## **GPU Programming**

\\

Introduction to High Performance Computing, High Performance Data Analytics and Large-Scale Machine Learning



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## **Lecture part**

- Review of GPU architecture
- Review of GPU programming and CUDA
- Some details of our training system, "Cyclone"

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### Covering:

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- Memory coalescing on GPUs
- Shared memory
- What is a warp and should you care

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- Familiarity with some common tasks in C, e.g. array allocation, pointers, etc.
- Can edit text files on a remote server, e.g. text-based (emacs or vim) or VS Code Remote
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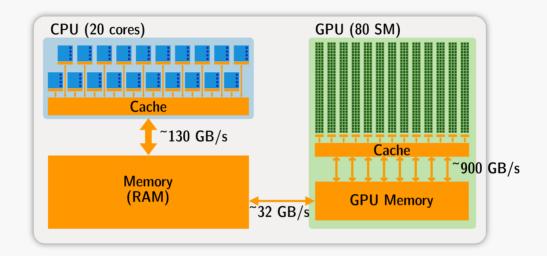
### Repo with slides (PDF) and exercises and URL for direct access to slides:

github.com/CaSToRC-Cyl/EuroCC2 training Nov23

https://slides.koutsou.net/EuroCC-2023-11-09

## **GPU** architecture

### At a very high level:

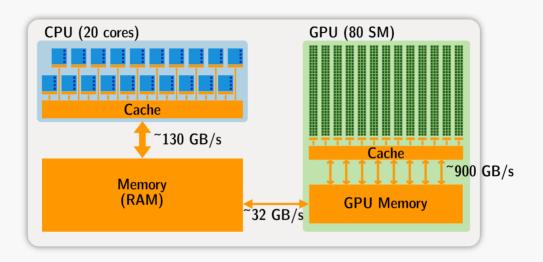


### **CPU**

- Few heavy cores
- Large memory
- Moderate BW to memory
- Optimized for serial execution

### **GPU**

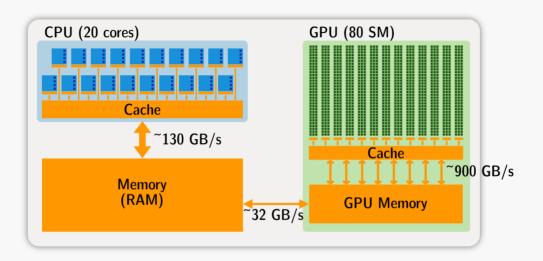
- Many light "cores"
- Smaller memory
- High BW to memory
- Optimized for parallel execution



### Some numbers from the GPU partition of our Cyclone cluster

### **NVIDIA V100** Volta GPUs

- 80 Streaming Multiprocessors (SM) per GPU
- 64 "cores" per SM
- GPU memory: 32 GBytes
- Memory bandwidth: 900~GB/s
- Peak performance: 7.8 Tflop/s (double precision)

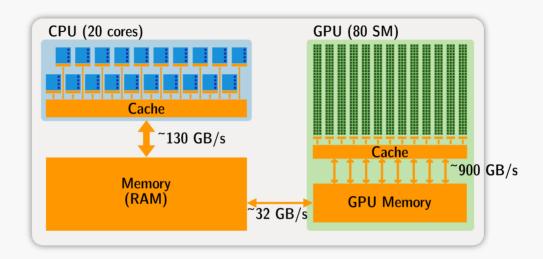


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### NVIDIA V100 Volta GPUs

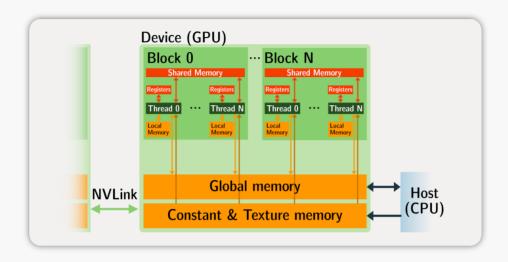
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We will come back to these numbers during the hands-on



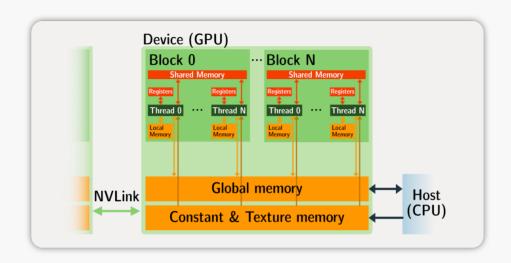
### "Offload" model of programming

- CPU starts program (runs main())
- CPU copies data to GPU memory (over e.g. PCIe, ~32 GB/s)
- CPU dispatches "kernels" for execution on GPU
  - Kernels read/write to GPU memory (~900 GB/s)
  - $\circ$  Kernels run on GPU threads (thousands) which share *fast* memory [O(10) times faster compared to GPU memory]
- Kernel completes; CPU copies data back from GPU (over e.g. PCIe, ~32 GB/s)



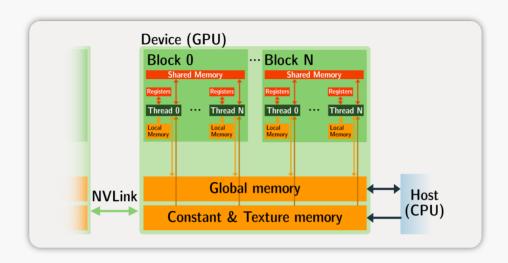
### **GPU** memory model (NVIDIA model)

- GPU threads: slow access to global, constant, and texture memory
- Each thread has registers (fast) and local memory (slow)
- Threads are grouped into blocks; Threads within the same block: shared memory (fast)
- Shared memory is limited. E.g. 96 KB per block for V100



### GPU memory model (NVIDIA model); some numbers for context

- Threads per block: 1024 (max)
- Register memory (per block): 64 KB
- Shared memory (per block): 96 KB
- Also, max. 255 registers per thread



### GPU memory model (NVIDIA model)

- Assumptions about execution order
  - Threads within the same block can be assumed to run concurrently
  - No assumption about the order by which blocks are executed

## **CUDA** programming model

### **NVIDIA** programming framework for NVIDIA GPUs

- Compute Unified Device Architecture
- C-like programming language for writing CUDA Kernels
  - Includes C/C++ and Fortran variants
  - Compiler for C/C++: nvcc
- Functions for transferring data to/from GPUs, starting kernels, etc.
- Some higher-level functionality also available (linear algebra, random number generations, etc.)
- Concepts generalizable to other accelerator programming frameworks (OpenCL, OpenACC, HiP, etc.)

#### **Nomenclature**

- "Host" is the CPU
- "Device" is the GPU

### Allocate memory on GPU

### err = cudaMalloc(&d\_ptr, size);

- Call from host (CPU)
- Allocate size bytes of memory on GPU and store the starting address in d\_ptr
- d\_ptr is a variable that holds an address to GPU memory i.e. a "device pointer"
- If err != cudaSuccess then something went wrong

### Free GPU memory

cudaFree(d\_ptr);

#### **Nomenclature**

- "Host" is the CPU
- "Device" is the GPU

### Copy data to GPU

#### cudaMemcpy(d\_ptr, ptr, size, cudaMemcpyHostToDevice);

- Call from host (CPU)
- Copy data on host pointed to by ptr to device at address pointed to by d\_ptr
- Device memory should have been allocated using cudaMalloc() to obtain d\_ptr

## Copy data from GPU

### cudaMemcpy(ptr, d\_ptr, size, cudaMemcpyDeviceToHost);

- Call from host (CPU)
- Copy data on device pointed to by d\_ptr to host at address pointed to by ptr
- Host memory should have been allocated using e.g. malloc() to obtain ptr

#### Declare a CUDA kernel

Example:

```
global__ void
func(int n, double a, double *x)
{
    return;
}
```

### Call a CUDA kernel

• Call from host. Example:

```
func<<<nblck, nthr>>>(n, a, x);
```

- nthr: number of threads per block; can be scalar or a dim3 type
- nblck: number of blocks; can be scalar or a dim3 type
- Example of dim3 type:

```
dim3 nthr(1024, 8, 8); /* No. of threads in (x, y, z) */
```

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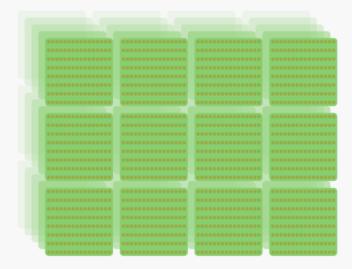
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#### Thread coordinates within kernel

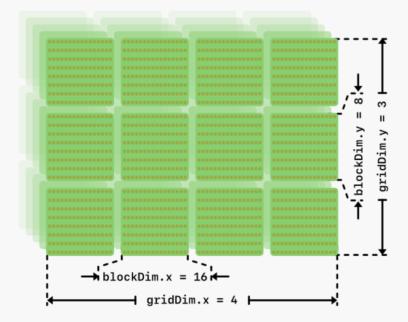
## Example:

```
global__ void
func(int n, double a, double *x)
{
   int idx = threadIdx.x + blockIdx.x*blockDim.x;
   return;
}
```

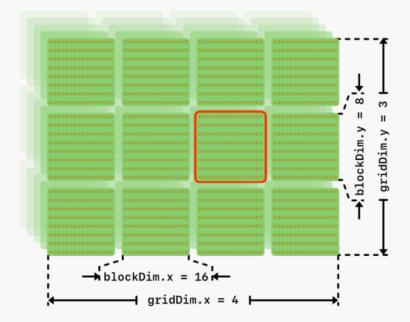
```
dim3 blcks( 4, 3, bz);
dim3 thrds(16, 8, tz);
func<<<blcks, thrds>>>(...);
```



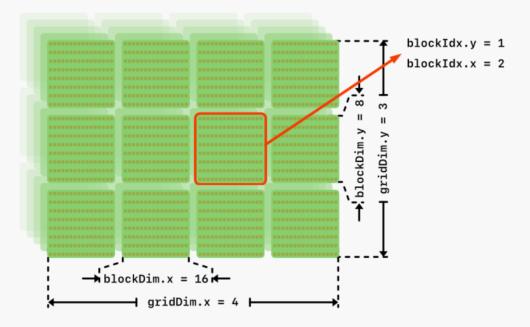
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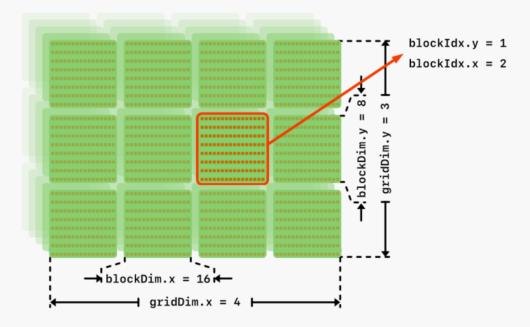
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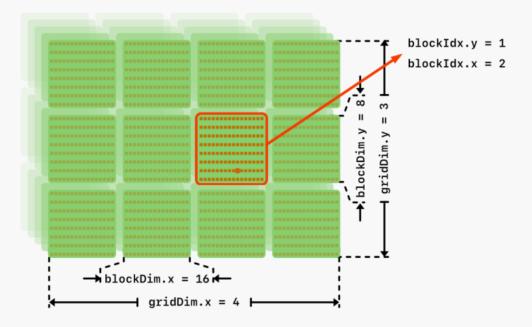
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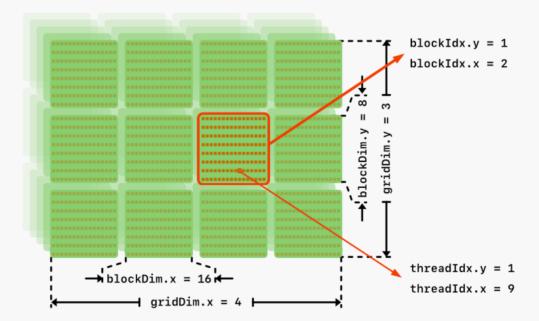
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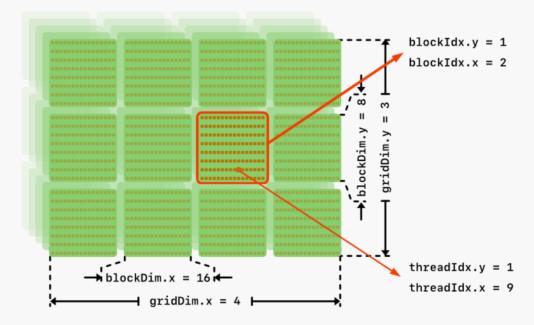


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Threads, blocks, grids

```
dim3 blcks( 4, 3, bz);
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func<<<blcks, thrds>>>(...);
```



### Variables available within kernel

- threadIdx.{x,y,z}
- blockIdx.{x,y,z}

- blockDim.{x,y,z}
- gridDim.{x,y,z}

Port a simple code to GPU and investigate performance

Sources: github.com/CaSToRC-CyI/EuroCC2 training Nov23/GPU programming/exercise/

• axpy.cu implements a so-called "axpy" operation (a-times-x-plus-y):

$$y_i \leftarrow a \cdot x_i + y_i, i = 0, \dots, n-1$$

with  $\alpha$  scalar and y and x vectors of length n.

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#### This will cover:

- Allocation of memory on the GPU;
- Transferring memory to/from GPU;
- Invoking kernels;
- Placement of threads and memory access

### File: axpy.cu

- Contains the C program we will begin with: axpy.cu
- Even though the file extension is .cu, the program contains no CUDA. Only OpenMP
- Allocates four arrays: x0[n], x1[n], y0[n], and y1[n], with n read from the command line
- x0 and y0 are initialized to random numbers
- x1 and y1 are initialized to x0 and y0 respectively
- The program:
  - o performs y0[:] = a\*x0[:] + y0[:] in the first part marked with A:
  - o performs y1[:] = a\*x1[:] + y1[:] in the second part marked with B:
  - reports the timing for part A and for B
  - o reports the difference between yo and y1

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Take some time to inspect axpy.cu before we compile and run

• Copy first exercise from this training's shared space:

```
[user@front02 ~]$ git clone https://github.com/CaSToRC-CyI/EuroCC2_training_Nov23.git
[user@front02 ~]$ cd EuroCC2_training_Nov23/GPU_programming/exercise
[user@front02 ]$ ls -1
axpy.cu
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Compile with nvcc including OpenMP:

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- Run on the CPUs of a GPU node
- Use srun to run interactively, e.g.:

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[user@front02 exercise] export OMP_NUM_THREADS=20
[user@front02 exercise] srun -n 1 --cpus-per-task=20 -p gpu -A da23 --gres=gpu:1 --reservation=data-analytics
CPU: nthr = 20  t0 = 0.0089 sec  P = 15.024 Gflop/s  B = 90.142 GB/s
CPU: nthr = 20  t0 = 0.0086 sec  P = 15.667 Gflop/s  B = 94.000 GB/s
Diff = 0.000000e+00
```

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Diff = 0.0000000e+00
```

Compare ~90 GB/s achieved vs ~130 GB/s peak memory bandwidth

#### Use a GPU to replace part B of the calculation

- Edits outside of main():
  - 1. Add the cuda\_runtime.h header file
  - 2. Add the GPU axpy kernel, naming it gpu\_axpy()
  - 3. Add a function similar to ualloc() that allocates memory on the GPU and checks whether an error occured
- Edits within main():
  - 1. Allocate arrays on GPU
  - 2. Copy x1[:] and y1[:] to GPU
  - 3. Call gpu\_axpy()
  - 4. Copy y1[:] from GPU

Edits outside of main() 1/3

• Add the cuda\_runtime.h header file on line 5:

#include <cuda\_runtime.h>

#### Edits outside of main() 2/3

• Add the GPU axpy kernel, naming it gpu\_axpy(), after the CPU axpy, around line 64:

```
/***
 * Do y <- a*x + y on the GPU
 ***/
   _global__ void
gpu_axpy(int n, float a, float *x, float *y)
{
   for(int i=0; i<n; i++)
    y[i] = a*x[i] + y[i];
   return;
}</pre>
```

#### Edits outside of main() 3/3

• At around line 30 add a function similar to ualloc() that allocates memory on the GPU and checks whether an error occurred

```
/***
 * Allocate memory on GPU; print error if not successful
    ***/
void *
gpu_alloc(size_t size)
{
    void *ptr;
    cudaError_t err = cudaMalloc(&ptr, size);
    if(err != cudaSuccess) {
        fprintf(stderr, "cudaMalloc() returned %d; quitting...\n", err);
        exit(-2);
    }
    return ptr;
}
```

#### Edits within main() 1/4

• Allocate arrays on GPU, within B part. Free arrays before closing B part:

```
/*
 * B: Run axpy(), return to y1, report performance
 */
 {
    /* Allocate GPU memory */
    float *d_x = (float *)gpu_alloc(n*sizeof(float));
    float *d_y = (float *)gpu_alloc(n*sizeof(float));
    ...
    cudaFree(d_x);
    cudaFree(d_y);
}
```

#### Edits within main() 2/4

• Copy x1[:] and y1[:] to GPU

```
cudaMemcpy(d_x, x1, sizeof(float)*n, cudaMemcpyHostToDevice);
cudaMemcpy(d_y, y1, sizeof(float)*n, cudaMemcpyHostToDevice);
```

#### Edits within main() 3/4

• Call gpu\_axpy(). For the moment use 1 thread and 1 block. Replace axpy(n, a, x, y) of part B with:

```
double t0 = stop_watch(0);
gpu_axpy<<<1, 1>>>(n, a, d_x, d_y);
t0 = stop_watch(t0);
```

Note we need to pass the *device pointers* since it is these pointers that point to the memory allocated on the GPU

#### Edits within main() 4/4

• Copy y1[:] from GPU:

```
/* Copy y1 back from GPU */
cudaMemcpy(y1, d_y, sizeof(float)*n, cudaMemcpyDeviceToHost);
```

• Also change:

```
printf(" CPU: nthr = %4d ...);
```

to:

```
printf(" GPU: ...);
```

and remove OpenMP parallel region.

#### Compile and run

• Compile as before:

```
[user@front02 exercise]$ nvcc -arch sm_70 -03 -Xcompiler -fopenmp -o axpy axpy.cu
```

• Run as before (I'm assuming OMP\_BIND, OMP\_PLACES, and OMP\_NUM\_THREADS were set before):

```
[user@front02 exercise]$ srun -n 1 --cpus-per-task=20 -p gpu -A da23 --gres=gpu:1 --reservation=data-analytic
CPU: nthr = 20    t0 = 0.0089 sec    P = 15.035 Gflop/s    B = 90.212 GB/s
GPU:         t0 = 0.0000 sec    P = 3198.579 Gflop/s    B = 19191.476 GB/s
Diff = 1.021564e-15
```

This performance is infeasible. What's going on?

#### Edits within main() 3/4

• The problem is here:

```
double t0 = stop_watch(0);
gpu_axpy<<<1, 1>>>(n, a, d_x, d_y);
t0 = stop_watch(t0);
```

- CUDA kernels return **immediately**; the kernel is still being executed on the device when stop\_watch(t0) is called. We are **not** timing the kernel execution time, but the time it takes to dispatch the kernel to the GPU.
- Correct this by adding cudaDeviceSynchronize(); after the CUDA kernel, which blocks until all running CUDA kernels are complete:

```
double t0 = stop_watch(0);
gpu_axpy<<<1, 1>>>(n, a, d_x, d_y);
cudaDeviceSynchronize();
t0 = stop_watch(t0);
```

• Compile and run again:

```
[user@front02 exercise]$ nvcc -arch sm_70 -03 -Xcompiler -fopenmp -o axpy axpy.cu [user@front02 exercise]$ srun -n 1 --cpus-per-task=20 -p gpu -A da23 --gres=gpu:1 --reservation=data-analytic CPU: nthr = 20 t0 = 0.0088 sec P = 15.199 Gflop/s B = 91.193 GB/s GPU: t0 = 3.9670 sec P = 0.034 Gflop/s B = 0.203 GB/s Diff = 1.021564e-15
```

• Compile and run again:

```
[user@front02 exercise]$ nvcc -arch sm_70 -03 -Xcompiler -fopenmp -o axpy axpy.cu
[user@front02 exercise]$ srun -n 1 --cpus-per-task=20 -p gpu -A da23 --gres=gpu:1 --reservation=data-analytic
CPU: nthr = 20  t0 = 0.0088 sec  P = 15.199 Gflop/s  B = 91.193 GB/s
GPU:        t0 = 3.9670 sec  P = 0.034 Gflop/s  B = 0.203 GB/s
Diff = 1.021564e-15
```

• This performance is of course extremely poor;

• Compile and run again:

```
[user@front02 exercise]$ nvcc -arch sm_70 -03 -Xcompiler -fopenmp -o axpy axpy.cu [user@front02 exercise]$ srun -n 1 --cpus-per-task=20 -p gpu -A da23 --gres=gpu:1 --reservation=data-analytic CPU: nthr = 20 t0 = 0.0088 sec P = 15.199 Gflop/s B = 91.193 GB/s GPU: t0 = 3.9670 sec P = 0.034 Gflop/s B = 0.203 GB/s Diff = 1.021564e-15
```

- This performance is of course extremely poor;
- We're using only one GPU thread in the kernel

#### Use more threads

• In this step, we will use 512 GPU threads. First, change the call to the GPU kernel:

```
double t0 = stop_watch(0);
gpu_axpy<<<1, 512>>>(n, a, d_x, d_y);
cudaDeviceSynchronize();
t0 = stop_watch(t0);
```

#### Use more threads

• In this step, we will use 512 GPU threads. First, change the call to the GPU kernel:

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double t0 = stop_watch(0);
gpu_axpy<<<1, 512>>>(n, a, d_x, d_y);
cudaDeviceSynchronize();
t0 = stop_watch(t0);
```

• Then we need to change the kernel. We need in each GPU thread to calculate which elements it will operate on:

```
/***
 * Do y <- a*x + y on the GPU
 ***/
   _global__ void
gpu_axpy(int n, float a, float *x, float *y)
{
   int ithr = threadIdx.x;
   int nthr = blockDim.x;
   int lt = n/nthr;
   for(int i=ithr*lt; i<(ithr+1)*lt; i++)
     y[i] = a*x[i] + y[i];
   return;
}</pre>
```

• With the above, each thread operated on n/nthr contiguous elements

• Compile and run again:

```
[user@front02 exercise]$ srun -n 1 --cpus-per-task=20 -p gpu -A da23 --gres=gpu:1 --reservation=data-analytic CPU: nthr = 8 t0 = 0.0064 sec P = 2.628 Gflop/s B = 15.765 GB/s GPU: t0 = 0.1316 sec P = 0.127 Gflop/s B = 0.765 GB/s Diff = 1.022961e-15
```

• Better than before, but still very poor performance. Can we do better?

#### **Optimized GPU memory access**

Always keep in mind that on GPUs, it is more optimal if contiguous threads access contiguous memory locations

This represents the order by which elements are accessed currently

- The same thread accesses continuous elements
- Very common approach on CPUs
- On GPUs, this results in so-called bank conflicts
- Suboptimal!

#### **Optimized GPU memory access**

Always keep in mind that on GPUs, it is more optimal if contiguous threads access contiguous memory locations

This represents an optimal data access pattern

- Different threads accesses continuous elements
- Each thread is served by a different memory bank

#### **Optimized GPU memory access**

Always keep in mind that on GPUs, it is more optimal if contiguous threads access contiguous memory locations

In our example:

```
/***
 * Do y <- a*x + y on the GPU
 ***/
    global__ void
gpu_axpy(int n, float a, float *x, float *y)
{
    int ithr = threadIdx.x;
    int nthr = blockDim.x;
    for(int i=0; i<n; i+=nthr)
        y[i+ithr] = a*x[i+ithr] + y[i+ithr];
    return;
}</pre>
```

• Compile and run:

```
[user@front02 exercise]\$ srun -n 1 --cpus-per-task=20 -p gpu -A da23 --gres=gpu:1 --reservation=data-analytic CPU: nthr = 20 t0 = 0.0097 sec P = 13.788 Gflop/s B = 82.730 GB/s GPU: t0 = 0.0665 sec P = 2.018 Gflop/s B = 12.111 GB/s Diff = 1.021564e-15
```

#### Blocks and threads

Now let's use blocks. Let's use as many blocks and threads as we can

- Upper limit of 1024 threads
- Upper limit of  $2^{31} 1$  blocks

```
double t0 = stop_watch(0);
int nthr = 1024;
gpu_axpy<<<n/nthr, nthr>>>(n, a, d_x, d_y);
cudaDeviceSynchronize();
t0 = stop_watch(t0);
```

```
/***
 * Do y <- a*x + y on the GPU
 ***/
    global__ void
gpu_axpy(int n, float a, float *x, float *y)
{
    int ithr = threadIdx.x;
    int nthr = blockDim.x;
    int iblk = blockIdx.x;
    int idx = ithr + iblk*nthr;
    y[idx] = a*x[idx] + y[idx];
    return;
}</pre>
```

#### **Blocks and threads**

• Compile and run:

```
[user@front02 exercise]\$ srun -n 1 --cpus-per-task=20 -p gpu -A da23 --gres=gpu:1 --reservation=data-analytic CPU: nthr = 20 t0 = 0.0088 sec P = 15.188 Gflop/s B = 91.129 GB/s GPU: t0 = 0.0011 sec P = 119.930 Gflop/s B = 719.578 GB/s Diff = 1.021564e-15
```

• ~720 GB/s is ~80% of peak bandwidth (which is 900 GB/s)

#### **Blocks and threads**

• Compile and run:

```
[user@front02 exercise]$ srun -n 1 --cpus-per-task=20 -p gpu -A da23 --gres=gpu:1 --reservation=data-analytic

CPU: nthr = 20 t0 = 0.0088 sec P = 15.188 Gflop/s B = 91.129 GB/s

GPU: t0 = 0.0011 sec P = 119.930 Gflop/s B = 719.578 GB/s

Diff = 1.021564e-15
```

- ~720 GB/s is ~80% of peak bandwidth (which is 900 GB/s)
- Try varying the number of threads per block. E.g. with 512 threads I got ~730 GB/s.