# Evaluation task for Awkward Array GSoC project

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#### Abstract

In this report, we demonstrate a sample GPU kernel designed to be an alternative to the CPU backend for array operations in the Awkward Array project. In particular, we implement a CUDA translation of the <code>awkward\_listarray\_compact\_offsets</code> CPU kernel, and show how parallelizing the code on a GPU could significantly increase the speed of computation. The motivation of this report is to serve as an evaluation to enable me to qualify to work on the project titled "Awkward Array GPU kernels" under the mentorship of Jim Pivarski and David Lange.

#### 1 Introduction

The sample CPU kernel (directly taken from the Awkward Array codebase) to be translated is defined below -

```
template <typename C, typename T>
ERROR awkward_listarray_compact_offsets(T* tooffsets, const C* fromstarts, const C
    * fromstops, int64_t startsoffset, int64_t stopsoffset, int64_t length) {
    tooffsets[0] = 0;
    for (int64_t i = 0; i < length; i++) {
        C start = fromstarts[startsoffset + i];
        C stop = fromstops[stopsoffset + i];
        if (stop < start) {
            return failure("stops[i] < starts[i]", i, kSliceNone);
        }
        tooffsets[i + 1] = tooffsets[i] + (stop - start);
    }
    return success();
}</pre>
```

Translating this particular CPU kernel serves as a good test because parallelizing this CPU kernel involves overcoming the loop carried dependency in the above code -

```
tooffsets[i + 1] = tooffsets[i] + (stop - start);
```

where tooffsets[i + 1] depends on toffsets[i].

The above algorithm can be visualized to illustrate its linear nature -



Figure 1: Visualizing sequential CPU algorithm

#### 1.1 Implementation of GPU algorithm

The evaluation task required creating a parallel algorithm as an alternative for the sequential CPU algorithm. There are several ways to do this -

- 1. OpenCL [1], which is a more uniform interface to write parallel algorithms to be executed on multiple backends, but is slower than CUDA and losing popularity.
- 2. CUDA [2], which is an easier and more performant interface for executing parallel algorithms only on an NVIDIA GPU.

Due to the availability of an NVIDIA GPU and performance requirements, we have decided to use CUDA to write and execute our parallel algorithms.

We use some CUDA intrinsics like thrust [3] in our sample code for the evaluation task which we might not use in the production code to avoid locking our parallel implementation to CUDA and allow expansion to newer interfaces in the future.

### 2 Sequential algorithm on an Nvidia GPU

#### 2.1 Sequential algorithm on a single thread

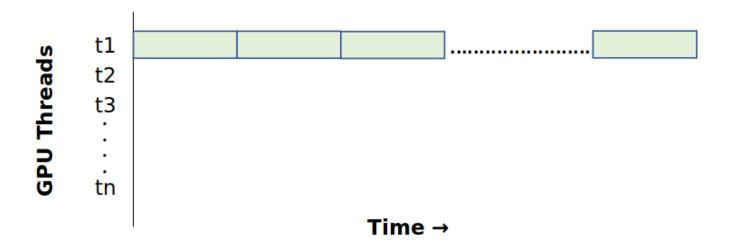


Figure 2: Visualizing GPU threads at work

```
device
void success_cuda(Error* err) {
  err.str = nullptr;
  err.identity = kSliceNone;
  err.attempt = kSliceNone;
  err.extra = 0;
}
void failure_cuda(const char* str, int64_t identity, int64_t attempt, Error* err)
  err.str = str;
  err.identity = identity;
  err.attempt = attempt;
  err.extra = 0;
}
template <typename C, typename T>
void awkward_listarray_compact_offsets(T* tooffsets, const C* fromstarts, const C*
    from stops, int64_t startsoffset, int64_t stopsoffset, int64_t length, Error*
```

```
err) {
    shared
            int flag [1];
  int idx = threadIdx.x + (blockIdx.x * blockDim.x);
  if (idx = 0) {
    tooffsets[0] = 0;
    flag[0] = 0;
  if (idx < length) {
    if (idx == 0) {
      for (int i = 0; i < length; i++) {
        C start = fromstarts[startsoffset + i];
        C stop = from stops [stopsoffset + i];
        if (stop < start) {
          failure\_cuda("stops[i] < starts[i]", i, kSliceNone, err);
          flag[0] = 1;
        tooffsets[i + 1] = tooffsets[i] + (stop - start);
    }
  }
    _{\rm syncthreads}();
  if (flag[0] != 1) {
    success_cuda(err);
}
```

In the above code block<sup>1</sup> we execute the algorithm only for a single GPU thread (in a for loop) while the other threads sit idle.

```
if (idx == 0) {
...
}
```

where the variable idx is the thread index.

Since CUDA GPU kernels have to be of the return type void, we use additional device functions, success\_cuda and failure\_cuda to record whether the kernel behaved as expected.

```
__device__
void success_cuda(Error* err) {
  err.str = nullptr;
  err.identity = kSliceNone;
  err.attempt = kSliceNone;
```

<sup>&</sup>lt;sup>1</sup>Full benchmarking code: https://github.com/reikdas/GSoC-Proposal-2020/blob/master/testgsoc/naivesingle.cu

We use a flag which has been implemented using shared memory and the CUDA intrinsic thread synchronization

```
__shared__ int flag[1];
...
__syncthreads();
if (flag[0] != 1) {
...
}
```

to determine whether the success\_cuda() call should be executed once the failure\_cuda() call has already been reached by one of the GPU threads.

### 2.2 Sequential algorithm on multiple threads

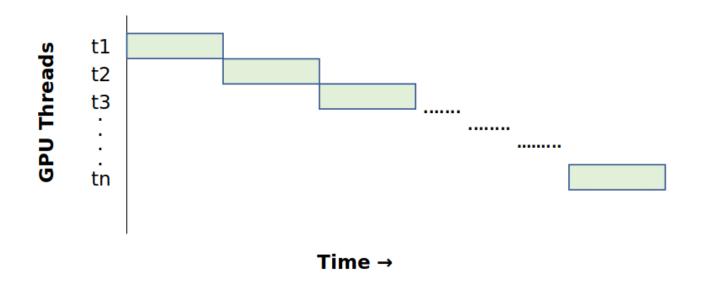


Figure 3: Visualizing GPU threads at work

```
_{
m device}
void success_cuda(Error* err) {
  err.str = nullptr;
  err.identity = kSliceNone;
  err.attempt = kSliceNone;
  err.extra = 0;
}
  _{
m device}
void failure_cuda(const char* str, int64_t identity, int64_t attempt, Error* err)
  err.str = str;
  err.identity = identity;
  err.attempt = attempt;
  err.extra = 0;
}
template <typename C, typename T>
___global___
```

```
void awkward_listarray_compact_offsets (T* tooffsets, const C* fromstarts, const C*
    from stops, int64_t startsoffset, int64_t stopsoffset, int64_t length, Error*
   err) {
    \underline{\text{shared}} int flag[1];
  int idx = threadIdx.x + (blockIdx.x * blockDim.x);
  if (idx < length) {
    if (idx == 0) {
      tooffsets[0] = 0;
      flag[0] = 0;
    for (int i = 0; i < length; i++) {
        _syncthreads();
      if (i = idx) 
        C start = fromstarts[startsoffset + i];
        C stop = from stops [stopsoffset + i];
        if (stop < start) {
          failure_cuda("stops[i] < starts[i]", i, kSliceNone, err);
          flag[0] = 1;
        tooffsets[i + 1] = tooffsets[i] + (stop - start);
    }
    syncthreads();
  if (flag[0] != 1) {
    success_cuda(err);
}
```

The sequential algorithm on multiple GPU threads implementation shown above<sup>2</sup> differs only slightly from the sequential algorithm on a single GPU thread implementation.

The main difference is that the multiple GPU threads implementation has all the threads execute the algorithm one after another (no loops), instead of all at the same time. When one thread is being executed, the other threads sit idly, then the next thread is executed while the previous and all other threads sit idle.

```
for (int i = 0; i < length; i++) {
    ...
    if (i == idx) {
        ...
    }
}</pre>
```

We use the CUDA intrinsic \_\_syncthreads() before each iteration of the loop to ensure that all the threads are at the same position and ready to be executed before each thread is executed.

<sup>&</sup>lt;sup>2</sup>Full benchmarking code: https://github.com/reikdas/GSoC-Proposal-2020/blob/master/testgsoc/naivemultiblocks.cu

## 3 Parallel algorithm on an Nvidia GPU

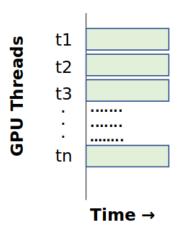


Figure 4: Visualizing GPU threads at work

```
#include <thrust/device_vector.h>
#include <thrust/scan.h>
#include "helper_cuda.h"
  _{
m device}
void success_cuda(Error* err) {
  err.str = nullptr;
  err.identity = kSliceNone;
  err.attempt = kSliceNone;
  err.extra = 0;
}
void failure_cuda(const char* str, int64_t identity, int64_t attempt, Error* err)
   {
  err.str = str;
  err.identity = identity;
  err.attempt = attempt;
  err.extra = 0;
}
template <typename T, typename C>
  _global__
```

```
void sub(T* output, const C* starter, const C* stopper, int64_t startsoffset,
   int64_t stopsoffset, int64_t n, Error* err) {
    \_shared\_\_ int flag[1];
  int thid = threadIdx.x + blockIdx.x * blockDim.x;
  if (thid == 0) flag [0] = 0;
  if (thid < n) {
    C start = starter[thid + startsoffset];
    C stop = stopper[thid + stopsoffset];
    if (stop < start) {
      failure_cuda("stops[i] < starts[i]", thid , kSliceNone, err);
      flag[0] = 1;
    output[thid] = stop - start;
   _syncthreads();
  if (flag[0] != 1) {
    success_cuda(err);
}
template <typename T, typename C>
void prefix sum (T* output, const C* arr, const C* arr2, int64 t startsoffset,
   int64_t stopsoffset, int64_t length, Error* err) {
  int block, thread;
  if (length > 1024) {
    block = (length / 1024) + 1;
    thread = 1024;
  }
  else {
    thread = length;
    block = 1;
 T* d_output;
 C* d arr, * d arr2;
  checkCudaErrors(cudaMalloc((void**)&d_output, length * sizeof(T)));
  checkCudaErrors(cudaMalloc((void**)&d_arr, length * sizeof(C)));
  checkCudaErrors(cudaMemcpy(d_arr, arr, length * sizeof(C),
     cudaMemcpyHostToDevice));
  checkCudaErrors(cudaMalloc((void**)&d_arr2, length * sizeof(C)));
  checkCudaErrors(cudaMemcpy(d_arr2, arr2, length * sizeof(C),
     cudaMemcpyHostToDevice));
  sub<T, C><<block, thread>>>(d_output, d_arr, d_arr2, startsoffset, stopsoffset,
      length, err);
  checkCudaErrors(cudaDeviceSynchronize());
  thrust::device vector <T> data(d output, d output+length);
```

```
thrust::device_vector<T> temp(data.size() + 1);
thrust::exclusive_scan(data.begin(), data.end(), temp.begin());
temp[data.size()] = data.back() + temp[data.size() - 1];
thrust::copy(temp.begin(), temp.end(), output);
checkCudaErrors(cudaFree(d_output));
checkCudaErrors(cudaFree(d_arr));
checkCudaErrors(cudaFree(d_arr2));
}
```

The above code<sup>3</sup> is the final resulting code for the evaluation task. We implement the given CPU algorithm as a parallel GPU algorithm using CUDA.

The CUDA kernel sub takes two arrays that reside on the GPU and returns a GPU array which stores the result of subtracting the value of the array at the position (indices + offset). We call sub from the function prefix\_sum which also calls Thrust's exclusive scan [4] algorithm on the resulting device array to get the required result.

```
thrust:: exclusive\_scan(data.begin(), data.end(), temp.begin());\\
```

and since the problem in question is not exactly an exclusive scan where the last cumulative sum number is not calculated, we need an additional step to calculate the required result -

```
temp [\,data.\,size\,(\,)\,] \ = \ data.\,back\,(\,) \ + \ temp [\,data.\,size\,(\,) \ - \ 1\,]\,;
```

We calculate the number of blocks and threads required by the sub CUDA kernel by -

```
if (length > 1024) {
    block = (length / 1024) + 1;
    thread = 1024;
}
else {
    thread = length;
    block = 1;
}
```

The prefix\_sum() function also has a lot of boilerplate code such as copying arrays from CPU to GPU, and back from the GPU to CPU whose execution time we do not take into consideration while benchmarking.

Arguably CUB's [5] exclusive scan algorithm is more optimized than Thrust's exclusive scan algorithm, but since Thrust comes installed with the CUDA development toolkit, we have decided to use Thrust in our sample code.

<sup>&</sup>lt;sup>3</sup>Full benchmarking code: https://github.com/reikdas/GSoC-Proposal-2020/blob/master/testgsoc/thrust.cu

### 4 Benchmarks

We have compared the performance of the various algorithms relative to each other<sup>4</sup>.

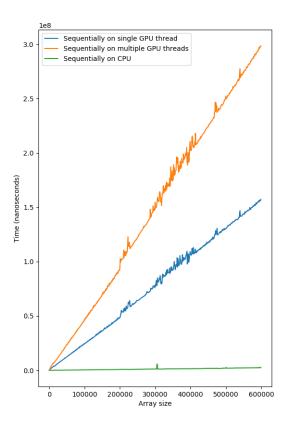


Figure 5: Comparing performance of sequential GPU algorithms with CPU algorithm

We can see that the sequential algorithms on the GPU run much slower than the sequential algorithm on the CPU<sup>5</sup>. This can be attributed to the inherent difference in the design of GPUs and CPUs. GPUs are designed to operate on independent data in parallel, whereas CPUs are optimized to execute a single stream of instructions as quickly as possible. So when we are executing a single instruction stream on the GPU, it is expected that the time for completion will be more than the time for completion of the same instructions on a CPU.

<sup>&</sup>lt;sup>4</sup>Benchmarking numbers: https://github.com/reikdas/GSoC-Proposal-2020/blob/master/testgsoc/data.txt generated using https://github.com/reikdas/GSoC-Proposal-2020/blob/master/testgsoc/plotter.py

<sup>&</sup>lt;sup>5</sup>We use an i7 8750H CPU for our tests

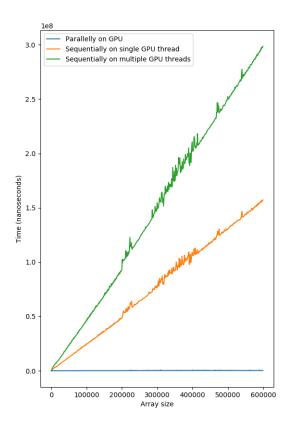


Figure 6: Comparing performance of sequential and parallel GPU algorithms

Similar to the previous plot, we can see that our parallelized GPU implementation is much faster than the implementation of our sequential algorithms on the GPU. This gives us an indication that our parallel implementation is correct and gives a clear speedup over the sequential algorithm when executed on the same hardware (GPU of our test machine in this case)<sup>6</sup>.

 $<sup>^6\</sup>mathrm{We}$ used an Nvidia Ge<br/>Force 1070 Max Q GPU for our tests

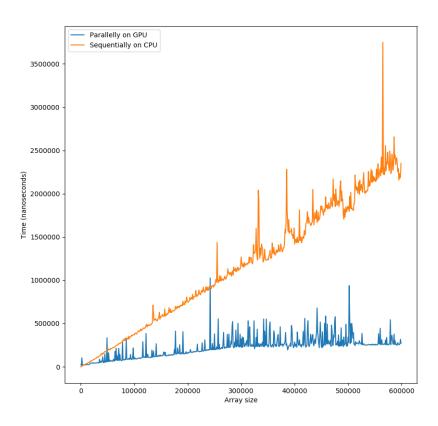


Figure 7: Comparing parallelized GPU performance vs CPU performance

This is the most important result of the evaluation task. Here, we can clearly visualize that the parallel algorithm we designed to be executed on the GPU is much faster than the given sequential algorithm on the CPU.<sup>7</sup>

# 5 Appendix

We have only demonstrated the relevant CUDA kernels and not the surrounding code or main functions which are required to execute the kernels. The entire collection of complete code samples can be found in the GitHub repository.

In our full sample code which we have used for benchmarking, we have used 2 header files which includes helper functions to facilitate CUDA error checking in our code -  $helper\_cuda.h$  and  $helper\_string.h$ . Both of these helper header files are included in the NVIDIA CUDA SDK under the samples/inc/ folder. In the Awkward Array production code we might use the same CUDA error checking header files or we might decide to write our own CUDA error checking code.

 $<sup>^{7}</sup>$ There are some non-uniform anamolous spikes in all iterations of recorded data

# References

- [1] https://www.khronos.org/opencl/.
- $[2] \ https://www.geforce.com/hardware/technology/cuda.$
- $[3] \ https://docs.nvidia.com/cuda/thrust/index.html.$
- $[4] \ https://docs.nvidia.com/cuda/thrust/index.htmlprefix-sums.$
- $[5] \ https://nvlabs.github.io/cub/.$