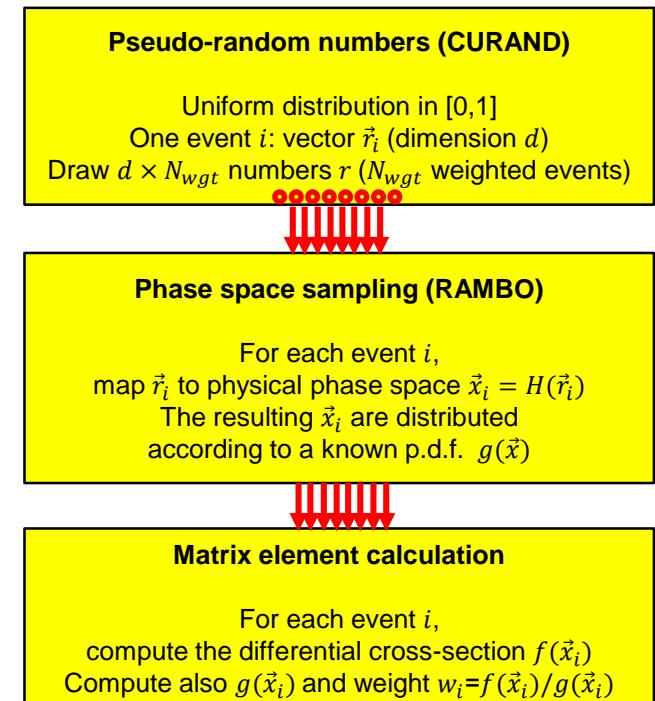


Overview

- My code is in https://gitlab.cern.ch/roiser/madgraph4gpu/-/tree/master/examples/gpu/eemumu_AV
 - Integrates my developments with some (not all...) of others' developments
 - Olivier's simplified IXX/OXX
 - Curand for random numbers
 - Not there yet: cucomplex, streams, graphs...
- General idea: single source, try to keep options open with #ifdefs
 - Cuda and c++ (symlink xxx.cc files as gxxx.cu files)
 - AOSOA (ASA), AOS, SOA memory layouts
 - Local/global/shared memory for the w[5][6] intermediate wavefunctions
 - Double and single precision (FPTYPE = double or float, CXTYPE = complex)
 - Curand generation on host and on device

Overall data flow

- Streamlined use of C-style arrays for inputs and outputs
 - Removing std::vector by itself is a large speedup in both c++ and cuda
 - Fix dimensions in advance
 - Not hardcoded: events per each iteration = ndim (=nthr*nblk i.e. #threads>#blocks)
 - Hardcoded: npar=4 (e+ e- mu+ mu-), np4=4 (E px py pz)
 - Each event also needs np4 random numbers
- Three main phases
 - Random number generation
 - Output: FPTYPE rnarray [ndim * npar * np4]
 - Need np4=4 random numbers per particle
 - (even if only 3 degrees of freedom)
 - Rambo: map random numbers to momenta
 - Input: rnarray [ndim * npar * np4]
 - Output: FPTYPE allmomenta[ndim * npar * np4]
 - Output: FPTYPE weights[ndim]
 - Sigmakin: from momenta to MEs
 - Input: allmomenta[ndim * npar * np4]
 - Output: FPTYPE MEs[ndim]
 - Physics validation: mean and stddev of MEs
 - Missing: proper computation with weights



Memory layout for allmomenta

- Current default: AOSOA (ASA): `allmomenta[npag][npar][np4][nepp]`
 - Partition ndim events per iteration into npag pages of nepp events per page
 - i.e. $\text{ndim} = \text{npg} * \text{nepp}$
 - Currently $\text{nepp}=32$ is hardcoded (#threads per GPU warp)
 - Eventually? Use $\text{nepp} = \# \text{threads per block}$ i.e. `allmomenta[nblk][npar][np4][nthr]`
- Alternative SOA: `allmomenta[npar][np4][ndim]`
- Alternative AOS: `allmomenta[ndim][npar][np4]`
- *Idea was to try and exploit memory coalescing (and SIMD?) in the GPU*
 - *BUT: no obvious performance benefit/penalty in any of these choices* ☹
- *Probably best to keep AOSOA anyway?*
 - `allmomenta[nblk][npar][np4][nthr]`?

Local memory and shared memory

- This refers to the intermediate wavefunctions $w[5][6]$ in each event
 - Where $w[0:3]$ are the $npar=4$ particles, $w[4]$ is a tmp for internal particles
- Current default: LOCAL $w[5][6]$ on the stack in each event
 - “local” is actually “thread-local” global memory
- Alternative SHARED: $sw[5][6][nthr]$ for all events in one GPU block
 - BUT 32 threads is $32*5*6*2(\text{complex})*8(\text{double})=15\text{kb}$, GPU has only 48kb
 - Probably non scalable, and worse performance than local (to be understood)
- Alternative GLOBAL: $gw[5][6][ndim]$ for all events in all GPU blocks
 - BUT this soon exhausts global memory (malloc and/or curand fail)
- *Idea was to try and exploit memory coalescing (and SIMD?) in the GPU*
 - *And also to reduce the number of registers and improve performance*
 - *BUT: LOCAL seems to have better performance without added issues*
- *Clearly best to keep LOCAL at the moment*
 - *Keep options in the code to do other things? (may also help in CPU SIMD?)*
 - *En passant: LOCAL “-p 16384 32 12” seems faster than “-p 2048 256 12”?*

Random number generation

- Curand can be used both on the CPU (host) and on the GPU (device)
 - It produces the same numbers in both options
 - Useful to get strict reproducibility of physics results
 - NB: the seeds are reinitialized at every new iteration of ndim events
- Currently: using the fastest curand generator, not the best for physics
 - Fastest: CURAND_RNG_PSEUDO_MTGP32
 - Best for physics (Lorenzo Moneta): CURAND_RNG_PSEUDO_MR32K3A
 - A factor 500 in speed between the two options
 - The point is only to get the GPU machine going, and same results on CPU

```
time ./check.exe -p 16384 32 12
```

```
*****  
NumIterations = 12
```

```
NumThreadsPerBlock = 32
```

```
NumBlocksPerGrid = 16384
```

```
-----  
FP precision = DOUBLE (nan=0)
```

```
Momenta memory layout = AOSOA[32]
```

```
Curand generation = HOST (C++ code)
```

```
-----  
NumberOfEntries = 12
```

```
TotalTimeInWaveFuncs = 1.741161e+01 sec
```

```
MeanTimeInWaveFuncs = 1.450968e+00 sec
```

```
StdDevTimeInWaveFuncs = 1.619446e-03 sec
```

```
MinTimeInWaveFuncs = 1.449295e+00 sec
```

```
MaxTimeInWaveFuncs = 1.449451e+00 sec
```

```
-----  
ProcessID: = 17797
```

```
NProcesses = 1
```

```
NumMatrixElementsComputed = 6291456
```

```
MatrixElementsPerSec = 3.613368e+05 sec^-1
```

```
*****  
NumMatrixElements(notNan) = 6291456
```

```
MeanMatrixElemValue = 1.394735e-02 GeV^0
```

```
StdErrMatrixElemValue = 3.034488e-06 GeV^0
```

```
StdDevMatrixElemValue = 7.611337e-03 GeV^0
```

```
MinMatrixElemValue = 6.071582e-03 GeV^0
```

```
MaxMatrixElemValue = 3.374925e-02 GeV^0
```

```
*****
```

Current CPP baseline

```
*****
```

```
0a ProInit : 0.000451 sec
```

```
0b MemAlloc : 0.053021 sec
```

```
0c GenCreat : 0.001038 sec
```

```
1a GenSeed : 0.000020 sec
```

```
1b GenRnGen : 0.359951 sec
```

```
2a Rambolni : 0.130261 sec
```

```
2b RamboFin : 2.077701 sec
```

```
3a SigmaKin : 17.411612 sec
```

```
4a DumpLoop : 0.015827 sec
```

```
9a DumpAll : 0.046769 sec
```

```
9b GenDestr : 0.000112 sec
```

```
9c MemFree : 0.001842 sec
```

```
    TOTAL : 20.098604 sec
```

```
*****
```

```
real 0m20.108s
```

```
user 0m20.060s
```

```
sys 0m0.045s
```



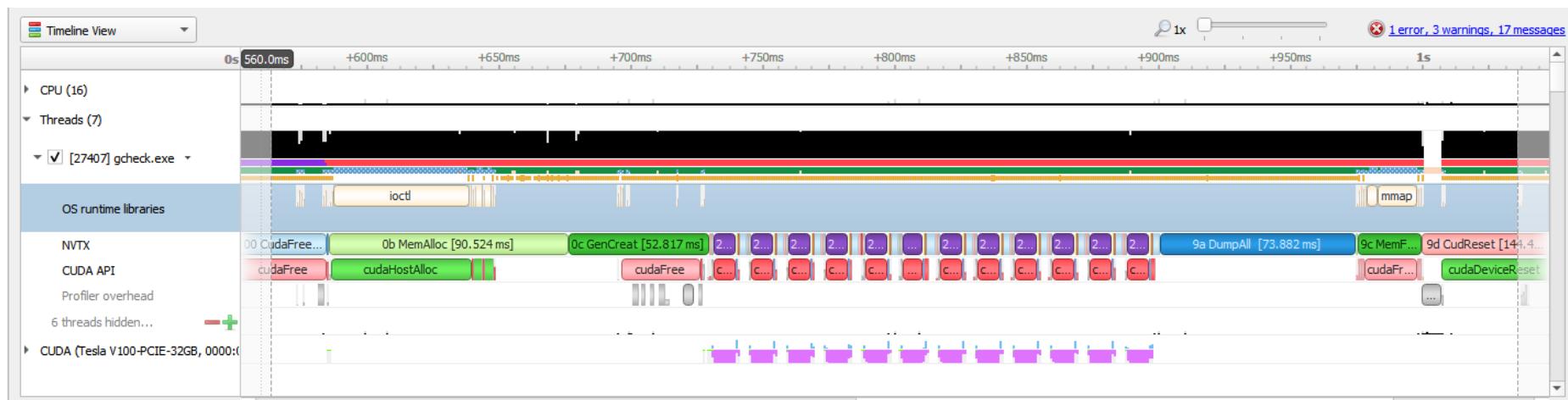
Current CUDA baseline

```
time ./gcheck.exe -p 16384 32 12
*****
NumIterations      = 12
NumThreadsPerBlock = 32
NumBlocksPerGrid   = 16384
-----
FP precision       = DOUBLE (nan=0)
Momenta memory layout = AOSOA[32]
Wavefunction GPU memory = LOCAL
Curand generation   = DEVICE (CUDA code)
-----
NumberOfEntries    = 12
TotalTimeInWaveFuncs = 1.257645e-02 sec
MeanTimeInWaveFuncs = 1.048037e-03 sec
StdDevTimeInWaveFuncs = 2.003851e-05 sec
MinTimeInWaveFuncs = 1.035607e-03 sec
MaxTimeInWaveFuncs = 1.040581e-03 sec
-----
ProcessID:         = 17851
NProcesses:        = 1
NumMatrixElementsComputed = 6291456
MatrixElementsPerSec = 5.002570e+08 sec^-1
*****
NumMatrixElements(notNaN) = 6291456
MeanMatrixElemValue     = 1.394735e-02 GeV^0
StdErrMatrixElemValue   = 3.034488e-06 GeV^0
StdDevMatrixElemValue   = 7.611337e-03 GeV^0
MinMatrixElemValue      = 6.071582e-03 GeV^0
MaxMatrixElemValue      = 3.374925e-02 GeV^0
*****
```

*Our “historical” throughput is 3.6E5 against 5.0E8 :
GPU ~1500 faster than one CPU core
Here “throughput” is ONLY sigmakin + copy MEs to host:
For unweighted evt generation should add copy of momenta!*

```
*****
00 CudaFree : 0.192427 sec
0a ProInit : 0.000582 sec
0b MemAlloc : 0.062740 sec
0c GenCreat : 0.017667 sec
1a GenSeed : 0.000017 sec
1b GenRnGen : 0.007964 sec
2a Rambolni : 0.000124 sec
2b RamboFin : 0.000062 sec
2c CpDTHwgt : 0.008801 sec
2d CpDTHmom : 0.106680 sec
3a SigmaKin : 0.000106 sec
3b CpDTHmes : 0.012470 sec
4a DumpLoop : 0.022217 sec
9a DumpAll : 0.046581 sec
9b GenDestr : 0.000321 sec
9c MemFree : 0.022609 sec
9d CudReset : 0.061855 sec
TOTAL : 0.563224 sec
*****
real 0m0.578s
user 0m0.211s
sys 0m0.344s
```





Miscellanea

- Various other items in random order on the todo list
 - Software issues
 - Namespaces
 - Variable names and where to hardcode them (also need short names)
 - General cleanup
 - Physics validation
 - Full chain: use weights for cross section and for unweighted event generation
 - C++ code
 - Complete new memory layouts, try vectorization?
 - Complex arithmetics and memory
 - Fewer registers with cucomplex?
 - Use RRRRIII layout rather than RIRIRIRI?
 - Helicities
 - Helicity loop as extra parallelism dimension?
 - Helicity masking (I implemented a very simple version in cuda)