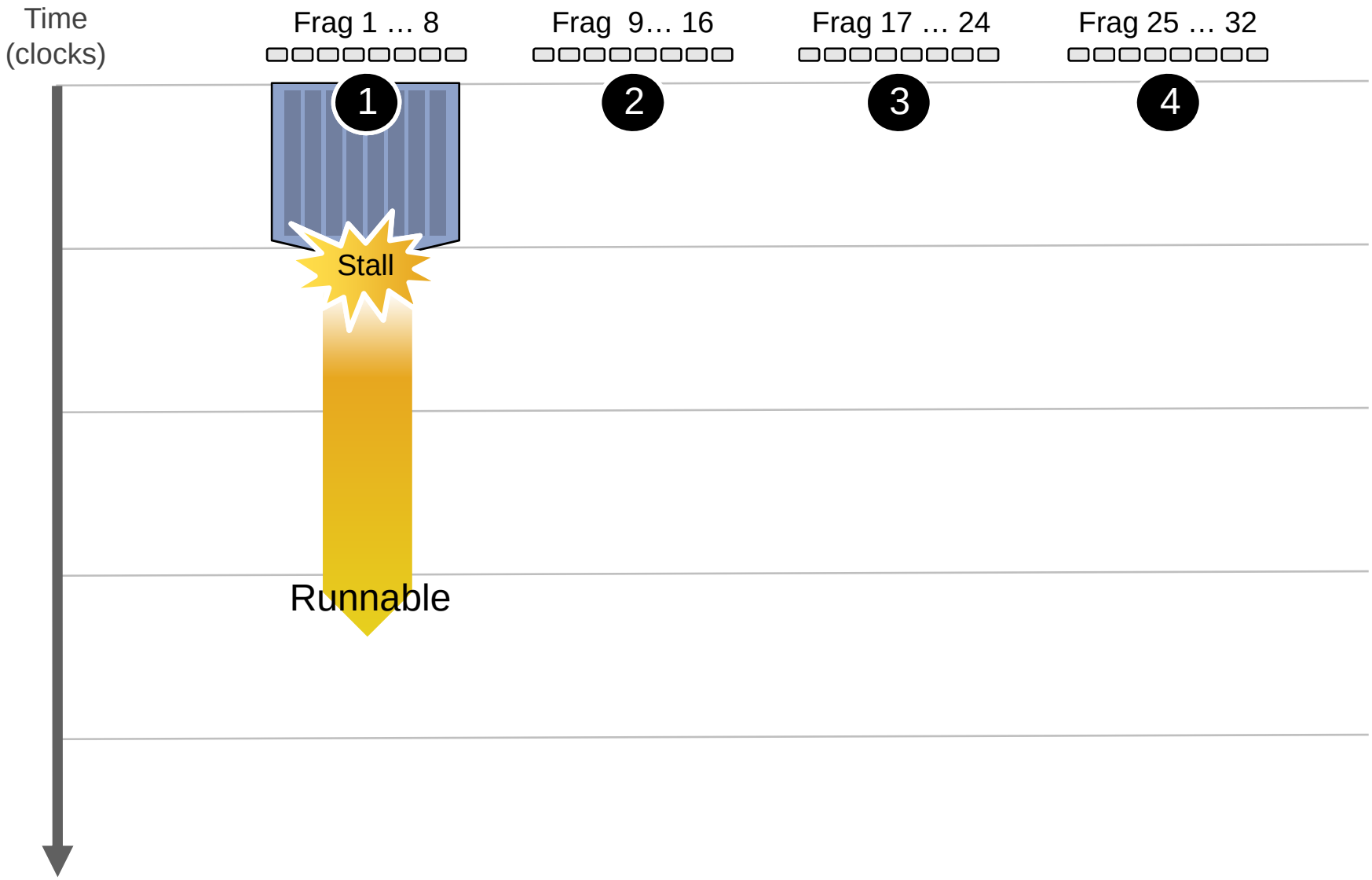




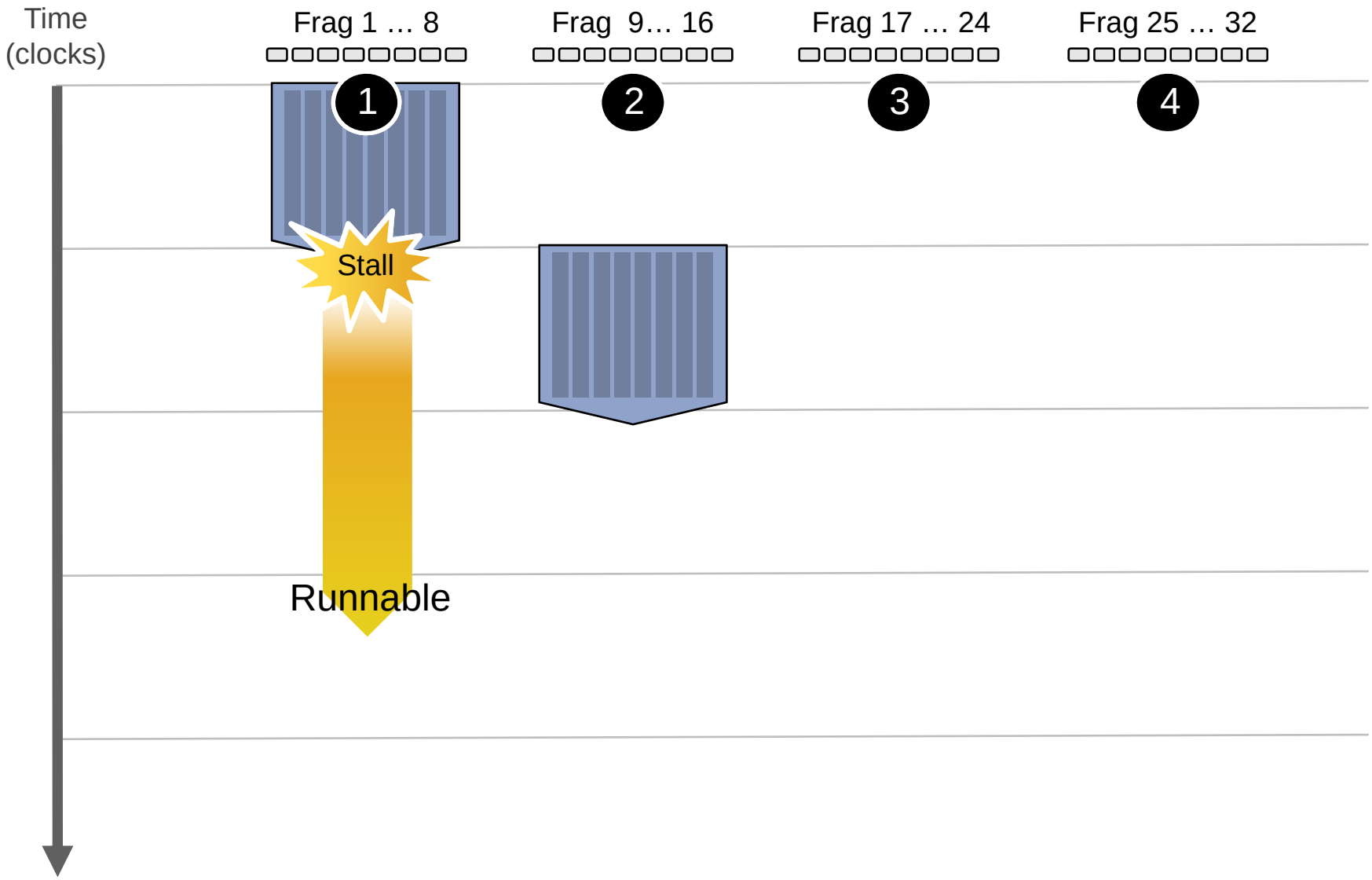
Hiding shader stalls



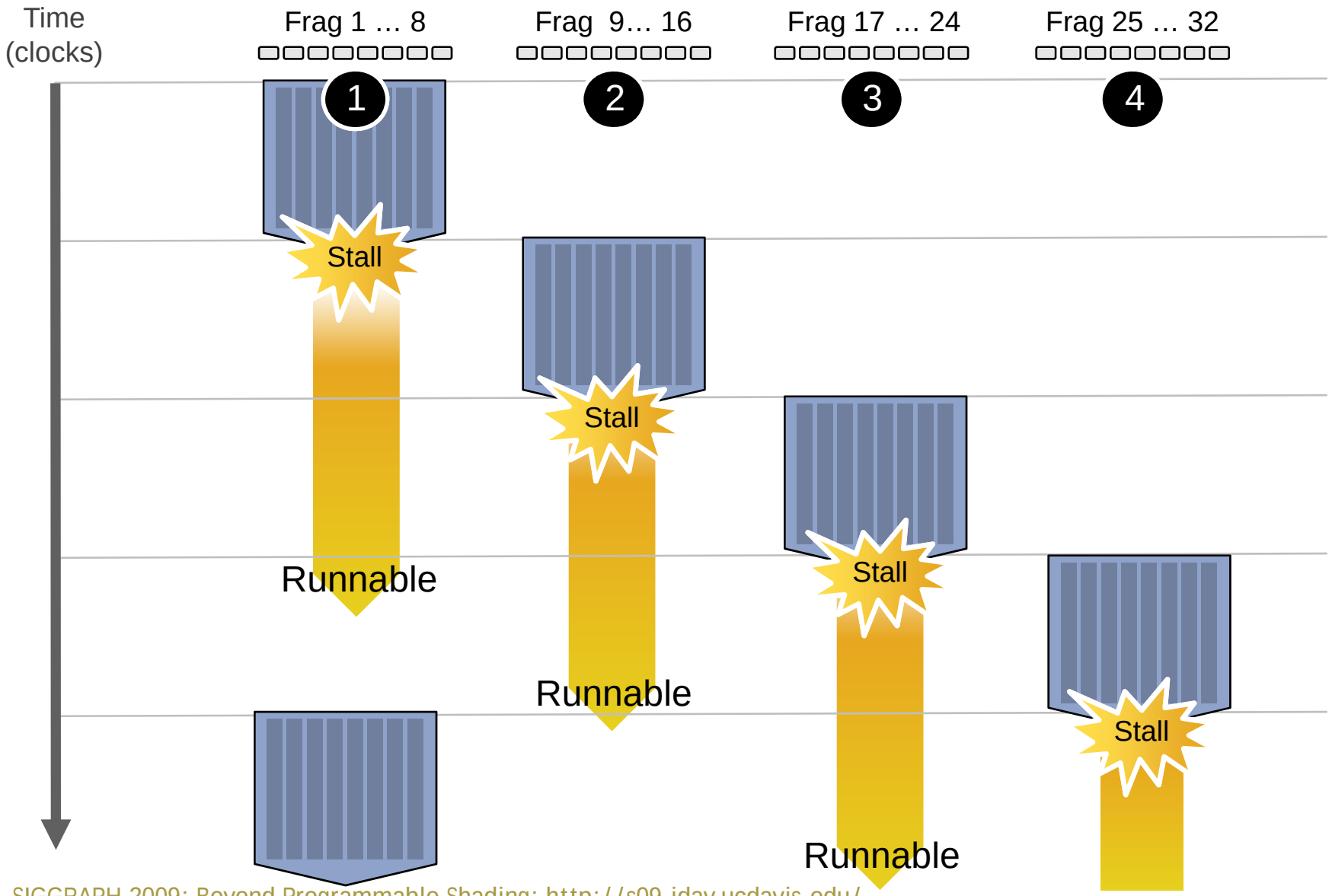
Hiding shader stalls



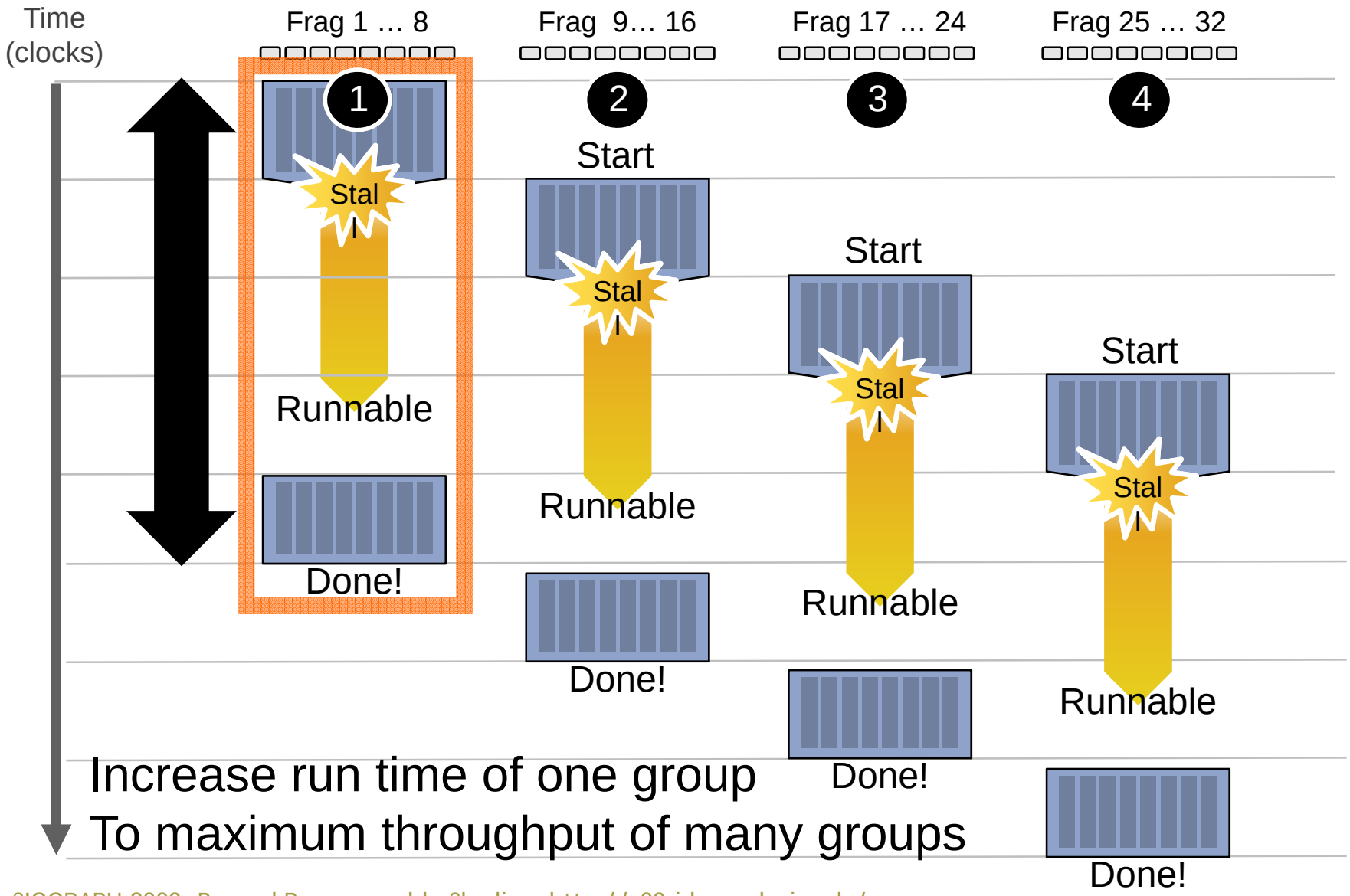
Hiding shader stalls



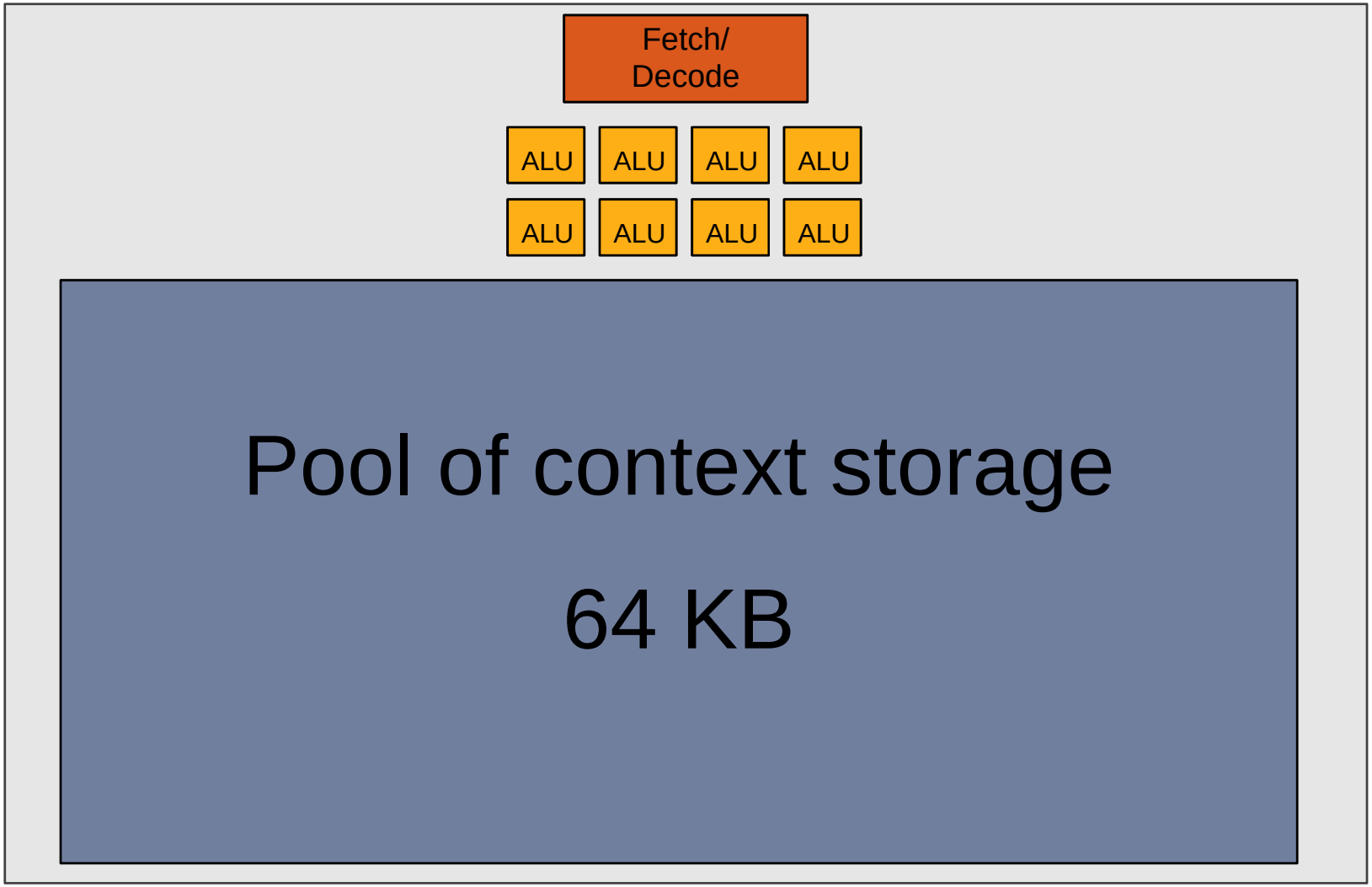
Hiding shader stalls



Throughput!

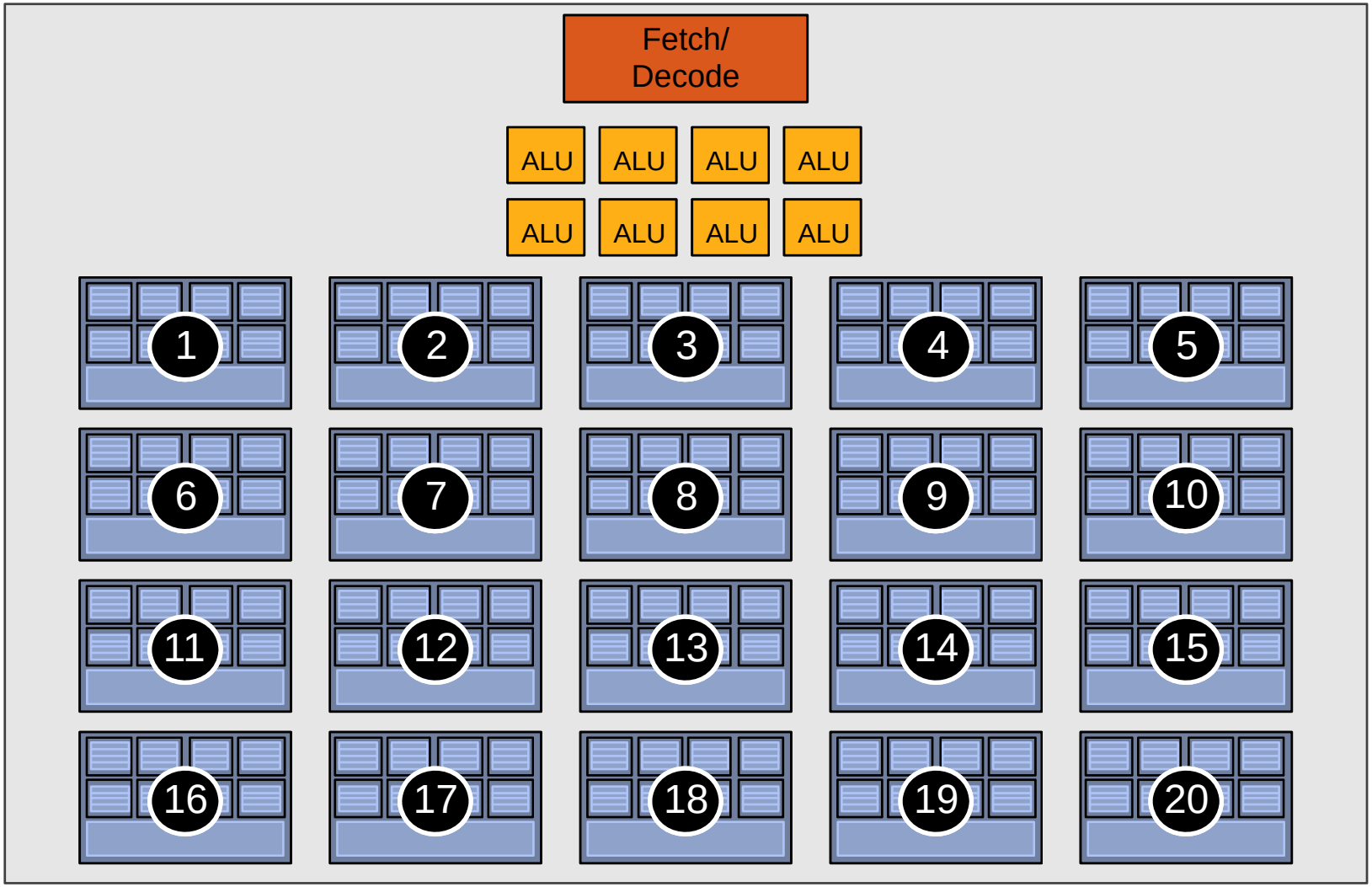

Time
(clocks)
Frag 1 … 8
Frag 9… 16
Frag 17 … 24
Frag 25 … 32
1
2
3
4
Stall
Runnable
Done!
Start
Stall
Runnable
Done!
Start
Stall
Runnable
Done!
Start
Stall
Runnable
Done!
Increase run time of one group
To maximum throughput of many groups

Storing contexts

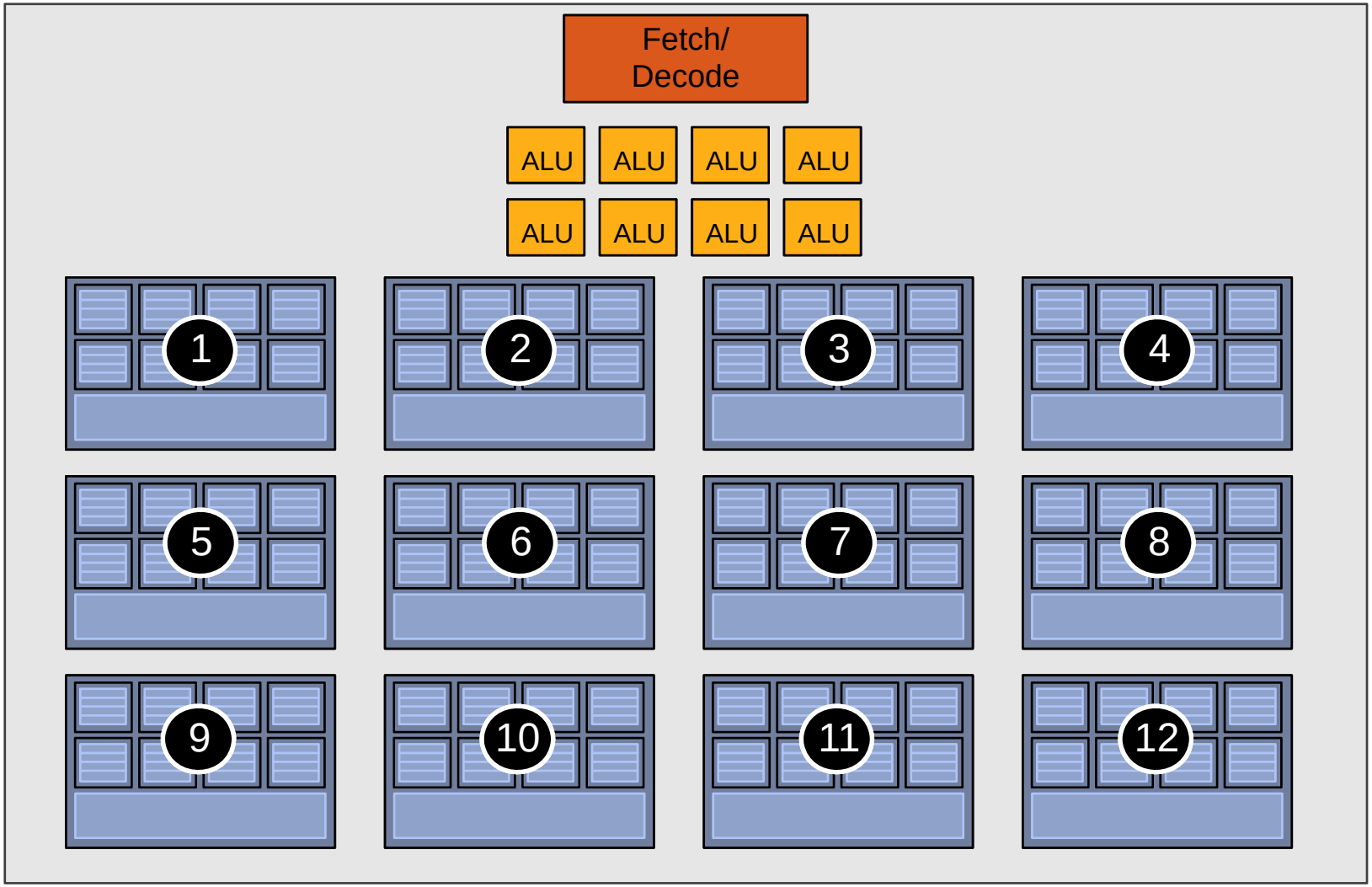


Twenty small contexts

(maximal latency hiding ability)

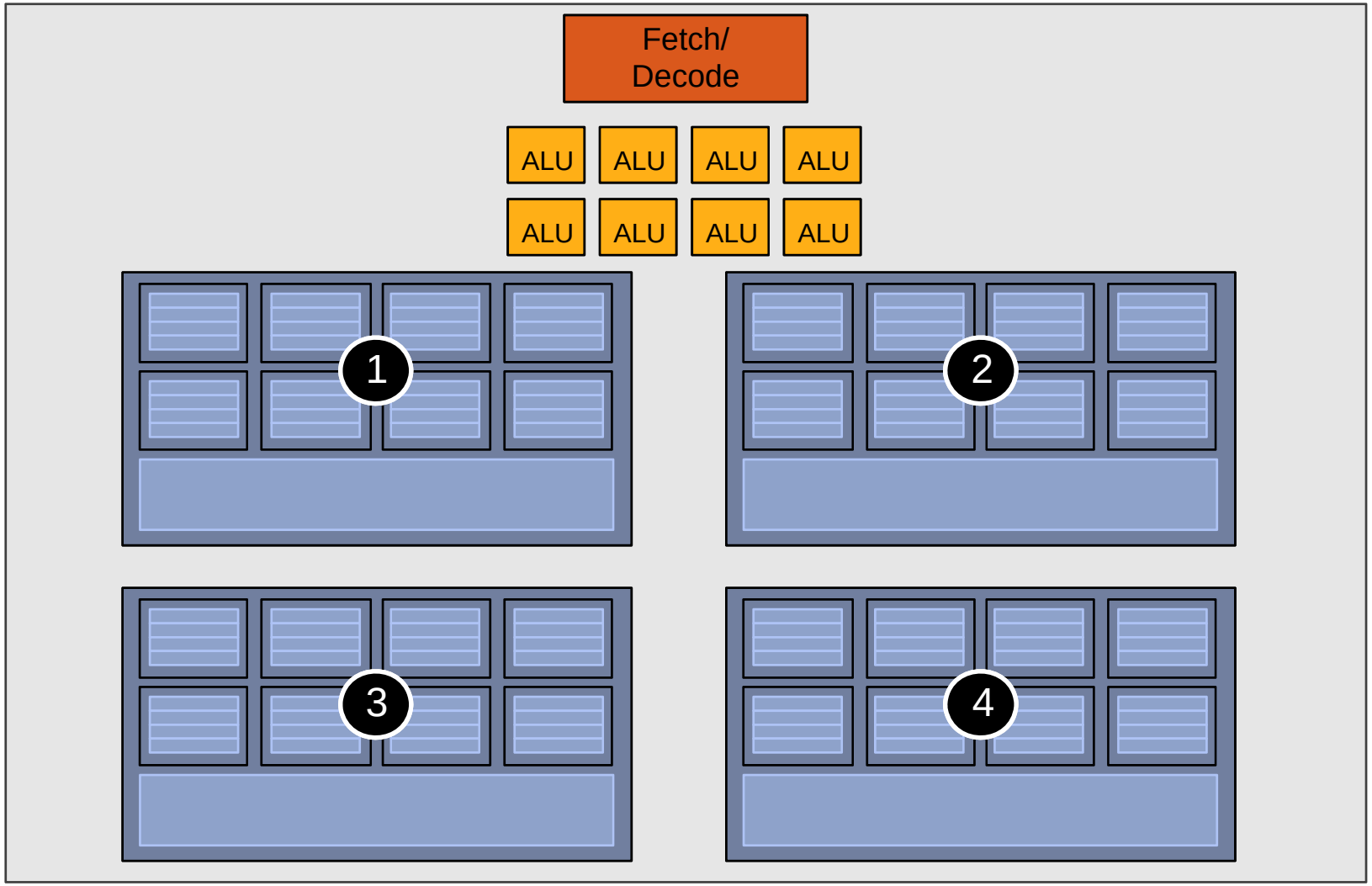


Twelve medium contexts



Four large contexts

(low latency hiding ability)



Clarification

Interleaving between contexts can be managed by HW or SW (or both!)

- NVIDIA / AMD Radeon GPUs
  - HW schedules / manages all contexts (lots of them)
  - Special on-chip storage holds fragment state
- Intel MIC/Larrabee
  - HW manages four x86 (big) contexts at fine granularity
  - SW scheduling interleaves many groups of fragments on each HW context
  - L1-L2 cache holds fragment state (as determined by SW)

My chip!

16 cores

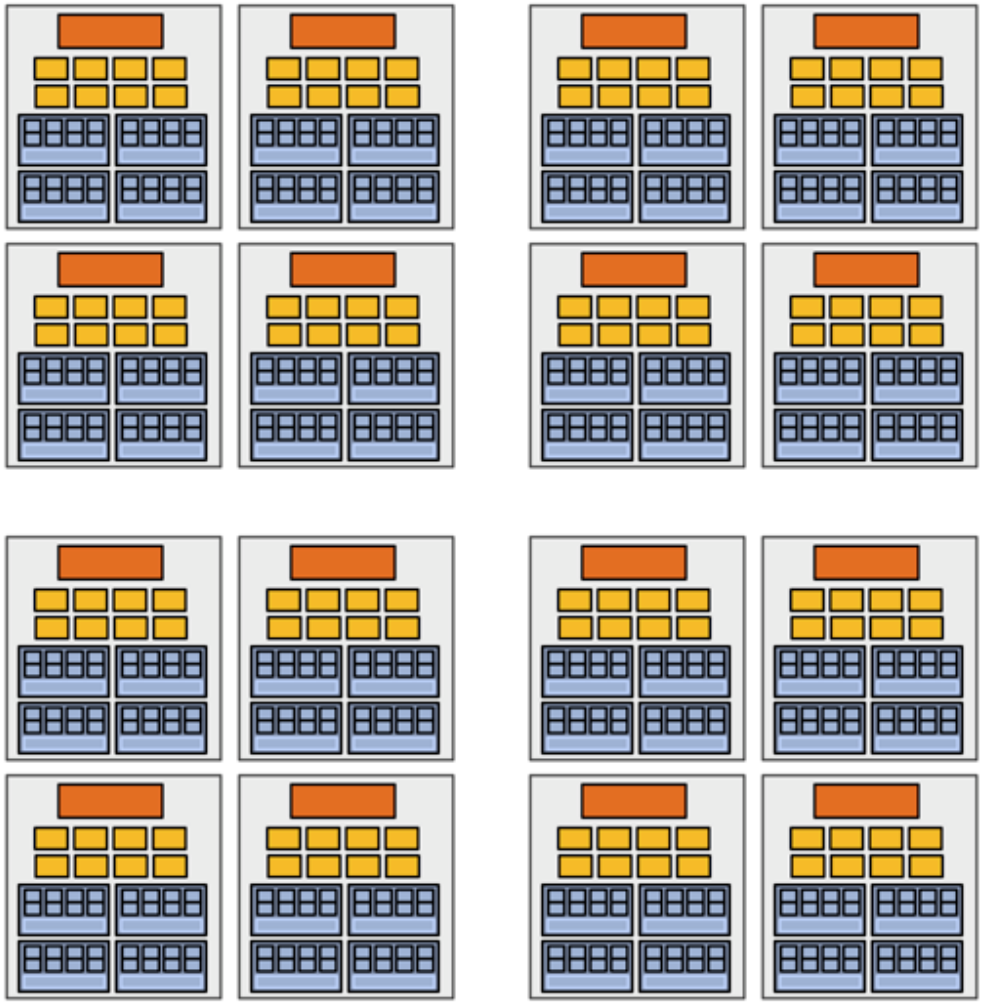
8 mul-add ALUs per core
(128 total)

16 simultaneous
instruction streams

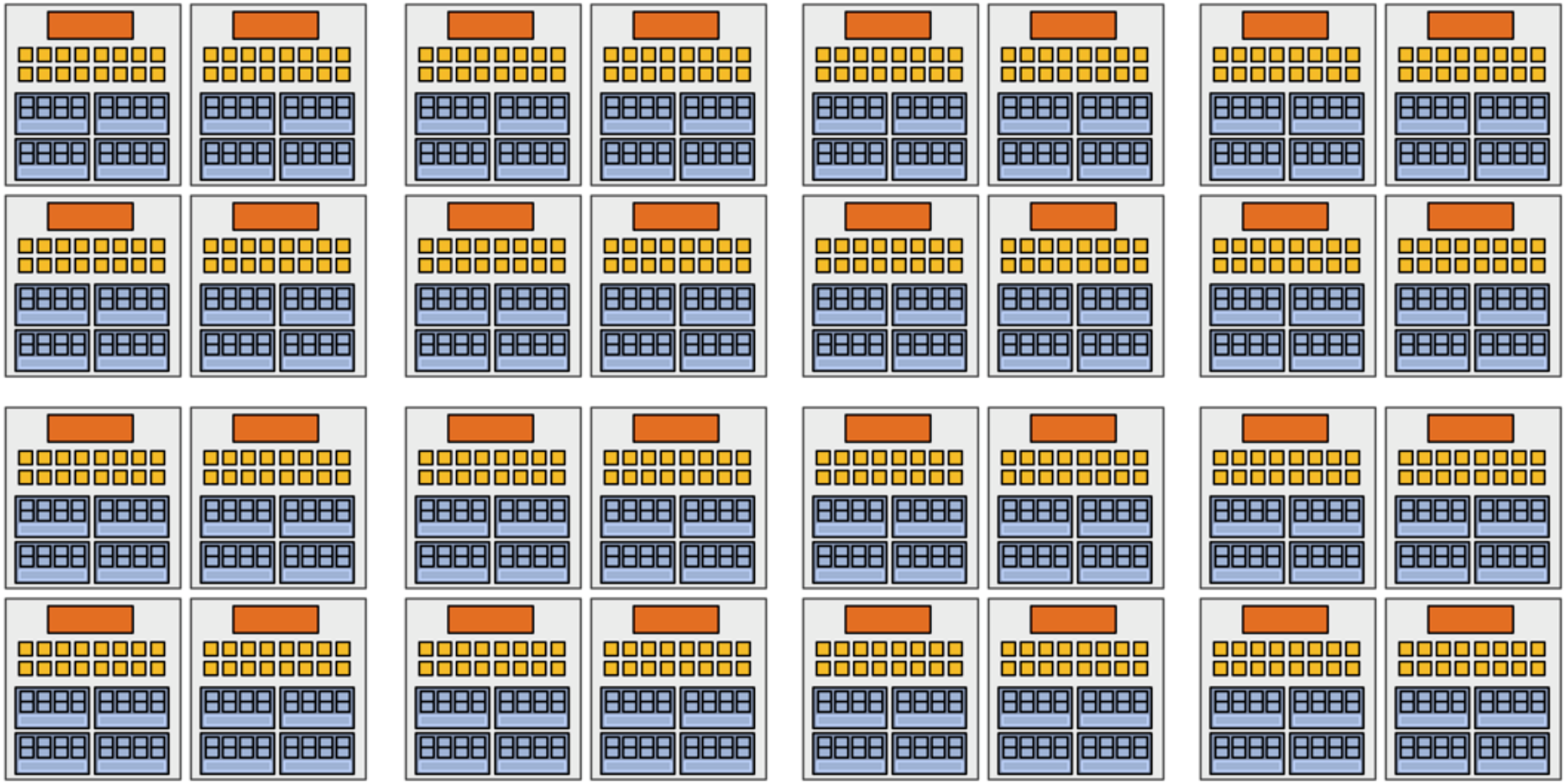
64 concurrent (but interleaved)
instruction streams

512 concurrent fragments

= 256 GFLOPs (@ 1GHz)



My “enthusiast” chip!



32 cores, 16 ALUs per core (512 total) = 1 TFLOP (@ 1 GHz)

Summary: three key ideas for high-throughput execution

1. Use many "slimmed down cores," run them in parallel

2. Pack cores full of ALUs (by sharing instruction stream overhead across groups of fragments)
   - Option 1: Explicit SIMD vector instructions
   - Option 2: Implicit sharing managed by hardware

3. Avoid latency stalls by interleaving execution of many groups of fragments
   - When one group stalls, work on another group

Thank you.