# **CME 213**

## Lecture 21: Atomics & Segmented Scan

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## **Overview**

- Atomics
- Segmented Scan

## **ATOMICS**

#### **The Problem**

- How do you do global communication?
- Finish a grid and start a new one
- Scan
- Atomics

- Coordinate by writing to a predefined memory location
  - Race condition! Updates can be lost

- What is the value of a in thread 0?
- What is the value of a in thread 1917?

- Thread 0 could have finished execution before 1917 started
- Or the other way around
- Or both are executing at the same time

Answer: not defined by the programming model, can be arbitrary

#### **Atomics**

CUDA provides atomic operations to deal with this problem

#### **Atomics**

- An atomic operation guarantees that only a single thread has access to a piece of memory while an operation completes
- The name atomic comes from the fact that it is uninterruptable
- No dropped data, but ordering is still arbitrary
- Different types of atomic instructions
- atomic{Add, Sub, Exch, Min, Max,
  Inc, Dec, CAS, And, Or, Xor}
- More types in fermi

#### **Example: Histogram**

```
// Determine frequency of colors in a picture
// colors have already been converted into ints
  Each thread looks at one pixel and increments
// a counter atomically
global void histogram(int* color,
                            int* buckets)
  int i = threadIdx.x
        + blockDim.x * blockIdx.x;
  int c = colors[i];
  atomicAdd(&buckets[c], 1);
```

## **Example: Workqueue**

```
// For algorithms where the amount of work per item
// is highly non-uniform, it often makes sense for
// to continuously grab work from a queue
global
void workq(int* work q, int* q counter,
           int* output, int queue max)
  int i = threadIdx.x
        + blockDim.x * blockIdx.x;
  int q index =
    atomicInc(q counter, queue max);
  int result = do work(work q[q index]);
  output[i] = result;
```

#### **Atomics**

- Atomics are slower than normal load/store
- You can have the whole machine queuing on a single location in memory
- Atomics unavailable on G80!

#### **Example: Global Min/Max (Naive)**

```
// If you require the maximum across all threads
// in a grid, you could do it with a single global
// maximum value, but it will be VERY slow
global
void global max(int* values, int* gl max)
  int i = threadIdx.x
        + blockDim.x * blockIdx.x;
  int val = values[i];
  atomicMax(gl max, val);
```

## **Example: Global Min/Max (Better)**

```
// introduce intermediate maximum results, so that
// most threads do not try to update the global max
global
void global max(int* values, int* gl max,
                 int *reg max,
                 int num regions)
  // i and val as before ...
  int region = i % num regions;
  if(atomicMax(&reg max[region],val) < val)</pre>
  {
    atomicMax(gl max, val);
```

#### Global Min/Max

- Single value causes serial bottleneck
- Create hierarchy of values for more parallelism
- Performance will still be slow, so use judiciously
- See next lecture for even better version!

## **Summary**

- Can't use normal load/store for inter-thread communication because of race conditions
- Use atomic instructions for sparse and/or unpredictable global communication
  - Scan is good for dense communication pattern and where ordering is needed
- Decompose data (very limited use of single global sum/max/min/etc.) for more parallelism

## **SEGMENTED SCAN**

- What it is:
  - Scan + Barriers/Flags associated with certain positions in the input arrays
  - Operations don't propagate beyond barriers
- Do many scans at once, no matter their size

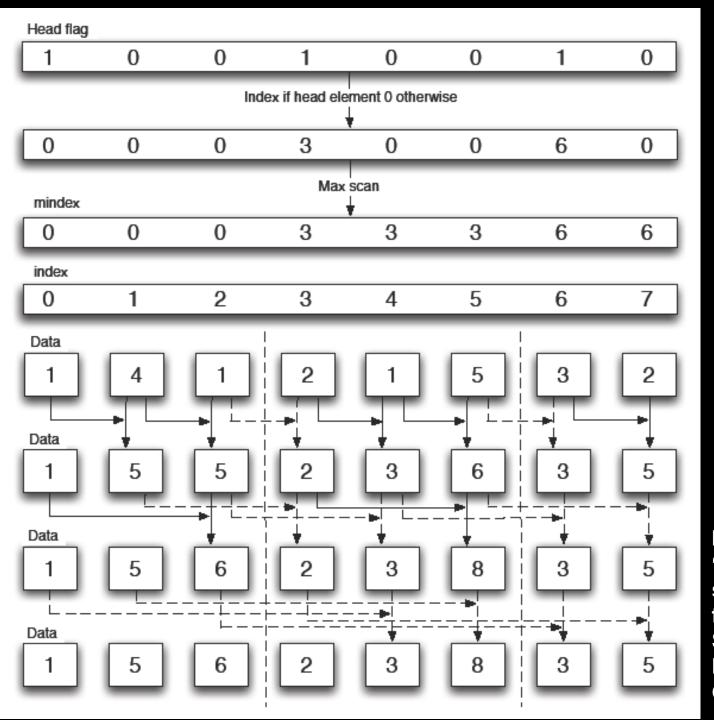


Image taken from "Efficient parallel scan algorithms for GPUs" by S. Sengupta, M. Harris, and M. Garland

```
global void segscan(int * data,
int * flags)
 shared int s data[BL SIZE];
 shared int s flags[BL SIZE];
 int idx = threadIdx.x + blockDim.x
* blockIdx.x;
// copy block of data into shared
// memory
s data[idx] = ...; s flags[idx] = ...;
syncthreads();
```

```
// choose whether to propagate
s data[idx] = s flags[idx] ?
  s data[idx] :
  s data[idx - 1] + s data[idx];
// create merged flag
s flags[idx] =
  s flags[idx - 1] | s flags[idx];
// repeat for different strides
```

Doing lots of reductions of unpredictable size at the same time is the most common use

Think of doing sums/max/count/any over arbitrary sub-domains of your data

- Common Usage Scenarios:
  - Determine which region/tree/group/object class an element belongs to and assign that as its new ID
  - Sort based on that ID
  - Operate on all of the regions/trees/groups/objects in parallel, no matter what their size or number

- Also useful for implementing divide-andconquer type algorithms
  - Quicksort and similar algorithms

## **Questions?**

## **Backup Slides**

#### **Example Segmented Scan**

```
int data[10] = \{1, 1, 1, 1, 1, 1, 1, 1, 1, 1\};
int flags[10] = \{0, 0, 0, 1, 0, 1, 1, 0, 0, 0\};
int step1[10] = \{1, 2, 1, 1, 1, 1, 1, 2, 1, 2\};
int flg1[10] = \{0, 0, 0, 1, 0, 1, 1, 1, 0, 0\};
int step2[10] = \{1, 2, 1, \frac{1}{1}, 1, 1, \frac{1}{2}, \frac{1}{1}, 2\};
int flg2[10] = \{0, 0, 0, 1, 0, 1, 1, 1, 0, 0\};
```

#### **Example Segmented Scan**

```
int step2[10] = {1, 2, 1, 1, 1, 1, 1, 2, 1, 2};
int flg2[10] = {0, 0, 0, 1, 0, 1, 1, 1, 0, 0};
....
```

int result[10] =  $\{1, 2, 3, 1, 2, 1, 1, 2, 3, 4\}$ ;