

# UNIFIED MEMORY ON KEPLER AND MAXWELL

Jiri Kraus, April 25th 2016

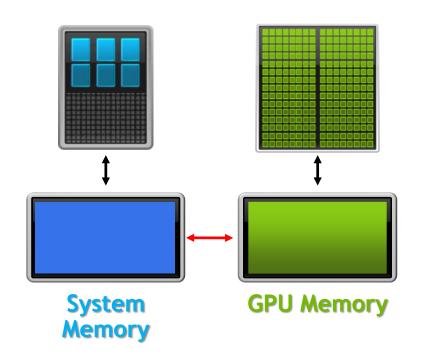


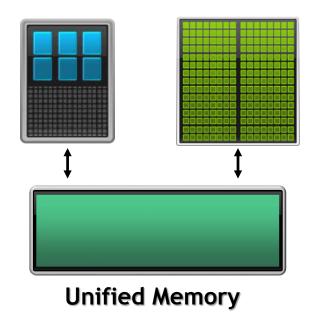
# **UNIFIED MEMORY**

Dramatically Lower Developer Effort

**Traditional Developer View** 

Developer View With Unified Memory





# **UNIFIED MEMORY**

#### Traditional Developer View

# Developer View With Unified Memory

```
void sortfile(FILE *fp, int N) {
                                            void sortfile(FILE *fp, int N) {
  char *data:
                                               char *data;
  char *data d;
  data = (char *)malloc(N);
                                               cudaMallocManaged( &data, N );
  cudaMalloc( &data d, N );
 fread (data, 1, N, fp);
                                               fread (data, 1, N, fp);
  cudaMemcpy ( data d, data, N,
              cudaMemcpyHostToDevice);
  qsort<<<...>>>(data, N, 1, compare);
                                               gsort<<<...>>>(data, N, 1, compare);
                                               cudaDeviceSynchronize();
  cudaMemcpy ( data, data d, N,
              cudaMemcpyDeviceToHost);
 use data(data);
                                               use data(data);
                                               cudaFree (data);
  cudaFree(data d); free(data); }
```

# **UNIFIED MEMORY**

#### Advantages

Unified Memory can be used in CPU and GPU code

No need for explicit device allocation(cudaMalloc) or memory copies (cudaMemcpy)

No need to fully understand data flow and allocation logic of application

Incremental profiler driven acceleration -> Data movement is just another optimization

#### UNIFIED MEMORY ON KEPLER AND MAXWELL

#### Implementations Details

Only heap and global data can be Unified Memory, e.g. no stack data:

```
qsort<<<...>>> (data Accessing Stack data from kernel not possible! kernel not possible);
```

Data is coherent only at kernel launch and sync points

Its not allowed to access unified memory in host code while a kernel is running. Doing so may result in a segmentation fault.

# CONCURRENT HOST DEVICE ACCESSES

CUDA\_LAUNCH\_BLOCKING

Its not allowed to access unified memory in host code while a kernel is running. Doing so may result in a segmentation fault:

```
ubuntu@ip-10-152-53-181:~/Pipelining$ ./pipelining
Bus error (core dumped)
```

CUDA\_LAUNCH\_BLOCKING can be used to check if missing synchronization is the issue:

```
ubuntu@ip-10-152-53-181:~/Pipelining$ CUDA_LAUNCH_BLOCKING=1 ./pipelining
Number of Primes = 151516
Runtime = 17.2459 ms
```

#### CONCURRENT HOST DEVICE ACCESSES

#### cuda-gdb

```
ubuntu@ip-10-152-53-181:~/Pipelining$ cuda-gdb ./pipelining
NVIDIA (R) CUDA Debugger
Reading symbols from /home/ubuntu/Pipelining/pipelining...(no debugging symbols
found) . . . done .
(cuda-qdb) run
Starting program: /home/ubuntu/Pipelining/./pipelining
[Thread debugging using libthread db enabled]
Using host libthread db library "/lib/x86 64-linux-gnu/libthread db.so.1".
[New Thread 0x7ffff5a0c700 (LWP 1451)]
[New Thread 0x7fffeffff700 (LWP 1452)]
[New Thread 0x7fffef7fe700 (LWP 1453)]
Program received signal CUDA EXCEPTION 15, Invalid Managed Memory Access.
0 \times 0000000000402b4f in generate data(int*, int, int) ()
(cuda-gdb)
```

# UNIFIED MEMORY ON KEPLER AND MAXWELL

#### Implementations Details

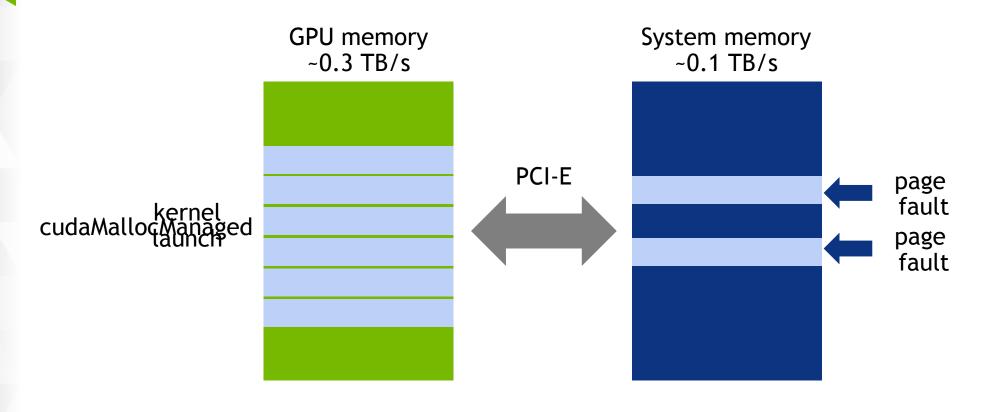
Allocation granularity of one page (4k) leads to allocation overhead (time and space) for small objects

Transfer overhead for small objects as data is transferred with the granularity of a single page

Number of allocations is limited to vm.max\_map\_count (defaults to 64k)

# UNIFIED MEMORY ON KEPLER AND MAXWELL

Implementations Details



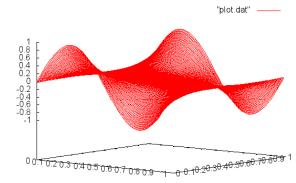
# **WEIGHTED JACOBI**

Solves the 2D-Laplace equation on a rectangle

$$\Delta u(x,y) = \mathbf{0} \ \forall \ (x,y) \in \Omega \backslash \delta \Omega$$

Dirichlet boundary conditions (constant values on boundaries) on left and right boundary

Periodic boundary conditions on top and bottom boundary



# **WEIGHTED JACOBI**

While not converged

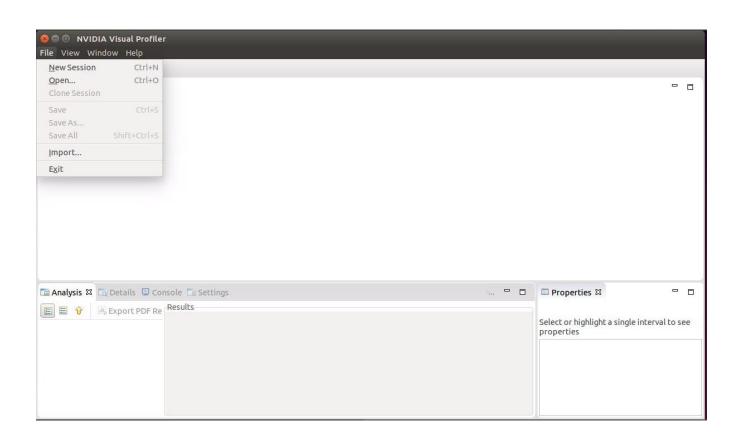
Do Jacobi step:

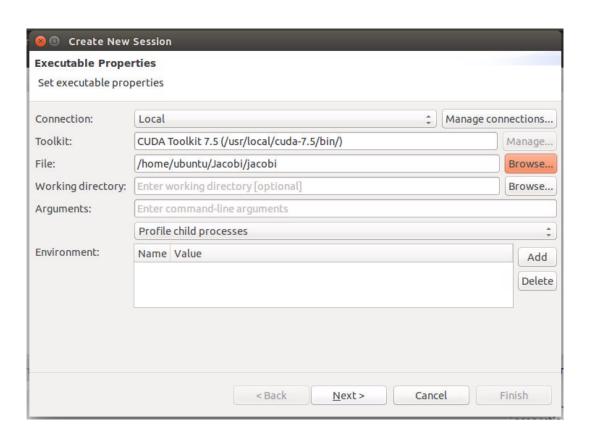
```
for (int iy=1; iy < ny-1; ++iy)</pre>
for (int ix=1; ix < nx-1; ++ix)
    float a new val = 0.25f*(a[ix-1][iy] + a[ix+1][iy] + a[ix][iy-1] + a[ix][iy+1]);
    a new[ix][iy] = weight*a new val+(1.0f-weight)*a[ix][iy];
```

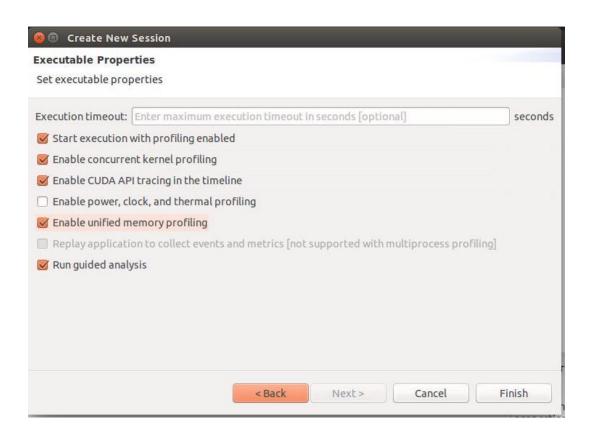
Swap a new and a

Apply periodic boundary conditions to a

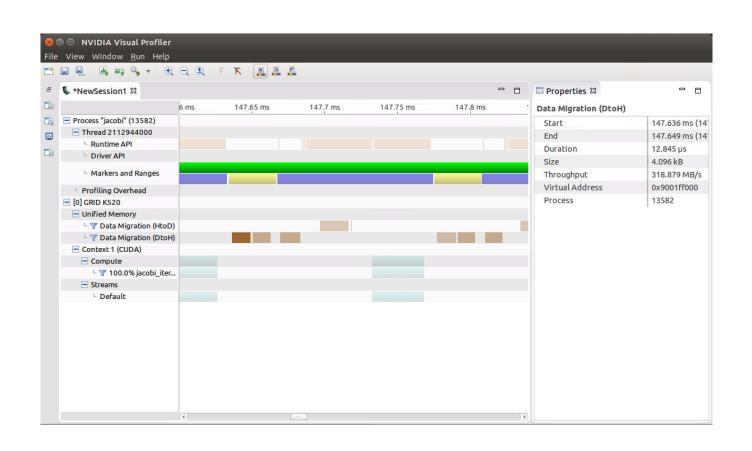
Next iteration



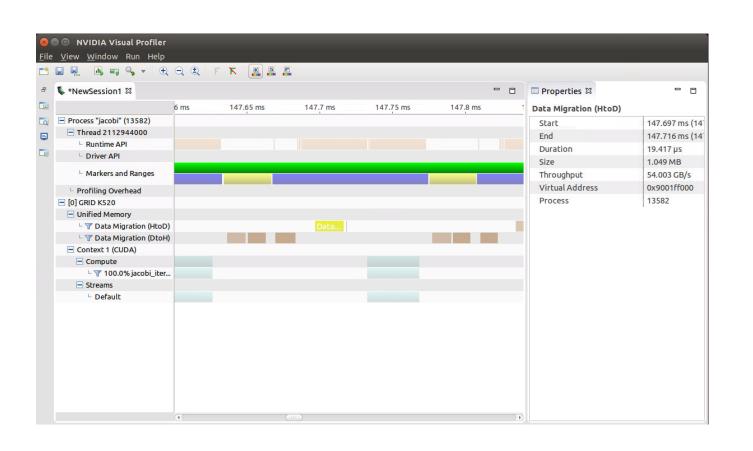




Host to Device (DtoH) Data Migration



Host to Device (HtoD) Data Migration



 $[\ldots]$ 

Runtime 0.111733 seconds.

1000

#### HANDS-ON

Task: Jacobi

Task: Avoid all data migrations within the while loop of the Jacobi solver

Apply boundary conditions with the provided GPU kernel (apply periodic bc)

Try to avoid the remaining data migrations? What's causing this?

```
-bash-4.2$ export CUDA MANAGED FORCE DEVICE ALLOC=1
-bash-4.2$ make
nvcc -DUSE NVTX -arch=sm 37 jacobi.cu -lnvToolsExt -o jacobi
./Jacobi
                                    Make Targets:
```

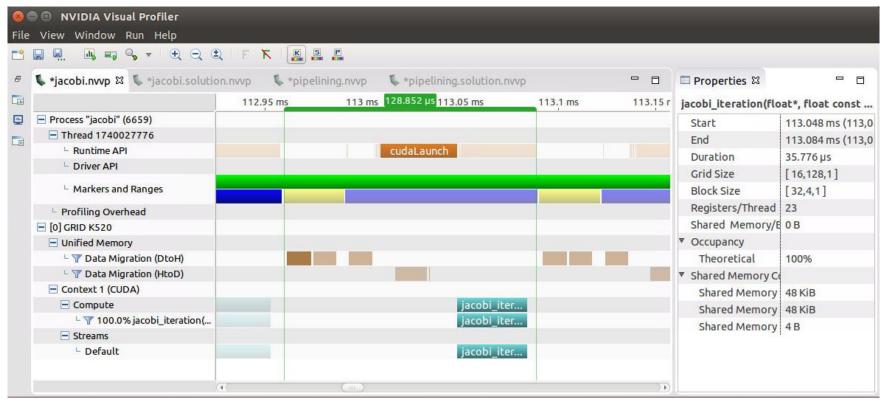
run jacobi (default) run: build jacobi binary jacobi: run with cuda-memcheck memcheck:

profile with nvprof profile:

Look for TODOs

Optional Task: What is the possible additional improvement?

#### **Initial Version**

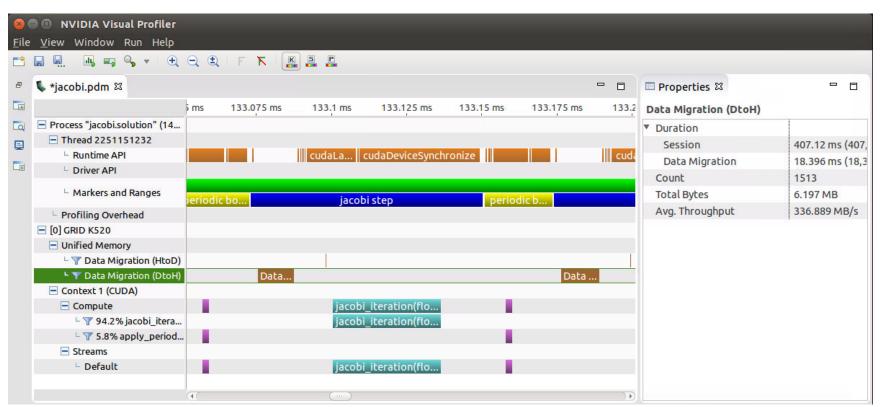


#### Apply boundary conditions GPU kernel

```
PUSH_RANGE("periodic boundary conditions",2)
//Apply periodic boundary conditions
apply_periodic_bc<<<dim3(nx/128),dim3(128)>>>(a,nx,ny);
CUDA_CALL(cudaGetLastError());
CUDA_CALL(cudaDeviceSynchronize());
POP_RANGE
```

See: <a href="https://devblogs.nvidia.com/parallelforall/cuda-pro-tip-generate-custom-application-profile-timelines-nvtx/">https://devblogs.nvidia.com/parallelforall/cuda-pro-tip-generate-custom-application-profile-timelines-nvtx/</a> for details on PUSH\_RANGE and POP\_RANGE macros.

#### Apply boundary conditions GPU kernel



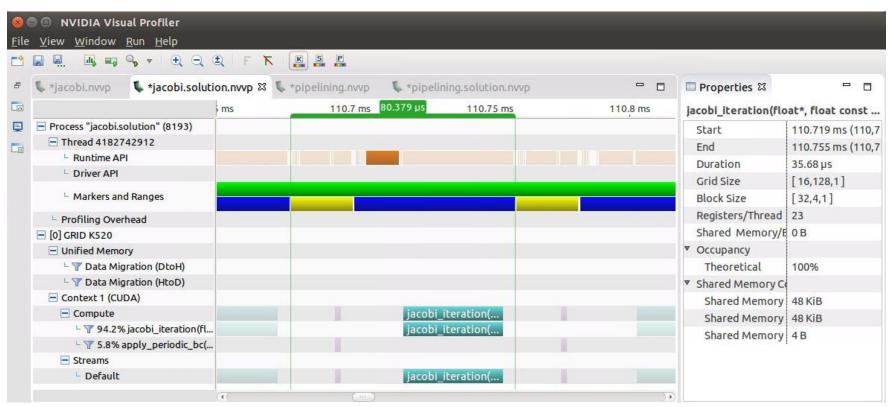
CPU reading weights[0] before launch triggers
DtoH Data Migration

See: <a href="https://devblogs.nvidia.com/parallelforall/cuda-pro-tip-generate-custom-application-profile-timelines-nvtx/">https://devblogs.nvidia.com/parallelforall/cuda-pro-tip-generate-custom-application-profile-timelines-nvtx/</a> for details on PUSH\_RANGE and POP\_RANGE macros.

#### Solution

```
const float weight = weights[0];
while ( iter <= iter max ) {</pre>
     jacobi iteration <<< dim3 (nx/32, ny/4), dim3 (32, 4) >>> (a_new, a, nx, ny, weight);
     CUDA CALL(cudaGetLastError());
     CUDA CALL(cudaDeviceSynchronize());
     std::swap(a,a new);
     //Apply periodic boundary conditions
     apply periodic bc << dim3(nx/128), dim3(128) >>> (a, nx, ny);
     CUDA CALL(cudaGetLastError());
     CUDA CALL(cudaDeviceSynchronize());
     if ( 0 == iter%100 )
                  std::cout<<iter<<std::endl; iter++;</pre>
```

#### Solution



#### **Results**

	RUNTIME (S)	HOST TO DEVICE - DATA MIGRATIONS*	DEVICE TO HOST - DATA MIGRATIONS*
Initial	0.112	982.334 MB	13.73047 MB
Final	0.040	2.003906 MB Optional Task: cudaDev	2.003906 MB

Optional Task: cudaDeviceSynchronize()
calls within while loop are no longer
necessary and can be removed for
further speedup

<sup>\*</sup>nvprof --unified-memory-profiling per-process-device ./jacobi