PPAR: CUDA basics

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This lecture: CUDA programming

We have seen some GPU architecture



Now how to program it ?

Outline

- GPU programming environments
- CUDA host side
- Parallel programming models: BSP, multi-BSP
- CUDA device side: threads, blocks, grids
- Expressing parallelism
 - Vector add example
- Managing communications
 - Parallel reduction example
- Re-using data
 - Matrix multiplication example

GPU development environments

For general-purpose programming (not graphics)

- Multiple toolkits
 - NVIDIA CUDA
 - Khronos OpenCL
 - Vulkan Compute
 - Microsoft DirectCompute
 - Google RenderScript
- Mostly syntactical variations
 - Underlying principles are the same
- In this course, focus on NVIDIA CUDA

Higher-level programming

- Directive-based
 - OpenACC
 - OpenMP 4.x
- Language extensions / libraries
 - Microsoft C++ AMP
 - Intel Cilk+
 - NVIDIA Thrust, CUB
- Languages
 - Intel ISPC

- - -

- Most corporations agree we need common standards...
 - But only if their own product becomes the standard!

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Hello World in CUDA

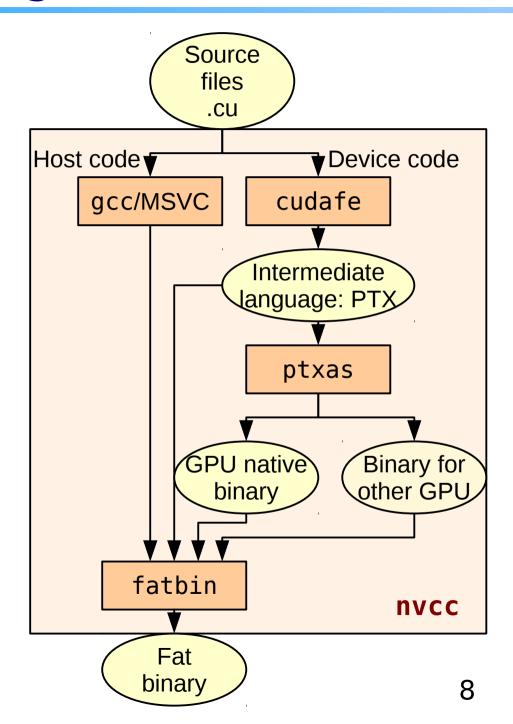
CPU "host" code + GPU "device" code

```
hello.cu:
__global__ void hello() {
    printf("Hello World!\n");
}
int main() {
    hello<<<1,1>>>();
    return 0;
}
Host code
```

Compiling a CUDA program

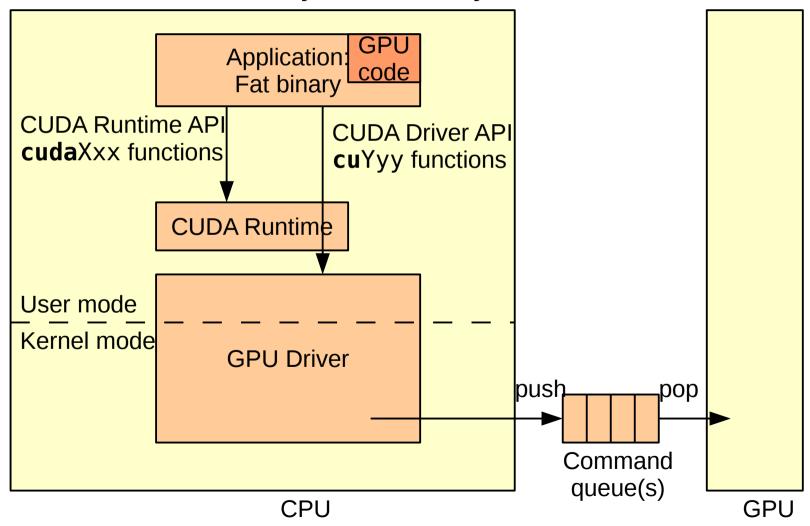
- Executable contains both host and device code
 - Device code in PTX and/or native
 - PTX can be recompiled on the fly (e.g. old program on new GPU)
- NVIDIA's compiler driver takes care of the process:

nvcc -o hello hello.cu

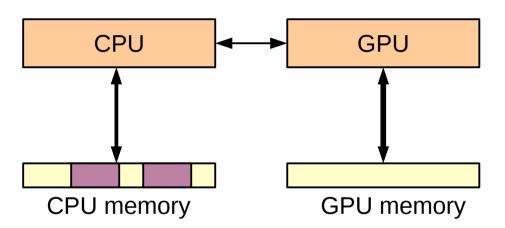


Control flow

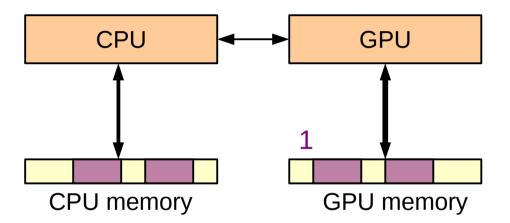
- Program running on CPUs
- Submit work to the GPU through the GPU driver
- Commands execute asynchronously



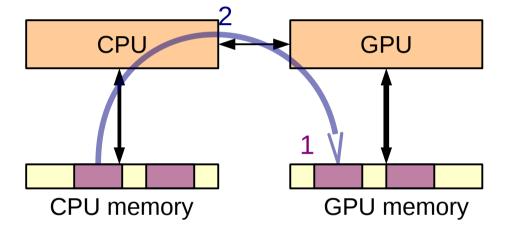
- Main program runs on the host
 - Manages memory transfers
 - Initiate work on GPU
- Typical flow



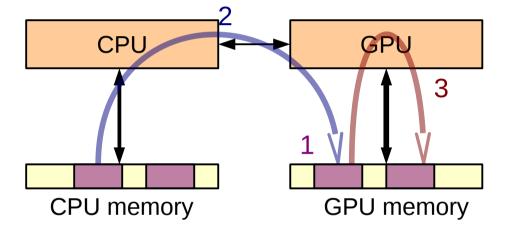
- Main program runs on the host
 - Manages memory transfers
 - Initiate work on GPU
- Typical flow
 - 1. Allocate GPU memory



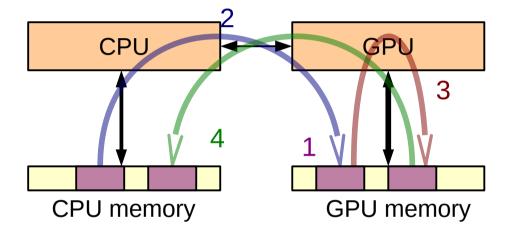
- Main program runs on the host
 - Manages memory transfers
 - Initiate work on GPU
- Typical flow
 - 1. Allocate GPU memory
 - 2. Copy inputs from CPU mem to GPU memory



- Main program runs on the host
 - Manages memory transfers
 - Initiate work on GPU
- Typical flow
 - 1. Allocate GPU memory
 - 2. Copy inputs from CPU mem to GPU memory
 - 3. Run computation on GPU



- Main program runs on the host
 - Manages memory transfers
 - Initiate work on GPU
- Typical flow
 - 1. Allocate GPU memory
 - 2. Copy inputs from CPU mem to GPU memory
 - 3. Run computation on GPU
 - 4. Copy back results to CPU memory



Example: a + b

- Our Hello World example did not involve the GPU
- Let's add up 2 numbers on the GPU
- Start from host code

```
int main()
{
    float ab[2] = {1515, 149}; // Inputs

    float c[1]; // Output
    // c[0] = ab[0] + ab[1];
    printf("c = %f\n", c[0]);
}
```

vectorAdd example: cuda/samples/0_Simple/vectorAdd

Step 1: allocate GPU memory

```
int main()
{
    float ab[2] = \{1515, 149\}, c[1]; // Inputs, in host mem
    // Allocate GPU memory
    float *d AB, *d C;
    cudaMalloc((void **)&d AB, 2*sizeof(float));
    cudaMalloc((void **)&d C, sizeof(float));
                                                  Allocate space
    Passing a pointer to the
                                                  for a, b and c
    pointer to be overwritten `
                                                  in GPU memory
                                                At the end,
    // Free GPU memory
                                                  free memory
    cudaFree(d AB);
    cudaFree(d C);
```

Step 2, 4: copy data to/from GPU memory

```
int main()
  float ab[2] = {1515, 149}, c[1]; // Inputs/outputs, CPU mem
  // Allocate GPU memory
  float *d AB, *d C;
  cudaMalloc((void **)&d AB, 2*sizeof(float));
  cudaMalloc((void **)&d C, sizeof(float));
  // Copy from CPU mem to GPU mem
  cudaMemcpy(d AB, ab, 2*sizeof(float), cudaMemcpyHostToDevice);
  // Copy results back to CPU mem
   cudaMemcpy(c, d C, sizeof(float), cudaMemcpyDeviceToHost);
  printf("c = %f\n", c[0]);
  // Free GPU memory
  cudaFree(d AB);
  cudaFree(d C);
```

Step 3: launch kernel

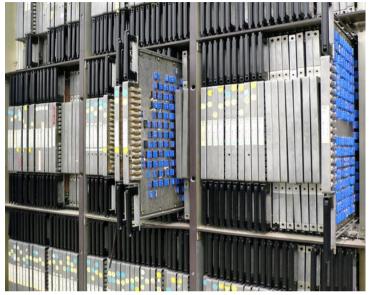
```
global void addOnGPU(float * ab, float * c)
     c[0] = ab[0] + ab[1];
int main()
                                                        Kernel is a function prefixed
   float ab[] = {1515, 159}; // Inputs, CPU mem
                                                            by global
   // Allocate GPU memory
   float *d AB, *d C;
                                                                Runs on GPU
   cudaMalloc((void **)&d AB, 2*sizeof(float));
   cudaMalloc((void **)&d C, sizeof(float));
   // Copy from CPU mem to GPU mem
                                                           Invoked from CPU code with
   cudaMemcpy(d AB, ab, 2*sizeof(float), cudaMemcpyHostToDevice);
                                                            <<>>> syntax
   // Launch computation on GPU
   add0nGPU<<<1, 1>>>(d_AB, d_C);
                                                          Note: we could have passed
                                                          a and b directly
              // Result on RU
   float c[1];
                                                          as kernel parameters
   // Copy results back to CPU mem
                             cudaMemcpyDeviceToHost);
   cudaMemcpy(c, d C, sizeof(float),
   printf("c = %f\n", c[0]);
   // Free GPU memory
   cudaFree(d AB);
                                                          What is inside the <<<>>>?
   cudaFree(d C);
}
```

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PRAM limitations

- PRAM model proposed in 1978
 - Inspired by SIMD machines of the time
- Assumptions
 - All processors synchronized every instruction
 - Negligible communication latency
- Useful as a theoretical model, but far from modern computers



ILLIAC-IV, an early SIMD machine

Modern architectures

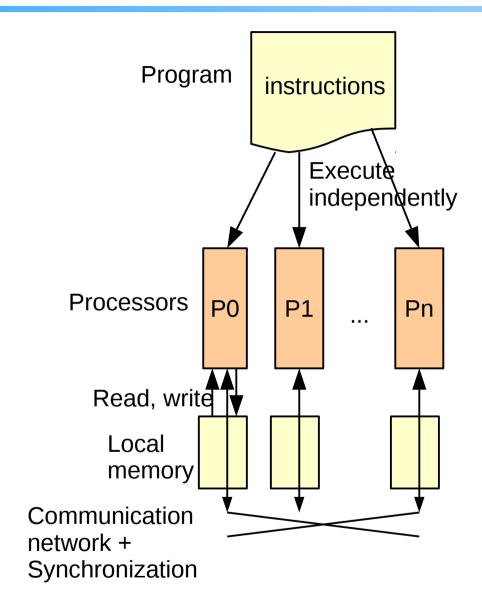
- Modern supercumputers are clusters of computers
 - Global synchronization costs millions of cycles
 - Memory is distributed
- Inside each node
 - Multi-core CPUs, GPUs
 - Non-uniform memory access (NUMA) memory
- Synchronization cost at all levels



Mare Nostrum, a modern distributed memory machine

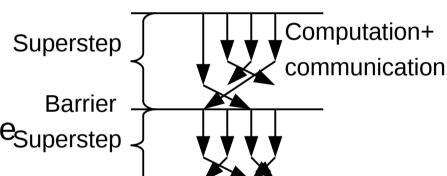
Bulk-Synchonous Parallel (BSP) model

- Assumes distributed memory
 - But also works with shared memory
 - Good fit for GPUs too, with a few adaptations
- Processors execute instructions independently
- Communications between processors are explicit
- Processors need to synchronize with each other



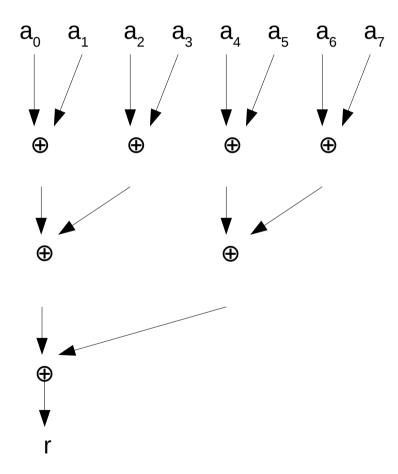
Superstep

- A program is a sequence of supersteps
- Superstep: each processor
 - Computes
 - Sends result
 - Receive data
- Barrier: wait until all processors have_{Superstep} finished their superstep
- Next superstep: can use data received in previous step



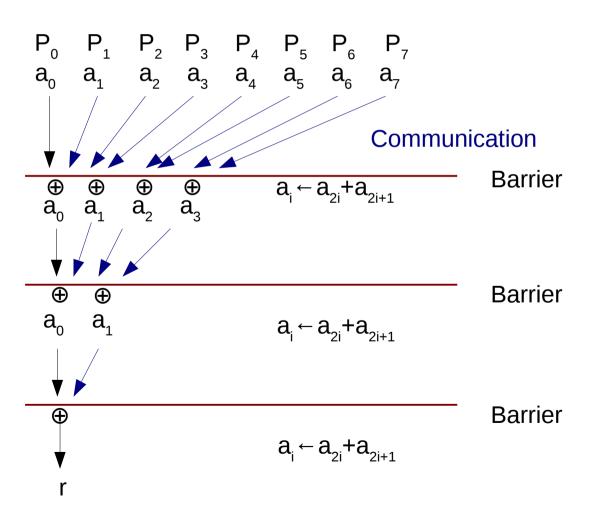
Example: reduction in BSP

Start from dependency graph



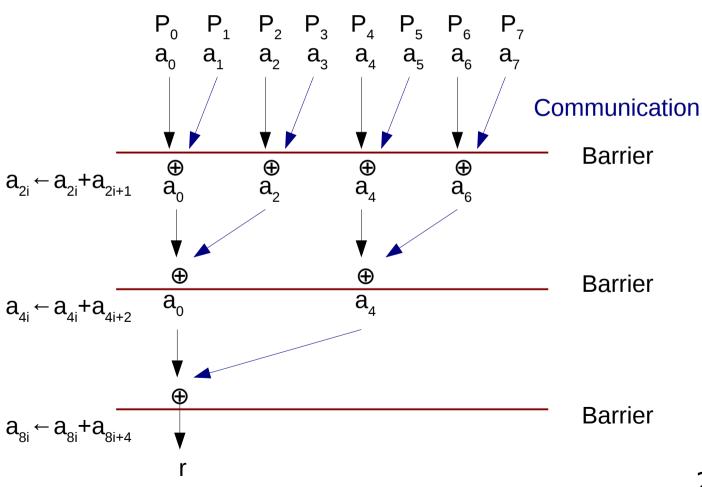
Reduction: BSP

Add barriers



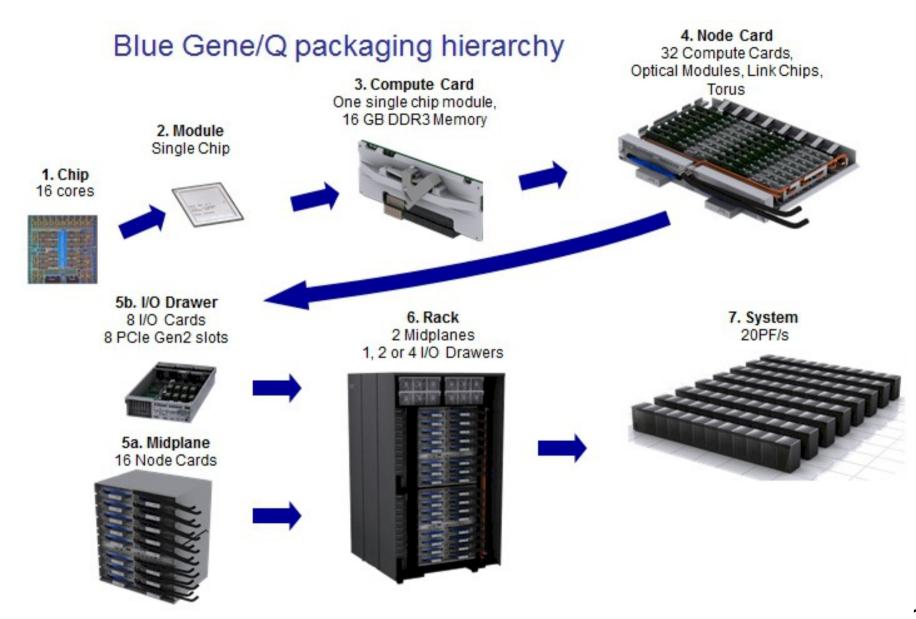
Reducing communication

- Data placement matters in BSP
- Optimization: keep left-side operand local



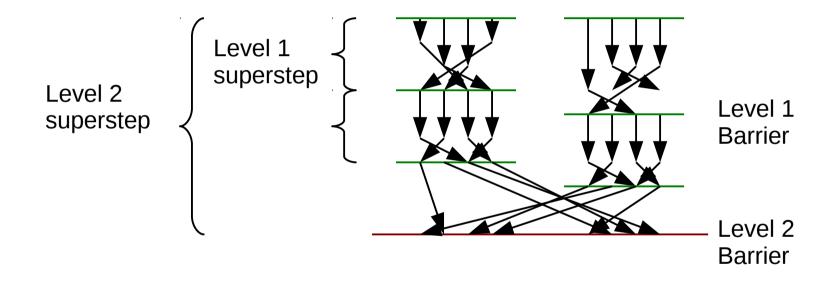
The world is not flat

It is hierarchical!



Multi-BSP model

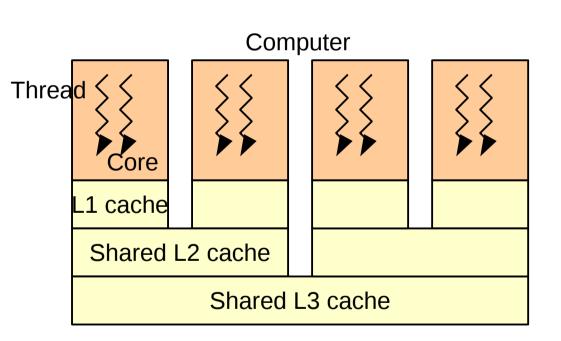
 Multi-BSP: BSP generalization with groups of processors in multiple nested levels

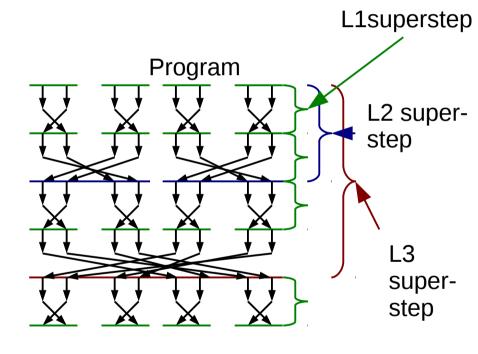


- Higher level: more expensive synchronization
- Arbitrary number of levels

Multi-BSP and multi-core

- Minimize communication cost on hierarchical platforms
 - Make parallel program hierarchical too
 - Take thread affinity into account

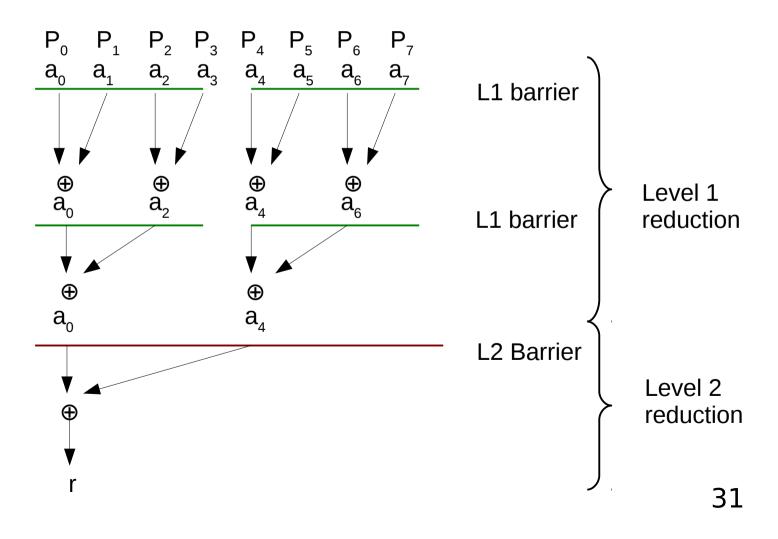




- On clusters (MPI): add more levels up
- On GPUs (CUDA): add more levels down

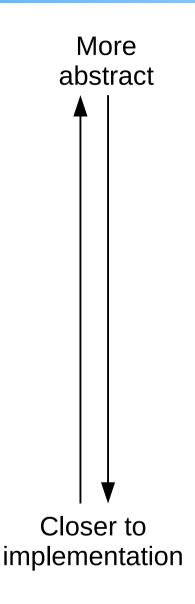
Reduction: multi-BSP

Break into 2 levels



Recap

- PRAM
 - Single shared memory
 - Many processors in lockstep
- BSP
 - Distributed memory, message passing
 - Synchronization with barriers
- Multi-BSP
 - BSP with multiple scales

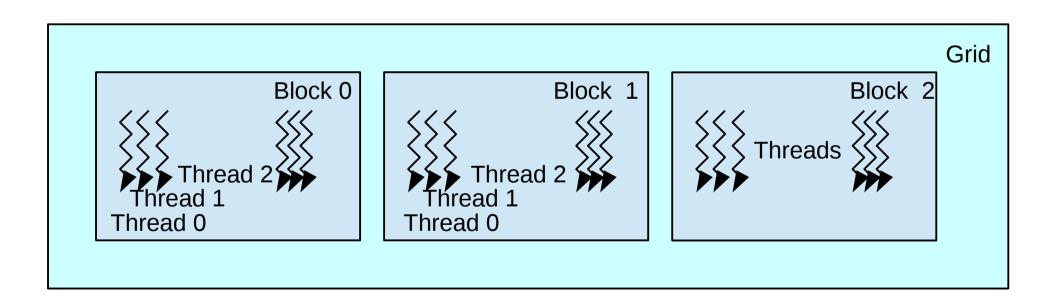


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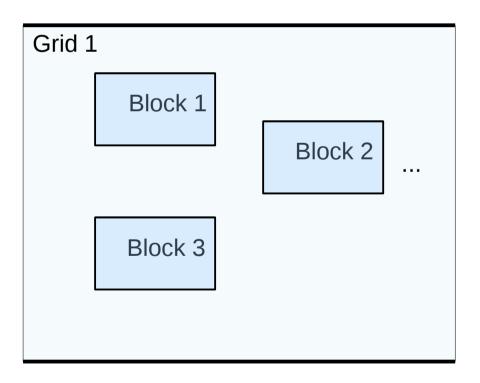
Workload: logical organization

- A kernel is launch on a grid: my_kernel<<<blooks, threads>>>(...)
- Two nested levels
 - Blocks
 - Threads



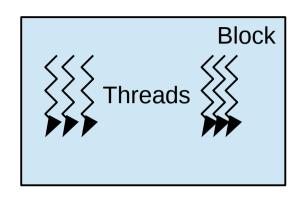
Outer level: grid of blocks

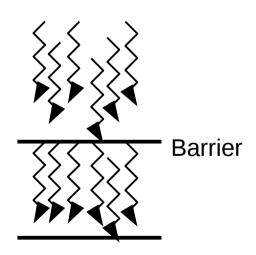
- Blocks also named Concurrent Thread Arrays (CTAs)
- No communication between blocks of the same grid
- Practically unlimited number of blocks / grid



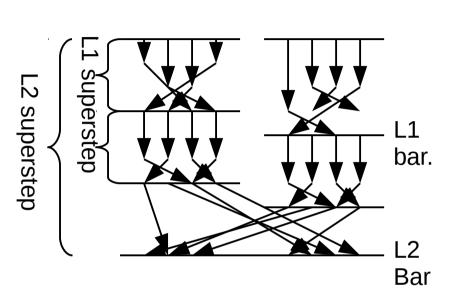
Inner level: threads

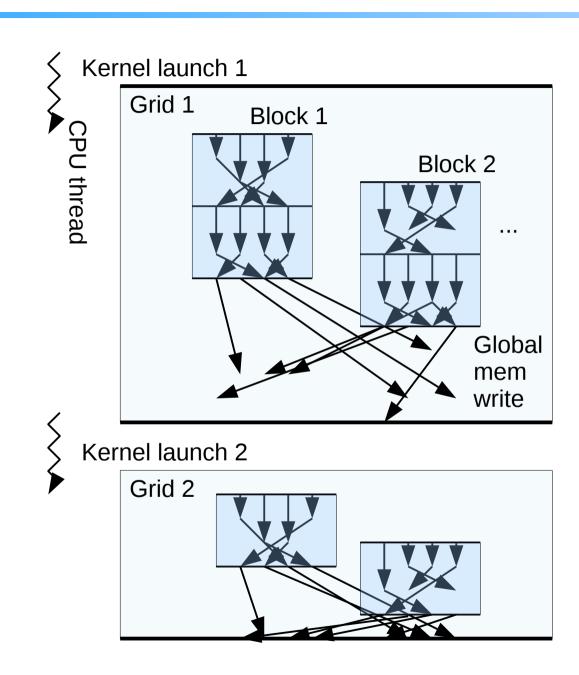
- Blocks contain threads
- All threads in a block
 - Run on the same SM: they can communicate
 - Run in parallel: they can synchronize
- Constraints on number of threads / block
 - Maximum:512 to 1024 depending on arch
 - Recommended: at least 64 threads for good performance
 - Recommended: multiple of the warp size (32)



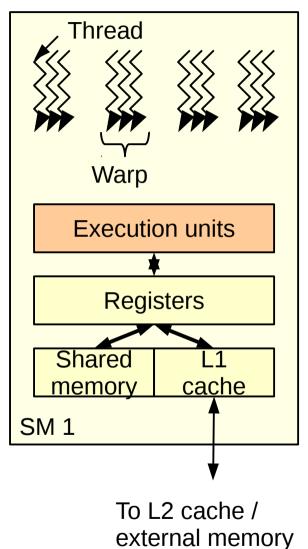


Multi-BSP and CUDA

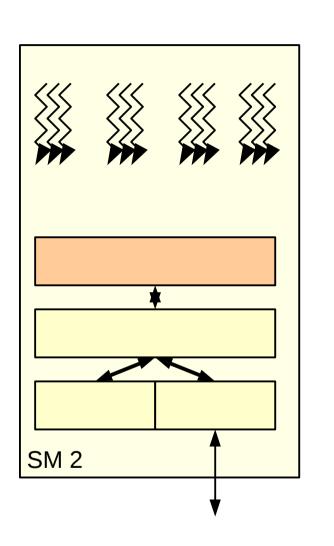




GPU physical organization



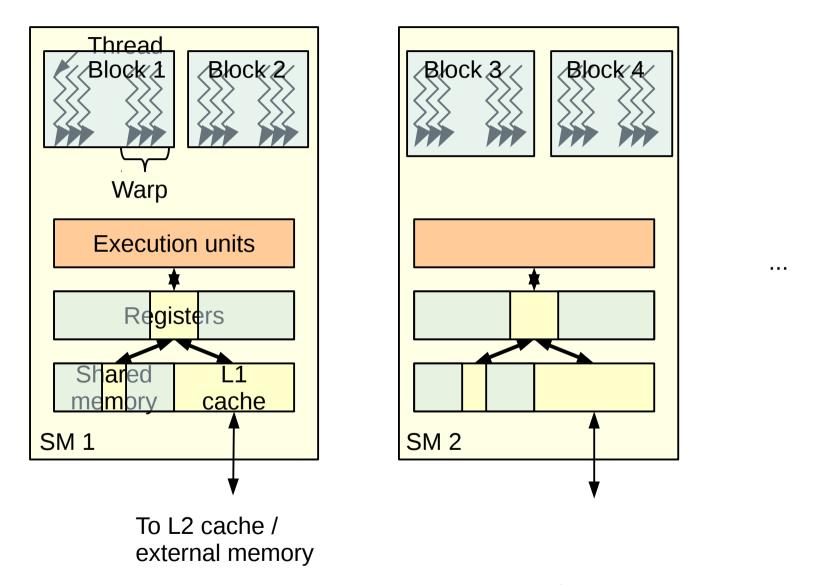
external memory



38

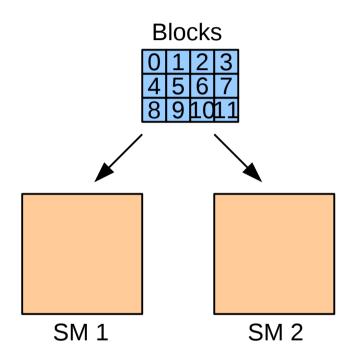
...

Mapping blocks to hardware resources

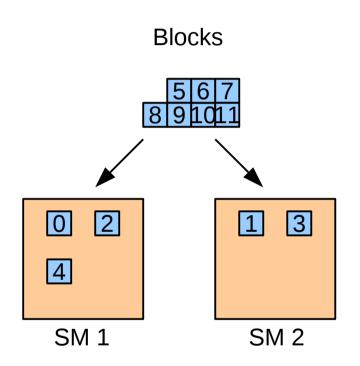


SM resources are partitioned across blocks

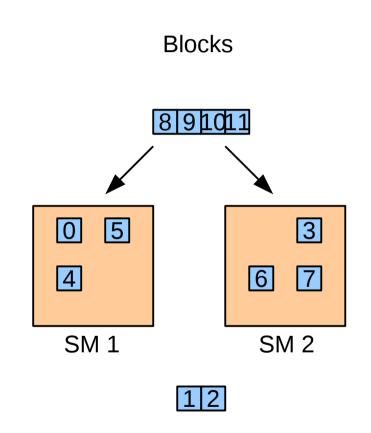
- Blocks may
 - Run serially or in parallel
 - Run on the same or different SM
 - Run in order or out of order
- Should not assume anything on execution order of blocks



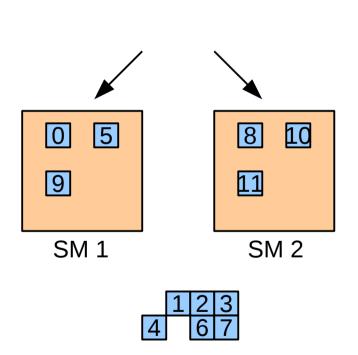
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Blocks

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Example: vector addition

- Addition example: only 1 thread
 - Now let's run a parallel computation
- Start with multiple blocks, 1 thread/block
 - Independent computations in each block
- No communication/synchronization needed

Host code: initialization

A and B are now arrays: just change allocation size

```
int main()
    int numElements = 50000;
    size t size = numElements * sizeof(float);
    float *h A = (float *)malloc(size);
    float *h B = (float *)malloc(size);
    float *h C = (float *)malloc(size);
    Initialize(h A, h B);
   // Allocate device memory
    float *d A, *d B, *d C;
    cudaMalloc((void **)&d_A, size);
    cudaMalloc((void **)&d_B, size);
    cudaMalloc((void **)&d C, size);
    cudaMemcpy(d A, h A, size, cudaMemcpyHostToDevice);
    cudaMemcpy(d_B, h_B, size, cudaMemcpyHostToDevice);
```

Host code: kernel and kernel launch

```
__global__ void vectorAdd2(float *A, float *B, float *C)
{
   int i = blockIdx.x;

   C[i] = A[i] + B[i];
}
```

Launch n blocks of 1 thread each (for now)

```
int blocks = numElements;
vectorAdd2<<<blooks, 1>>>(d_A, d_B, d_C);
```

Device code

- Block number i processes element i
- Grid of blocks may have up to 3 dimensions (blockIdx.x, blockIdx.y, blockIdx.z)
 - For programmer convenience: no effect on scheduling

Multiple blocks, multiple threads/block

Fixed number of threads / block: here 64

Host code

```
Not necessarily multiple of block size!
int threads = 64;
int blocks = (numElements + threads - 1) / threads; // Round up
vectorAdd3<<<blooks, threads>>>(d_A, d_B, d_C, numElements);
```

Device code

```
__global__ void vectorAdd3(const float *A, const float *B, float *C,
    int n)
{
        int i = blockIdx.x * blockDim.x + threadIdx.x;

        if(i < n) {
            C[i] = A[i] + B[i];
        }
}</pre>
Last block may have less work to do

}
```

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Barriers

- Threads can synchronize inside one block
 - Wait until all threads in the block have reached the barrier
- In C for CUDA:

```
__syncthreads();
```

 Needs to be called at the same place for all threads of the block

```
if(tid < 5) {
    ...
}
else {
    ...
}
__syncthreads();</pre>
```

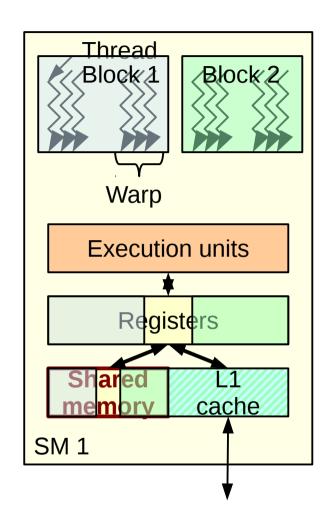
```
if(a[0] == 17) {
   __syncthreads();
}
else {
   __syncthreads();
}
Same condition
for all threads in the block
```

```
if(tid < 5) {
    __syncthreads();
}
else {
    __syncthreads();
}</pre>
```

Correct Correct Wrong

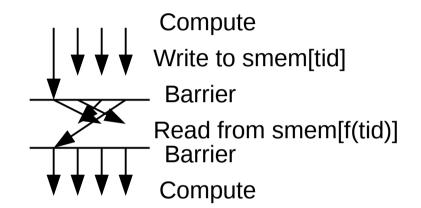
Shared memory

- Fast, software-managed memory
 - Faster than global memory
- Valid only inside one block
 - Each block sees its own copy
- Used to exchange data between threads
- Concurrent writes: one thread wins, but we do not know which one



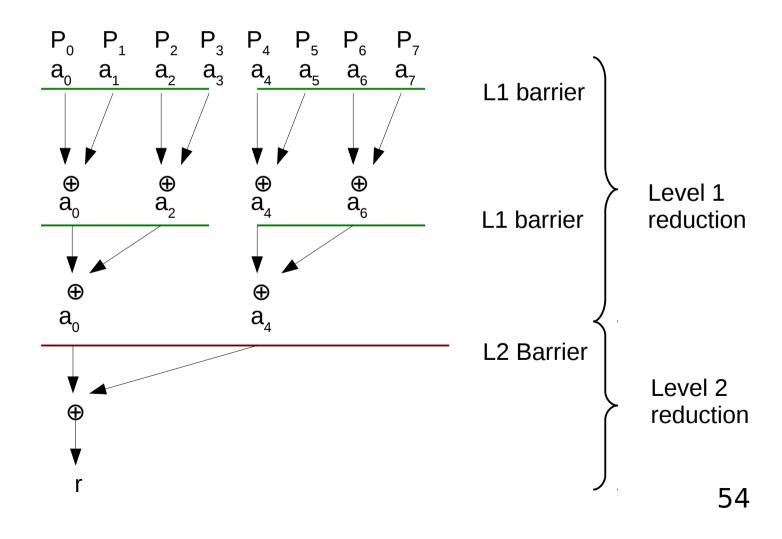
Thread communication: common pattern

- Each thread writes to its own location
 - No write conflict
- Barrier
 - Wait until all threads have written
- Read data from other threads



Example: parallel reduction

Algorithm for 2-level multi-BSP model



Reduction in CUDA: level 1

```
__global__ void reduce1(float *g_idata, float *g_odata, unsigned int n)
{
    extern __shared__ float sdata[];
                                                  Dynamic shared memory allocation:
                                                  will specify size later
    unsigned int tid = threadIdx.x;
    unsigned int i = blockIdx.x * blockDim.x + threadIdx.x;
    // Load from global to shared mem
    sdata[tid] = (i < n) ? g idata[i] : 0;
    __syncthreads();
    for(unsigned int s = 1; s < blockDim.x; s *= 2) {
        int index = 2 * s * tid;
        if(index < blockDim.x) {</pre>
            sdata[index] += sdata[index + s];
        __syncthreads();
    // Write result for this block to global mem
    if (tid == 0) g_odata[blockIdx.x] = sdata[0];
}
```

Reduction: host code

- Level 2: run reduction kernel again, until we have 1 block left
- By the way, is our reduction operator associative?

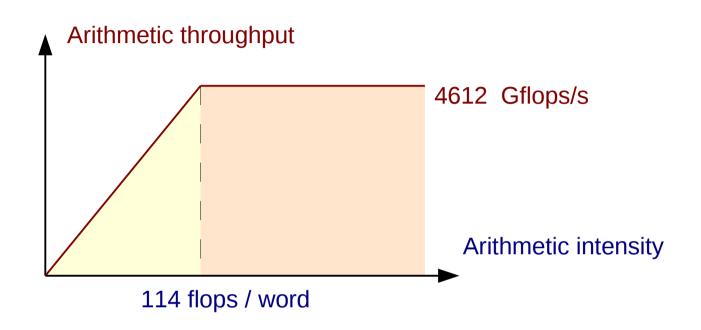
A word on floating-point

- Parallel reduction requires the operator to be associative
- Is addition associative?
 - On reals: yes
 - On floating-point numbers: no
 With 4 decimal digits:
 (1.234 + 123.4) 123.4 = 124.6 123.4 = 1.200
 1.234 + (123.4 123.4) = 1.234 + 0 = 1.234
- Consequence: different result depending on thread count

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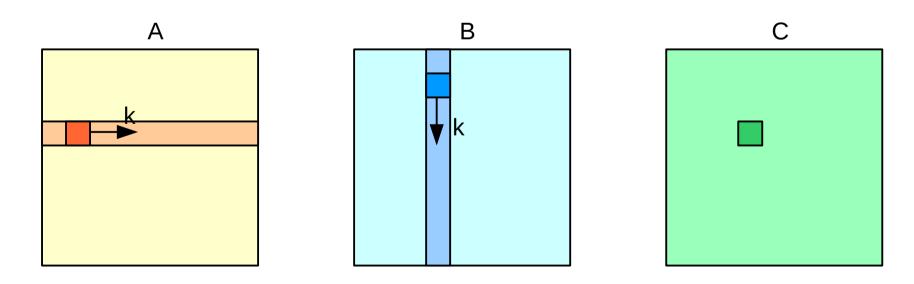
Arithmetic intensity



- Example from first lecture
 - NVIDIA GTX 980 needs ≥114 flops / word to reach peak performance
- How to reach enough arithmetic intensity?
 - Need to reuse values loaded from memory

Classic example: matrix multiplication

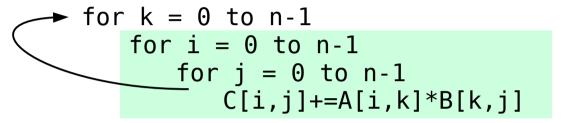
Naive algorithm

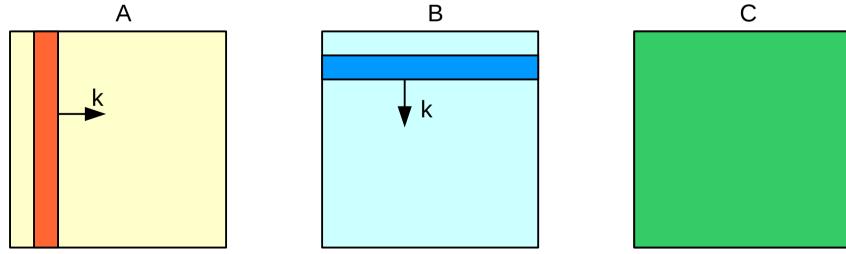


• Arithmetic intensity: 1:1 :(

Reusing inputs

Move loop on k up



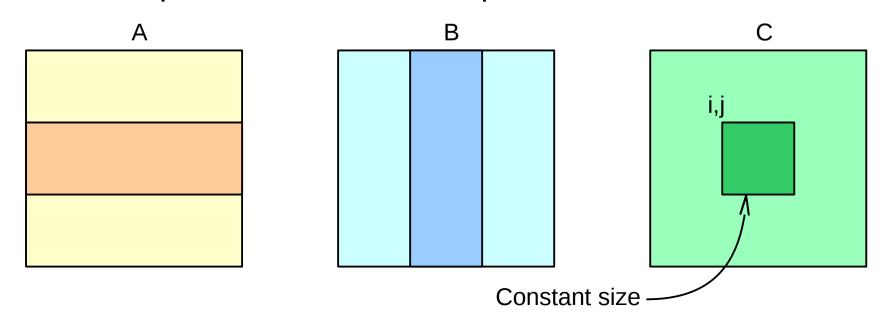


- Enable data reuse on inputs A and B
- But no more reuse on matrix C!

With tiling

Block loops on i and j

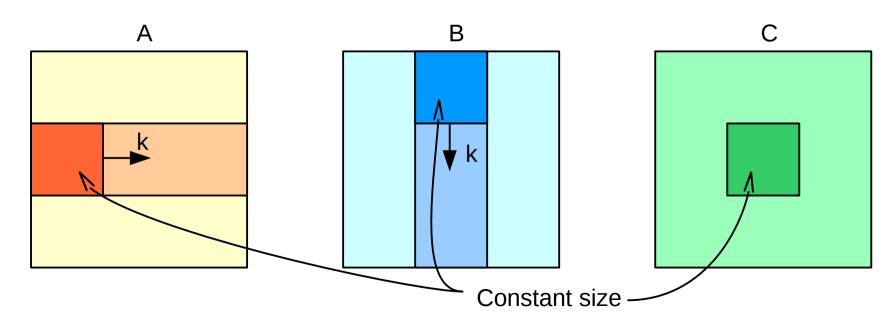
 For one block: product between horizontal panel of A and vertical panel of B



With more tiling

Block loop on k

```
for i = 0 to n-1 step 16
  for j = 0 to n-1 step 16
    for k = 0 to n-1 step 16
    for k2 = k to k+15
        for i2 = i to i+15
        for j2 = j to j+15
        C[i2,j2]+=A[i2,k2*B[k2,j2]
```



Pre-loading data

Arithmetic intensity?

```
for i = 0 to n-1 step 16
   for j = 0 to n-1 step 16
       c = \{0\}
       for k = 0 to n-1 step 16
           a = A[i..i+15,k..k+15]
                                                        Load submatrices a and b
           b = B[k..k+15,j..j+15]
           for k2 = 0 to 15
                                                       Multiply submatrices c = a \times b
              for i2 = 0 to 15
                  for j2 = 0 to 15
                      c[i2,j2]+=a[i2,k2]*b[k2,j2]
       C[i..i+15,j..j+15] = c
                                                     → Store submatrix c
                                   В
          Α
                                   b
    a
                                                            C
```

Breaking into two levels

Run loops on i, j, i2, j2 in parallel

Let's focus on threads

Level 1: SIMD (PRAM-style) version

- Each processor has ID (x,y)
 - Loops on i2, j2 are implicit

```
c[x,y] = 0
for k = 0 to n-1 step 16
a[x,y] = A[i+x,k+y]
b[x,y] = B[k+x,j+y]
for k2 = 0 \text{ to } 15
c[x,y] = a[x,k2]*b[k2,y]
C[i+x,j+y] = c[x,y]
```

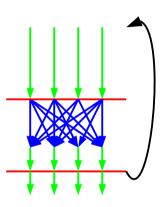
Read from other processors

How to translate to SPMD (BSP-style) ?

SPMD version

Place synchronization barriers

```
c[x,y] = 0
for k = 0 to n-1 step 16
   a[x,y] = A[i+x,k+y]
   b[x,y] = B[k+x,j+y]
   Barrier
   for k2 = 0 to 15
        c[x,y]+=a[x,k2]*b[k2,y]
   Barrier
C[i+x,j+y] = c[x,y]
```



Why do we need the second barrier?

Data allocation

- 3 memory spaces: Global, Shared, Local
 - Where should we put: A, B, C, a, b, c?

```
c[x,y] = 0
for k = 0 to n-1 step 16
  a[x,y] = A[i+x,k+y]
  b[x,y] = B[k+x,j+y]
  Barrier
  for k2 = 0 to 15
     c[x,y] += a[x,k2]*b[k2,y]
  Barrier
C[i+x,j+y] = c[x,y]
```

Data allocation

- Memory spaces: Global, Shared, Local
 - As local as possible

CUDA version

Straightforward translation

```
float Csub = 0;
                                              Precomputed base addresses
for(int a = aBegin, b = bBegin;
     a \le aEnd:
     a += aStep, b += bStep) {
    __shared__ float As[BLOCK_SIZE][BLOCK_SIZE]; \_ Declare shared memory
    __shared__ float Bs[BLOCK_SIZE][BLOCK_SIZE];
    As[ty][tx] = A[a + wA * ty + tx];
                                              Linearized arrays
    Bs[ty][tx] = B[b + wB * ty + tx];
    __syncthreads();
    for(int k = 0; k < BLOCK_SIZE; ++k)</pre>
        Csub += As[ty][k] * Bs[k][tx];
    syncthreads();
int c = wB * BLOCK SIZE * by + BLOCK SIZE * bx;
C[c + wB * ty + tx] = Csub;
```

Local memory

- Registers are fast but
 - Limited in size
 - Not addressable
- Local memory used for
 - Local variables that do not fit in registers (register spilling)
 - Local arrays accessed with indirection

```
int a[17];
b = a[i];
```

- Warning: local is a misnomer!
 - Physically, local memory usually goes off-chip

Device functions

- Kernel can call functions
- Need to be marked for GPU compilation

```
__device__ int foo(int i) {
}
```

A function can be compiled for both host and device

```
__host__ __device__ int bar(int i) {
}
```

- Device functions can call device functions
 - Older GPUs do not support recursion

Recap

- Memory management:
 Host code and memory / Device code and memory
- Writing GPU Kernels
- Dimensions of parallelism: grids, blocks, threads
- Memory spaces: global, local, shared memory

Next time: advanced features and optimization techniques

References and further reading

- CUDA C Programming Guide
- Mark Harris. Introduction to CUDA C. http://developer.nvidia.com/cuda-education
- David Luebke, John Owens. Intro to parallel programming.
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- Paulius Micikevicius. GPU Performance Analysis and Optimization. GTC 2012. http://on-demand.gputechconf.com/gtc/2012/presentations/S05 14-GTC2012-GPU-Performance-Analysis.pdf