

■ The matmult f.nvcc driver is provided

```
matmult f.nvcc type m n k [bs]
where m, n, k are the parameters defining the matrix sizes, bs is the
optional blocksize for the block version, and type can be one of:
       - the native/naive version
nat.
lib
       - the library version (note that this now calls a multithreaded
library)
       - the first qpu version
qpu1
gpu2 - the second gpu version
gpu3 - the third gpu version
gpu4 - the fourth gpu version
gpu5 - the fifth gpu version
gpu6 - the sixth gpu version
qpulib - the CUBLAS library version
as well as blk, mnk, nmk, ... (the permutations).
```

See README for more (also week 1 README)



- Driver uses dlsym for dynamic linking (C99 library) therefore all functions in your shared library must have C naming convention
- Use extern "C" {}

```
extern "C" {
    #include <cblas.h> // The "C" headers also inside
...
    void matmult_lib(...)
    {
        ...
    }
...etc.
}
```



- New cuBlas API used in the assignment
 - □ #include <cublas v2.h>
- Creating a handle (to cuBlas context)
 - Important when using multiple host threads and multiple GPUs and makes it reentrant (non-blocking)
 - First creation has a large overhead!!

```
cublasStatus_t stat;
cublasHandle_t handle;
stat = cublasCreate(&handle);
```

■ The matmult_f.nvcc driver creates a handle (and destroys it again) to "wake up" cuBlas



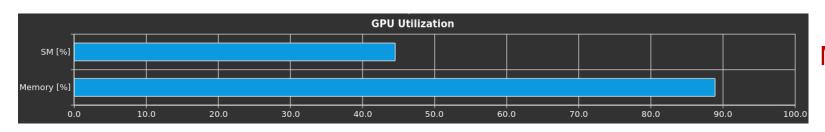
- For large matrices please use
 - □ MFLOPS MAX IT=1 ./matmult f.nvcc ...
- For benchmarks please use
 - □ MATMULT COMPARE=0
- For running on 1 core on CPU
 - ☐ MKL NUM THREADS=1
- Overflow for large matrices (Please note!!)
 - You should have the same overflow for all methods!

```
n-62-12-19(hhbs) $ ./matmult_f.nvcc gpulib 2048 2048 2048 98304.000 1178532.905 300 # matmult_gpulib
```

Overflow - not an actual error!

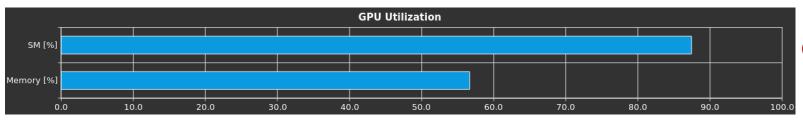


- Profiler (if available) may help to explain results
 - matmult_f.nvcc gpu2 1024 1024 1024



Memory bound!

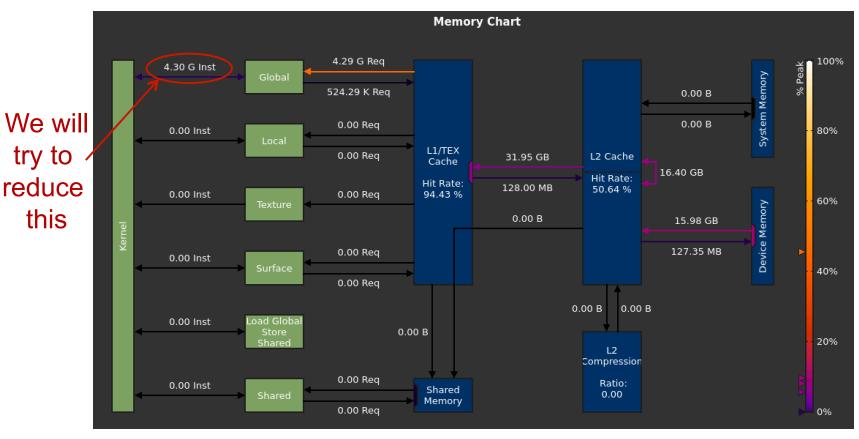
□ matmult f.nvcc gpulib 1024 1024 1024



Compute bound!



- Profiler (if available) may help to explain results
 - □ matmult f.nvcc gpu2 4096 4096 4096



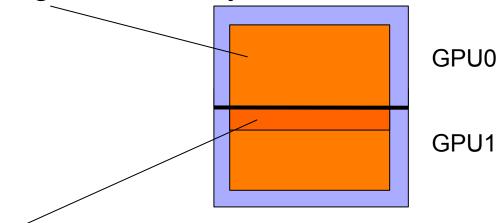
this

GPU Poisson problem



- Multi-GPU version using Peer-to-peer access
 - □ Split task into two top and bottom
 - Interior points can be updated from global memory
 - Border points must read "peer values" from other GPU

Read from global memory



Available from other GPU

Submitting GPU batch jobs



- Benchmark runs should always be submitted to queue hpcintrogpu
- See https://www.hpc.dtu.dk/?page_id=2759
- Maximum wall-clock time on jobs is 1 hour!
- For jobs using two GPUs use num=2
- For CPU-only jobs please do not request GPUs
- The -G flag sets debug lines into your code
 - This reduces the performance drastically
 - Remove it for performance tuning!!!

Reports / Analysis



- Assess your performance (Gflops / Bandwidth)
 - Profiler gives these numbers ("Throughput")
- Speed-up calculations (fair)
 - □ If using 1 GPU, compare with 1 CPU!
 - □ numactl --cpunodebind=0 to bind threads to cpu 0
- Tuning considerations
- Profiler analysis (if available)
- Relevant comments and observations
- Please keep the report format as close to the assignment questions as possible
- Please try to have one section in the report for and acceptance in the ancience at the leaves and it