

Event Generators: Automated Computation of Leading Order QCD Matrix Elements Part 2

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Outline

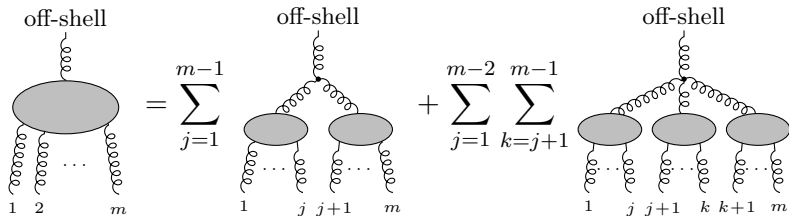
- 1) Reminder: Color Decomposition and Berends-Giele Recursion
- 2) Computation of Observables and Helicity Treatment
- 3) Data and Results (Christopher)

Reminder: Computation of Scattering Amplitudes at LO

Matrix elements can be computed using *color decomposition*:

$$\mathcal{A} = \sum_{P \in S_{n-1}} c_P A_P, \quad |\mathcal{A}|^2 = \sum_{\substack{P \in S_{n-1} \\ P' \in S_{n-1}}} A_P^\dagger C_{PP'} A_{P'}$$

A_P are the *partial amplitudes*, which depend on momenta and helicities of the external particles, and their permutation given by P . They can be computed via the Berends-Giele recursion relations



Computation of Observables

We cannot measure matrix elements! Instead we measure observables \mathcal{O} :

$$\langle \mathcal{O} \rangle \propto \int d\phi_{n-2} \mathcal{O}(\phi_n) \overline{|\mathcal{A}(\phi_n)|^2}$$

Integration over final state momenta (phase space) performed by Monte Carlo integration methods.

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Monte Carlo integration:

Evaluate integrand at N randomly chosen phase space points ϕ_{n-2} and take average as estimate for the integral. (N large!)

$$\int f(\vec{x}) d\vec{x} \approx \frac{1}{N} \sum_{i=1}^N f(\vec{x}_i)$$

Helicity Treatment

Most colliders measure *unpolarized* cross sections, not sensitive to particle helicities/spins!

↪ Unpolarized amplitude by summation over helicities:

$$\overline{|\mathcal{A}|^2} = \frac{1}{4} \underbrace{\sum_{\lambda_1} \sum_{\lambda_2} \cdots \sum_{\lambda_n}}_{2 \cdot 2 \cdots 2 = 2^n \text{ helicities}} |\mathcal{A}_{\lambda_1 \lambda_2 \cdots \lambda_n}|^2$$

⇒ Have to evaluate matrix element 2^n times, once for every helicity configuration of external particles!

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⇒ very slow, but:

We can do better!

Helicity Sampling

Necessary property for polarization vectors:

$$\sum_{\lambda=\pm} \epsilon_{\mu}^{\lambda} (\epsilon_{\nu}^{\lambda})^{*} = \epsilon_{\mu}^{+} (\epsilon_{\nu}^{+})^{*} + \epsilon_{\mu}^{-} (\epsilon_{\nu}^{-})^{*}$$

Idea: Use different polarization to achieve same result w/o sum!

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Naive attempt:

$$\begin{aligned} \epsilon_{\mu} &= \epsilon_{\mu}^{+} + \epsilon_{\mu}^{-} \\ \epsilon_{\mu} (\epsilon_{\nu})^{*} &= \epsilon_{\mu}^{+} (\epsilon_{\nu}^{+})^{*} + \epsilon_{\mu}^{-} (\epsilon_{\nu}^{-})^{*} + \epsilon_{\mu}^{+} (\epsilon_{\nu}^{-})^{*} + \epsilon_{\mu}^{-} (\epsilon_{\nu}^{+})^{*} \end{aligned}$$

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Better idea:

$$\epsilon_{\mu}(\theta) = e^{i\theta} \epsilon_{\mu}^{+} + e^{-i\theta} \epsilon_{\mu}^{-}, \quad 0 \leq \theta < 2\pi$$

$$\frac{1}{2\pi} \int_0^{2\pi} d\theta [\epsilon_{\mu}(\theta) \epsilon_{\nu}^{*}(\theta)] = \frac{1}{2\pi} \int_0^{2\pi} d\theta [\epsilon_{\mu}^{+} (\epsilon_{\nu}^{+})^{*} + \epsilon_{\mu}^{-} (\epsilon_{\nu}^{-})^{*} + e^{2i\theta} \epsilon_{\mu}^{+} (\epsilon_{\nu}^{-})^{*} + e^{-2i\theta} \epsilon_{\mu}^{-} (\epsilon_{\nu}^{+})^{*}]$$

Helicity Sampling

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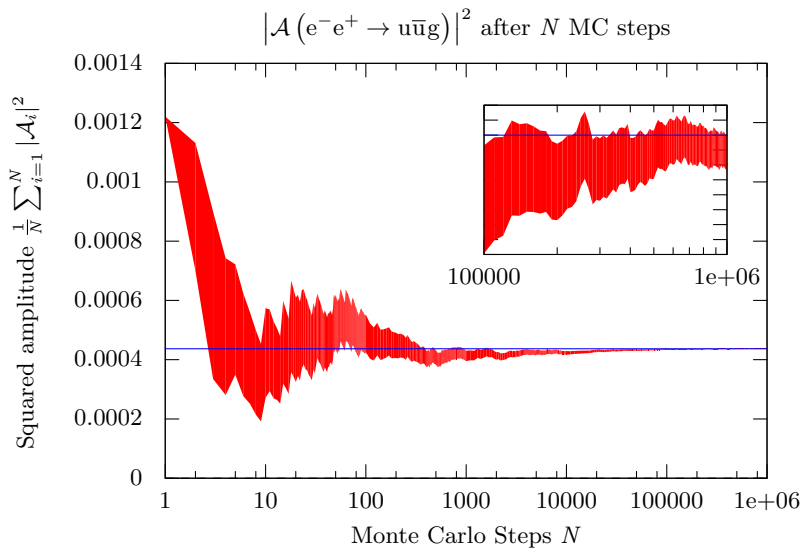
$$\epsilon_{\mu}(\theta) = e^{i\theta} \epsilon_{\mu}^{+} + e^{-i\theta} \epsilon_{\mu}^{-}, \quad 0 \leq \theta < 2\pi$$

$$\frac{1}{2\pi} \int_0^{2\pi} d\theta [\epsilon_{\mu}(\theta) \epsilon_{\nu}^{*}(\theta)] = \sum_{\lambda=\pm} \epsilon_{\mu}^{\lambda} (\epsilon_{\nu}^{\lambda})^{*}$$

⇒ Integration can be combined with phase space integral!

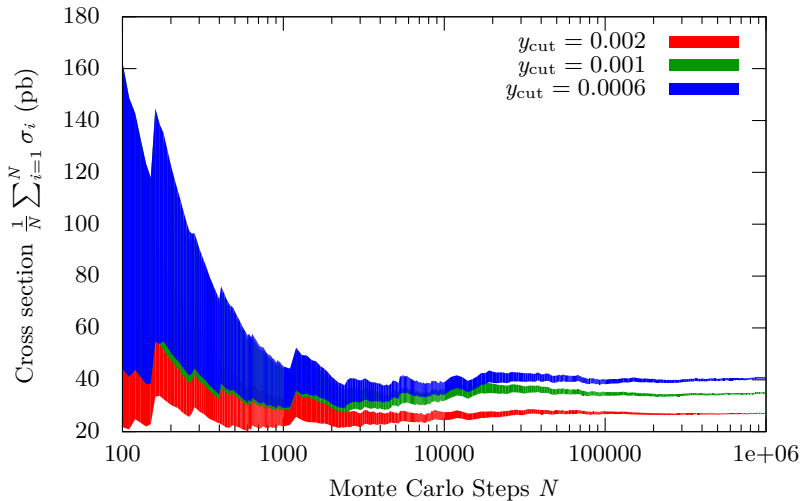
Saves factor of 2^n evaluations of squared matrix elements!

MC Integration (Helicity)



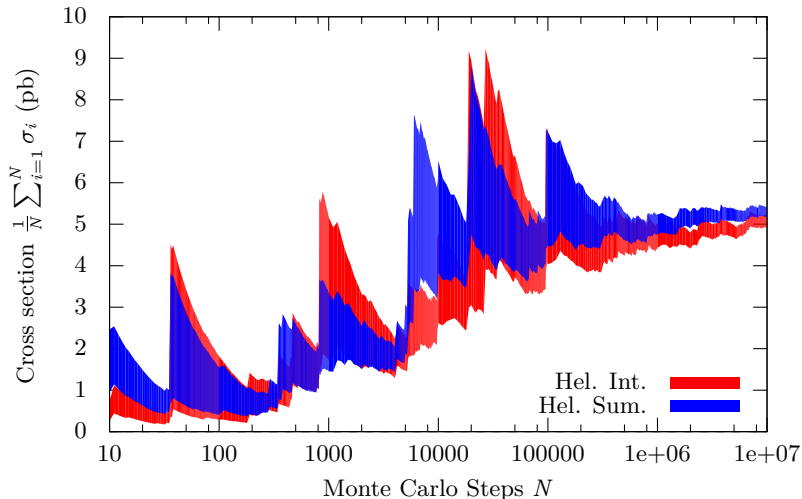
MC Integration (Helicity+Phase Space)

$\sigma(e^-e^+ \rightarrow u\bar{u}g)$ after N MC steps



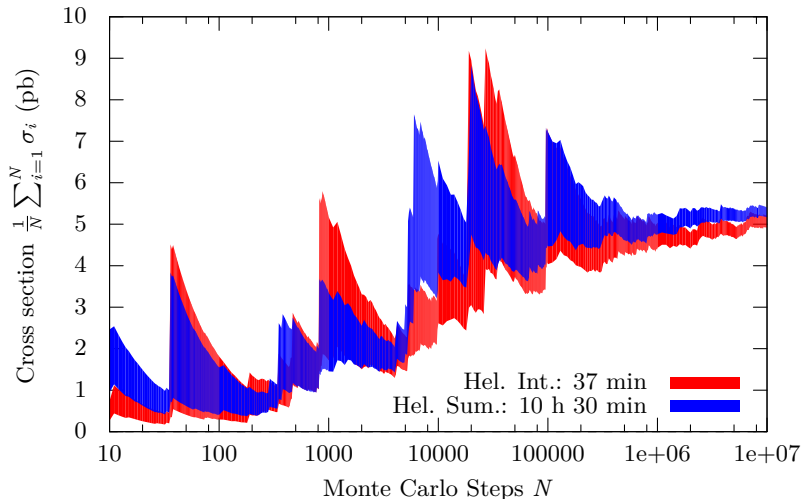
MC Integration (Helicity+Phase Space)

$\sigma_{\text{LO}}(e^-e^+ \rightarrow u\bar{u}ggg)$ after N MC steps, $y_{\text{cut}} = 0.0006$

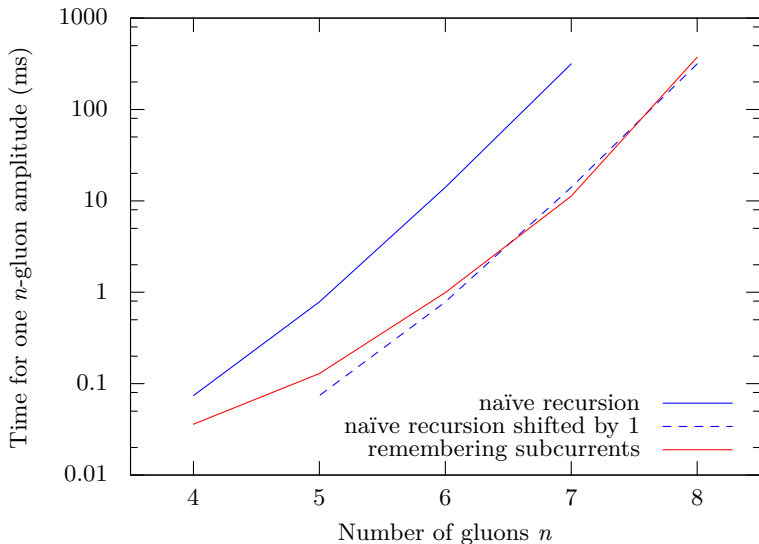


MC Integration (Helicity+Phase Space)

$\sigma_{\text{LO}}(e^-e^+ \rightarrow u\bar{u}ggg)$ after N MC steps, $y_{\text{cut}} = 0.0006$



Storing Subcurrents: What's the Benefit for Gluon Processes: $gg \rightarrow (n - 2)g$?



Comparison with MadGraph 5 (MG5)

- ▶ Comparison of our program with MG5 is a bit like comparing apples and oranges, but timings give “a feeling”
- ▶ MG5 is not automated: you need to ...
 - ▶ Specify process first (e.g. $gg \rightarrow ggggg$) in MG5's interface which then will generate Fortran code
 - ▶ Compile generated program (may take a lot of time and memory)
 - ▶ Execute the program (very fast).
- ▶ Our program is automated: a user specifies a process and our program outputs the result. There are no intermediate steps

Comparison with MG5: Timings

- ▶ Computing time for LEP amplitude of the type $|\mathcal{A}(e^-e^+ \rightarrow q\bar{q} + ng)|^2$ in milliseconds:

n	MG5 ¹	CFD+BG ²	
		Hel. Sum.	Hel. Int.
0	4(2)	0.7180(45)	0.05189(71)
1	4(2)	1.5623(50)	0.06110(20)
2	5(2)	3.7940(41)	0.08882(20)
3	7(2)	18.171(10)	0.2847(10)
4	70(5)	213.80(82)	1.876(21)
5	564(10)	3950.9(2.0)	45.25(13)
6	? ³	272270(98)	1240.09(30)

¹Execution time only

²CFD = Color Flow Decomposition, BG = Berends-Giele-Type Recursion

³Compilation aborted after 4h using more than 8 GB memory