

MadFlow: towards the automation of Monte Carlo simulation on GPU for particle physics processes

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Outline

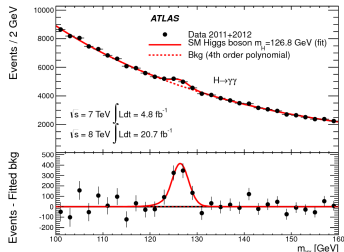
- 1 Motivation
 - Introduction
 - How can we do better
 - Tensors, tensors everywhere
- 2 The Flow suite: VegasFlow, PDFFlow, and MadFlow
 - The what, the where and the how
- 3 Benchmarks and examples
 - PDF interpolation
 - Automatic cross section integration
- 4 Conclusions

Parton-level Monte Carlo generators

Behind most predictions for LHC phenomenology lies the numerical computation of the following integral:

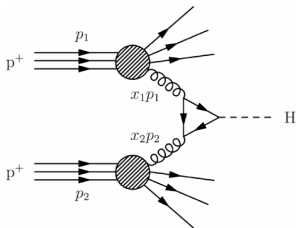
$$\int dx_1 dx_2 f_1(x_1, q^2) f_2(x_2, q^2) |M(\{p_n\})|^2 \mathcal{J}_m^n(\{p_n\})$$

- $f(x, q)$: Parton Distribution Function
- $|M|$: Matrix element of the process
- $\{p_n\}$: Phase space for n particles.
- \mathcal{J} : Jet function for n particles to m .

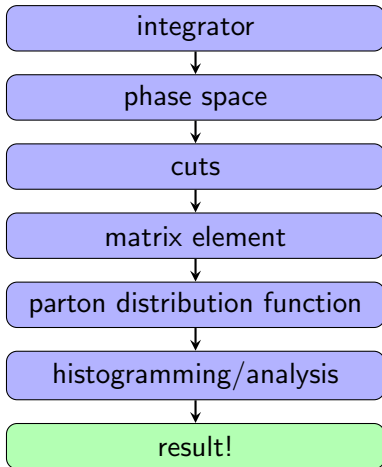


Parton-level Monte Carlo generators ingredients:

$$\int dx_1 dx_2 f_1(x_1, q^2) f_2(x_2, q^2) |M(\{p_n\})|^2 \mathcal{J}_m^n(\{p_n\})$$



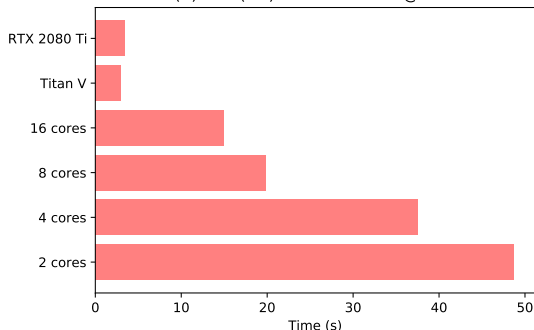
The integrals are usually computed numerically using CPU-expensive Monte Carlo generators.



GPU computing

Monte Carlo simulations are highly parallelizable, which make them a great target for GPU computation.

Float-64 performance comparison for a MC integral
Intel(R) Core(TM) i9-9980XE CPU @ 3.00GHz



Quick Example:

n -dimensional gaussian function

$$I = \int dx_1 \dots dx_n e^{x_1^2 + \dots + x_n^2}$$

Every event is independent of all other events!

GPU computation can increase the performance of the integrator by more than an order of magnitude.

Why is then GPU computing not more widespread in HEP?

Most of the more advance theoretical calculations still rely exclusively on CPU.

✗ Diminishing returns

- Huge CPU-optimized Fortran 77/90 or C++ codebases.
- Publication-ready results are easily obtained expanding existing code.
- It's catch-22: porting the code becomes more and more complicated.

✗ Lack of expertise

- CPU expertise is not necessarily applicable to GPU programming.
- New programming languages: Cuda? OpenCL?
- Low-reward situation when trying to achieve previous performance.

✗ Lack of tools

- Many ready-made tools for CPU.
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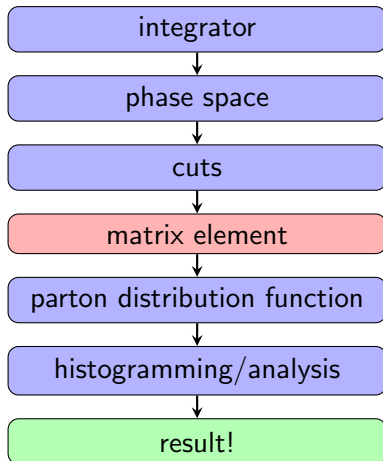
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Lack of Tools

Running on a CPU:

Worry only about what you are interested in. For instance, if we want an NNLO computation for $H \rightarrow j$ and we have a $Z \rightarrow j$ computation we only need to change the matrix elements.



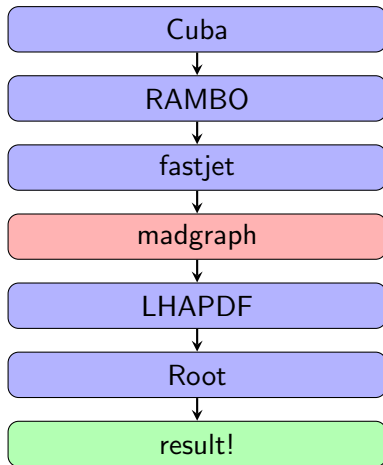
Lack of Tools

Running on a CPU:

Even if we don't already have some obscure and private fortran-based framework already built, there exists a complete tool set for producing results.

- ✓ PDF providers
- ✓ Phase space generators
- ✓ Integrator libraries...

some of which can still provide that sweet 70s' Fortran taste



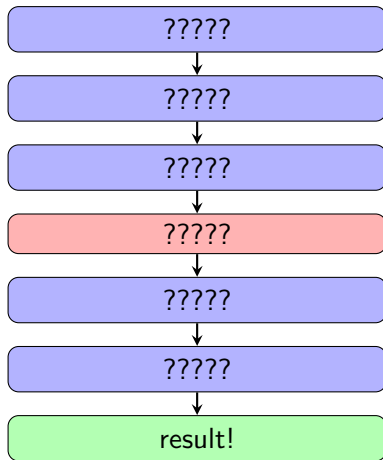
Lack of Tools

Running on a GPU:

There is no such tool set yet



so it needs to be written from scratch



Filling up the box

- 1 A phase space generator which takes an array of (`n_events`) random numbers and returns (`n_events`) an array of phase space points space points.

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- 3 An integrator framework able to send and receive batches to and from the GPU ✓
- 4 A tool for the evaluation of Matrix Elements (at tree or loop level) parallelized on the received phase (tree ✓, loop tbd)
- 5 Analysis tools, experiment simulation, jet algorithms...

Filling up the box: moving with the flow

The flow suite focus on speed and efficiency for both the computer and the developer

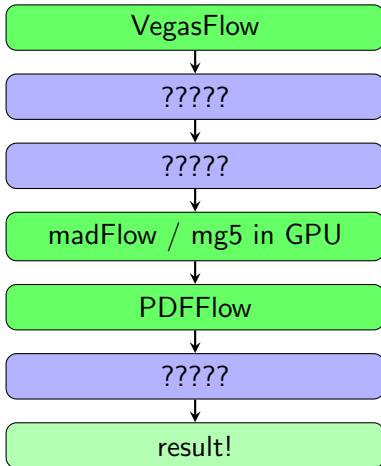
- Python and TF based engine
- Compatible with other languages: Cuda, C++
- Seamless CPU and GPU computation out of the box
- Easily interfaceable with NN-based integrators

Source code available at:

github.com/N3PDF/VegasFlow

github.com/N3PDF/PDFFlow

github.com/N3PDF/madflow

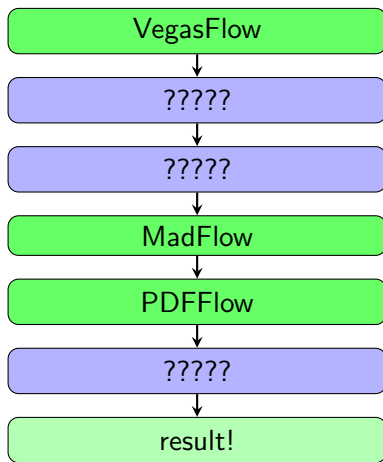


Interface with Madgraph's matrix generation

As a first step towards a full parton-level fixed-order Monte Carlo generator we interface the "flow" suite with Madgraph's matrix element generator.

We take advantage of ALOHA to produce tensorized versions of the matrix elements that can be efficiently run in GPU. With TensorFlow we can easily run in AMD and Nvidia cards.

We aim to be modular enough that different ME providers can be used or even combined (work towards a pure CUDA implementation underway)

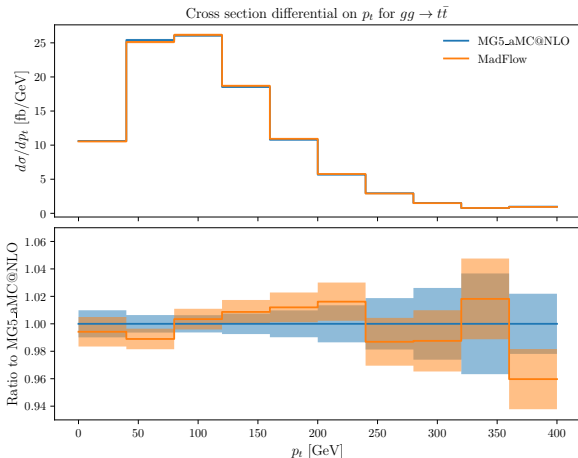


Phenomenological Results

Exact same ME and
feynman diagrams ✓

Using “RamboFlow” as
phase space ✗

Perfect compatibility ✓



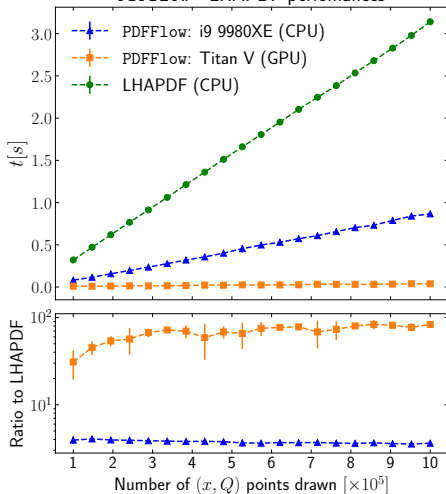
Benchmarks and examples

To wrap it up, we will see some examples and benchmarks that show how the parallelization (and tensorization!) of calculations can speed them up enormously.

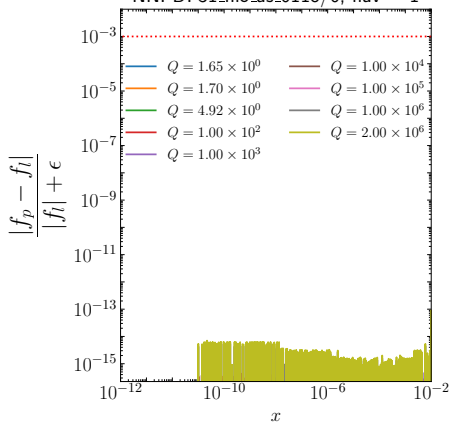
- ✓ Parallel PDF interpolation
- ✓ A completely automatic LO calculation, CPU vs GPU
- ✓ Generation of unweighted events

LHAPDF vs PDFFlow

PDFFlow - LHAPDF performances

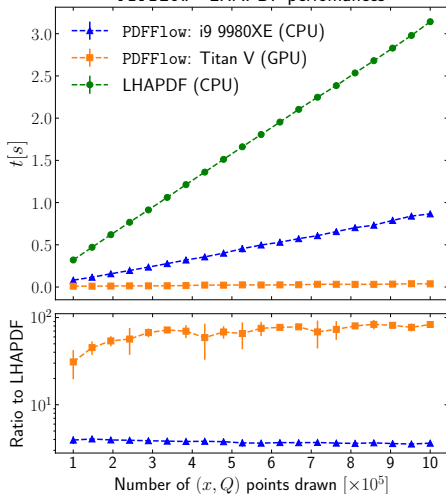


NNPDF31_nlo_as_0118/0, flav = 1

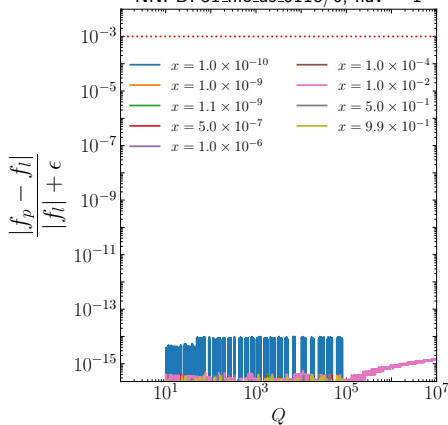
Interpolation in x for fixed q .

LHAPDF vs PDFFlow

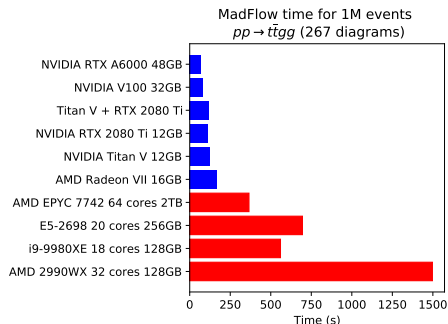
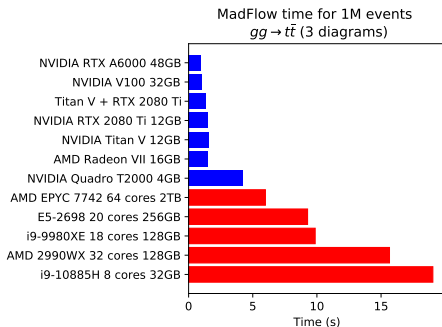
PDFflow - LHAPDF performances



NNPDF31_nlo_as_0118/0, flav = 1

Interpolation in q for fixed x .

MadFlow Vs Madgraph LO



- ✓ PhaseSpace for VegasFlow: a GPU version of RAMBO
- ✓ There's even room for improvement if a clever phase space were to be used!

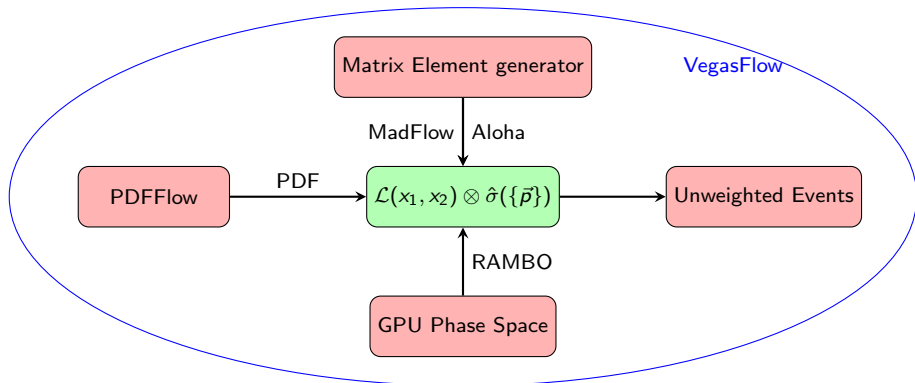
- ✓ Almost a factor ten between professional-grade CPUs (AMD EPYC) and consumer (gamer) GPUs (RTX 2080)

The MadFlow prototype

open source released

the code for madflow is available at

<https://github.com/n3pdf/madflow>



Summary

- GPU computation is increasingly gaining traction in many areas of science but it is still not heavily used in particle physics phenomenology.
- Is competitive with CPU for MC simulations.
- A lot of effort on GPU-based computations.
- ✓ VegasFlow, PDFFlow and MadFlow provide a framework to run in any device.
- ✓ Generate all the different pieces (ME, PS, PDFs, integration algorithm) needed for fixed order calculations.

Available open source

the code for madflow is available at

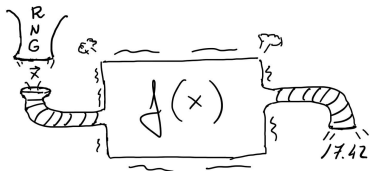
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Thanks!

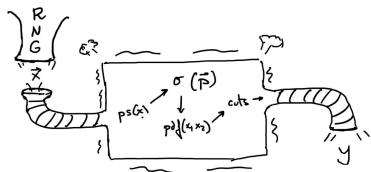
Act in parallel: CPU

The way we do Monte Carlo calculations in CPU already allows for a certain degree of parallelization

$$I = \frac{1}{N} \sum f(\vec{x}_i)$$



(the function $f(\vec{x})$ might be arbitrarily complicated)



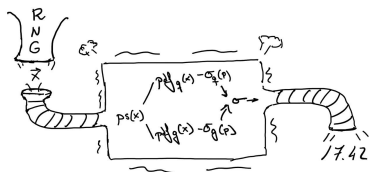
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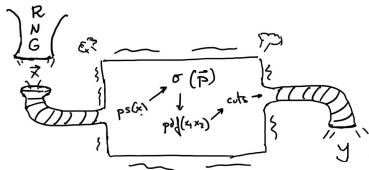
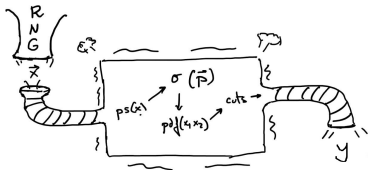
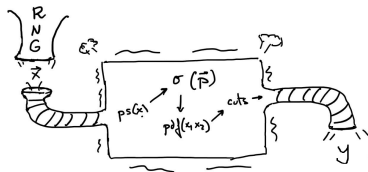
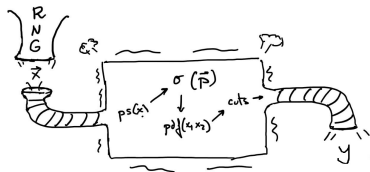


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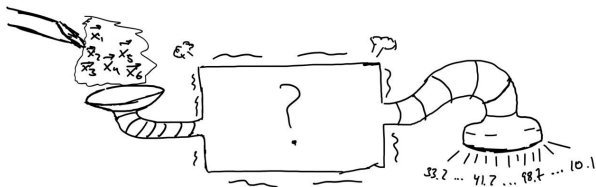
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Act in parallel: GPU

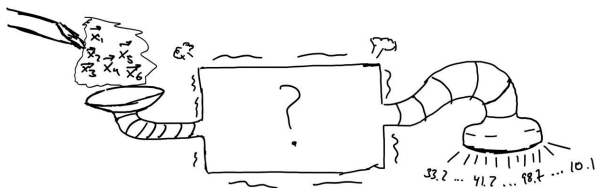
What can we do then in these machines?



We need a completely different machine, which takes a different input and a different output

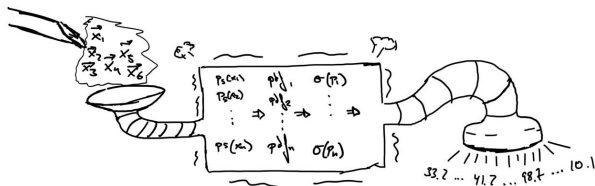
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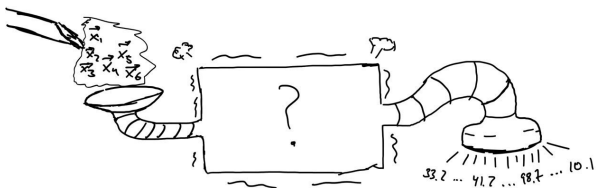
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All operations must act on all inputs at once!



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