

WHIZARD

Introduction & Tutorial



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Theoretical Challenges

Matrix Elements

Adaptive Monte Carlo

Usage

Exercise

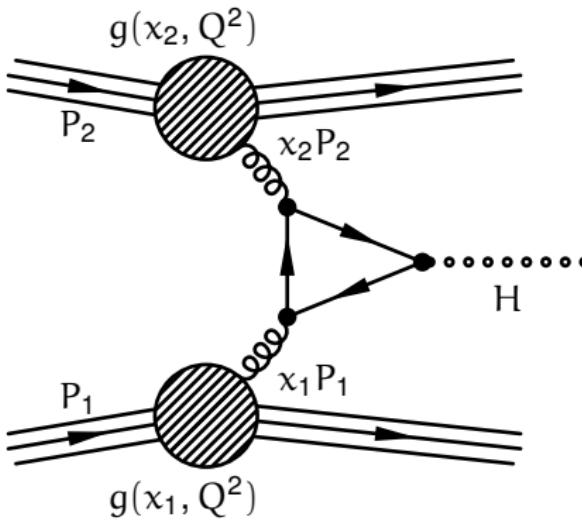
- ▶ efficiently and reliably compute scattering probabilities

$$|\langle q_1, q_2, \dots; \text{out} | T | p_1, p_2; \text{in} \rangle|^2$$

with lots of gauge cancellations among contributions

- ▶ for a multitude of physics models with qualitatively different particle content and interactions
 - ▶ standard model
 - ▶ supersymmetric extensions of the SM
 - ▶ SM with anomalous couplings
 - ▶ SM with extended gauge sector
 - ▶ SM with strongly interacting gauge bosons
 - ▶ additional space dimensions
 - ▶ ...
- ▶ such that their parameter space can be scanned and compared with experimental observations
- ▶ efficiently sample the multi particle phase space
 - ▶ scattering probabilities typically have many overlapping narrow peaks and integrable boundary singularities

- ▶ **Caveat:** not everything can be calculated in **perturbation theory** when hadrons (i. e. **strongly interacting particles**) are in play
- ▶ e. g. Higgs production at the LHC depends not only on the $gg \rightarrow H$ cross section, but also on the composition of the protons:



- ▶ asymptotic freedom and factorization allow to separate

$$\sigma(s) = \sum_{i_1 i_2} \int dx_1 dx_2 D_{i_1}(x_1, \mu) D_{i_2}(x_2, \mu) \hat{\sigma}(x_1 x_2 s; \mu)$$

independent of the scale μ !

- ▶ weakly coupled short distance/high energy phenomena, calculable in perturbation theory
 - ▶ hard scattering cross sections $\hat{\sigma}(\hat{s}; \mu)$

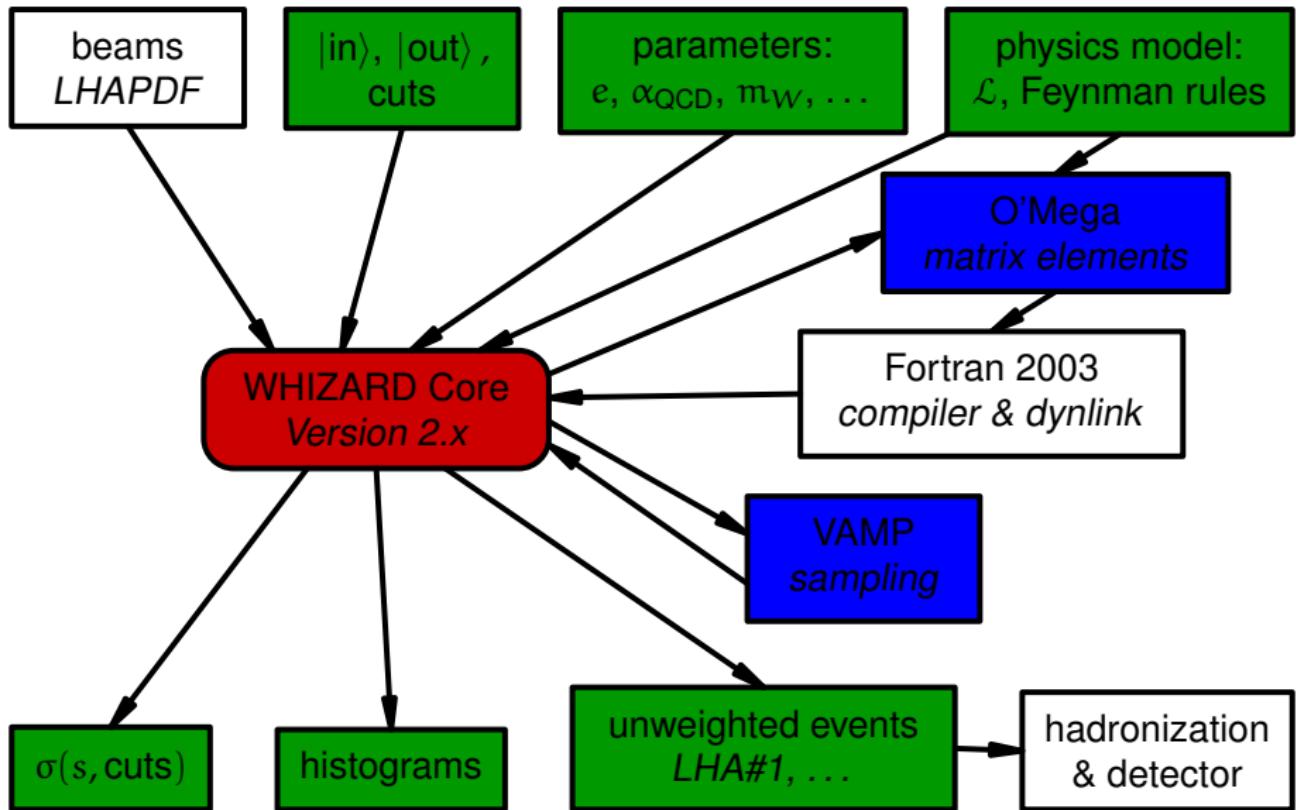
- ▶ universal strongly coupled long distance/low energy phenomena, described by parametrizations
 - ▶ parton distributions $D_{i_j}(x_j; \mu)$

and/or phenomenological models

- ▶ fragmentation and hadronization

- ▶ a series of Les Houches Accords defines interfaces implementing this separation

∴ studies of new physics can concentrate on the hard interactions!



- desired: a **computer program** implementing the function

$$(\mathcal{L}, \{\text{incoming}\}, \{\text{outgoing}\}) \mapsto \mathcal{M}(\alpha_1, \dots; p_1, \dots; s_1, \dots)$$

where

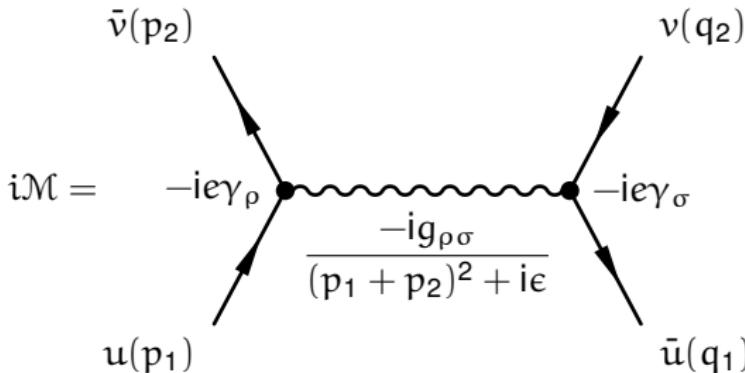
- \mathcal{L} : **Lagrangian** (or **Feynman rules**) of a **model** (SM, MSSM, ...)
- $\mathcal{M}(\alpha_1, \dots; p_1, \dots; s_1, \dots)$: a function

$$\underbrace{\mathbf{R} \times \dots \times \mathbf{R}}_{\text{masses, couplings, ...}} \times \underbrace{\mathbf{V}^+ \times \dots \times \mathbf{V}^+}_{\text{4-momenta (forward light cones)}} \times \underbrace{\mathbf{Z} \times \dots \times \mathbf{Z}}_{\text{helicities, colors}} \rightarrow \underbrace{\mathbf{C}}_{\text{amplitude}}$$

in a form that can be evaluated **numerically**, typically as C, C++ or Fortran code in that can be compiled and linked to Monte Carlo phase space integrators and generators

- NB: in some cases only $\mathcal{L} \rightarrow \sum |\mathcal{M}(\alpha_1, \dots; p_1, \dots; s_1, \dots)|^2$ is required. It is often better defined (**infrared/collinear cancellations**) and sometimes more compact (**spin/polarization sums**).
- first robust and usable examples in the early 1990s: **CompHEP**, **FeynArts**, **Grace**, **MadGraph**, ...

- for simplicity: $e^+ e^- \rightarrow \mu^+ \mu^-$ at PETRA (i. e. QED, mostly)
- just one Feynman diagram



- analytical expression

$$\begin{aligned}
 i\mathcal{M} &= \bar{v}(p_2)(-ie\gamma^\rho)u(p_1) \frac{-ig_{\rho\sigma}}{(p_1 + p_2)^2 + i\epsilon} \bar{u}(q_1)(-ie\gamma^\sigma)v(q_2) \\
 &= ie^2 \frac{1}{s} [\bar{v}(p_2)\gamma_\rho u(p_1)] [\bar{u}(q_1)\gamma^\rho v(q_2)]
 \end{aligned}$$

- ▶ corresponding Fortran95 code
(using a library for **vector** and **spinor products** and **states**)

```

pure function eleposmuamu (k, s) result (amp)
  real(kind=omega_prec), dimension(0:,:), intent(in) :: k
  integer, dimension(:, ), intent(in) :: s
  complex(kind=omega_prec) :: amp
  type(momentum) :: p1, p2, p3, p4
  type(spinor) :: muo_4, ele_1
  type(conjspinor) :: amu_3, pos_2
  type(vector) :: gam_12
  type(momentum) :: p12
  p1 = - k(:,1) ! incoming e-
  p2 = - k(:,2) ! incoming e+
  p3 =   k(:,3) ! outgoing m-
  p4 =   k(:,4) ! outgoing m+
  ele_1 = u (mass(11), - p1, s(1))           ! u s1(k1)
  pos_2 = vbar (mass(11), - p2, s(2))          ! v̄ s2(k2)
  amu_3 = ubar (mass(13), p3, s(3))            ! ū s3(k3)
  muo_4 = v (mass(13), p4, s(4))               ! v s4(k4)
  p12 = p1 + p2
  gam_12 = pr_feynman(p12, + v_ff(qlep,pos_2,ele_1)) ! (1/s) ēv(k2)γμu(k1)
  amp = 0
  amp = amp + gam_12*(+ v_ff(qlep,amu_3,muo_4))    ! (1/s) ēv(k2)γμu(k1) ēū(k3)γμv(k4)
  amp = - amp ! 2 vertices, 1 propagators
end function eleposmuamu

```

- ▶ the usual rules for **manual** calculations are **algorithmic**
∴ can be implemented in a computer program

The number of tree Feynman diagrams w/ n legs grows like a **factorial**, e.g. in ϕ^3 -theory: $F(n) = (2n - 5)!! = (2n - 5) \cdot (2n - 7) \cdot \dots \cdot 3 \cdot 1$

n	$F(n)$	$P(n)$
4	3	3
5	15	10
6	105	25
7	945	56
8	10 395	119
9	135 135	246
10	2 027 025	501
11	34 459 425	1 012
12	654 729 075	2 035
13	13 749 310 575	4 082
14	316 234 143 225	8 177
15	7 905 853 580 625	16 368

- ▶ computational **costs** grow beyond all reasonable limits
- ▶ **gauge cancellations** cause loss of precision

Number of possible momenta in tree diagrams grows only **exponentially**

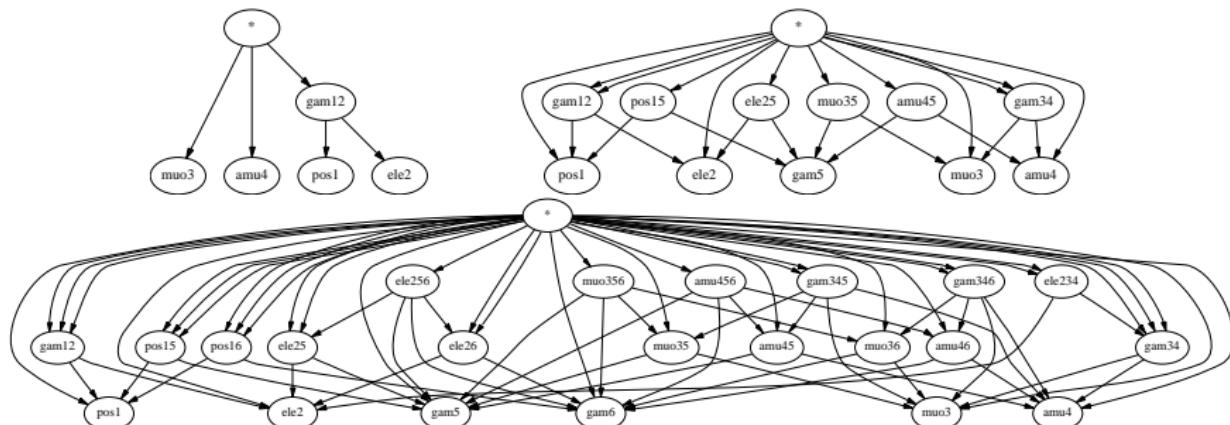
$$P(n) = \frac{2^n - 2}{2} - n = 2^{n-1} - n - 1$$

∴ Feynman diagrams **redundant** for many external particles!

∴ Replace the forest of tree diagrams by the **Directed Acyclical Graph (DAG)** of the algebraic expression.

$$ab(ab + c) = \begin{array}{c} \times \\ / \quad \backslash \\ a \quad b \end{array} + \begin{array}{c} \times \\ / \quad \backslash \\ a \quad b \end{array} c = \begin{array}{c} \times \\ \diagup \quad \diagdown \\ a \quad b \end{array} + \begin{array}{c} \times \\ \diagup \quad \diagdown \\ a \quad b \end{array} c$$

- simplest examples: $e^+e^- \rightarrow \mu^+\mu^-$, $e^+e^- \rightarrow \mu^+\mu^-\gamma$ and $e^+e^- \rightarrow \mu^+\mu^-\gamma\gamma$ (only QED)



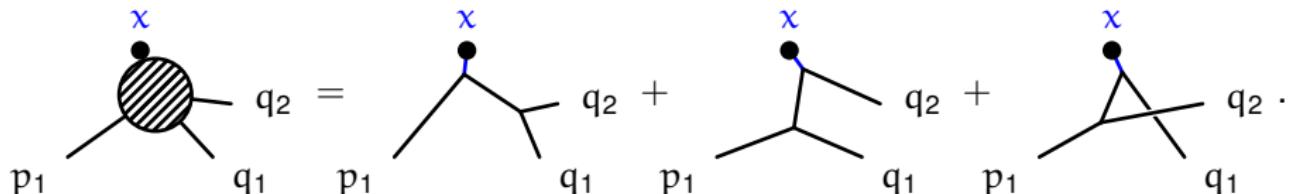
Efficient tree amplitudes

- ▶ Berends-Giele Recursion Relations [Berends, Giele]
 - ▶ manual calculations
- ▶ HELAS [Hagiwara et al.],
 - ▶ manual partial common subexpression elimination
- ▶ Madgraph [Stelzer et al.], AMEGIC++, COMIX [Krauss et al.]:
 - ▶ partial common subexpression elimination
 - ∴ **partial** elimination of redundancy
- ▶ ALPHA [Caravaglios & Moretti]:
 - ▶ tree level scattering amplitude is Legendre transform of Lagragian
 - ▶ can be performed **numerically**, using only $P^*(n)$ independent variables
- ▶ HELAC [Papadopoulos et al.]:
 - ▶ ALPHA algorithm can be reformulated as recursive **numerical** solution of Schwinger-Dyson equations
- ▶ O'Mega [TO et al.]:
 - ▶ systematic elimination of **all** redundancies
 - ▶ symbolic, generation of compilable code

One particle off-shell wave functions (**1POWs**) are obtained from by applying the LSZ reduction formula to all but one line:

$$W(x; p_1, \dots, p_n; q_1, \dots, q_m) = \\ \langle \phi(q_1), \dots, \phi(q_m); \text{out} | \Phi(x) | \phi(p_1), \dots, \phi(p_n); \text{in} \rangle .$$

E.g. $\langle \phi(q_1), \phi(q_2); \text{out} | \Phi(x) | \phi(p_1); \text{in} \rangle$ in ϕ^3 -theory at tree level



- the set of **all** 1POWs at tree level grows **exponentially** and can be constructed **recursively** from other 1POWs at tree level.

There exists a well defined set of **keystones K** that allow to express the sum of **Feynman diagrams** through **1POWs**:

$$T = \sum_{i=1}^{F(n)} D_i = \sum_{k,l,m=1}^{P(n)} K_{f_k f_l f_m}^3(p_k, p_l, p_m) W_{f_k}(p_k) W_{f_l}(p_l) W_{f_m}(p_m)$$

Even for vector particles, the 1POWs are ‘almost’ physical objects and satisfy simple **Ward Identities** in unbroken gauge theories

$$\frac{\partial}{\partial x_\mu} \langle \text{out} | A_\mu(x) | \text{in} \rangle_{\text{amp.}} = 0$$

and in spontaneously gauge theories in R_ξ -gauge

$$\frac{\partial}{\partial x_\mu} \langle \text{out} | W_\mu(x) | \text{in} \rangle_{\text{amp.}} = \xi_W m_W \langle \text{out} | \phi_W(x) | \text{in} \rangle_{\text{amp.}} .$$

- ▶ code for matrix elements can optionally be instrumented to check these Ward identities, testing the **consistency** a particular model and the **numerical stability** of expressions.

Amplitudes can be continued off-shell:

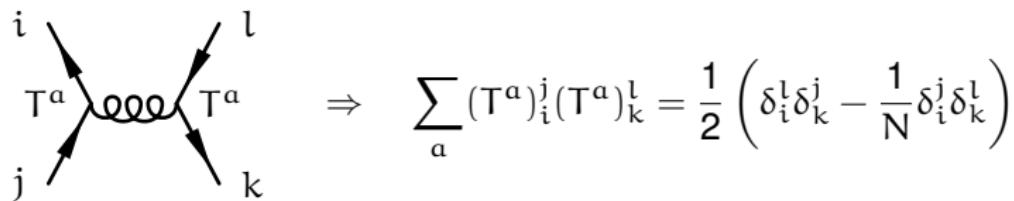
- ▶ **Slavnov-Taylor Identities** can be checked numerically by adding **operator insertions** implementing BRS transformations.

Slightly simplified Model.T signature that **all** models must implement:

```
module type Model.T =
sig
  type flavor (* all quantum numbers *)
  val flavor_symbol : flavor -> string
  val conjugate : flavor -> flavor (* antiparticles *)
  val lorentz : flavor -> Coupling.lorentz (* spin *)
  val fermion : flavor -> int (* fermion, boson, antifermion *)
  val width : flavor -> Coupling.width (* scheme, not value! *)
  type gauge (* parametrized gauges *)
  val gauge_symbol : gauge -> string
  val propagator : flavor -> gauge Coupling.propagator
  type constant (* coupling constants *)
  val constant_symbol : constant -> string
  val fuse2 : flavor -> flavor ->
    (flavor * constant Coupling.t) list (*  $A_\mu(p_{12}) \leftarrow g\bar{\Psi}(p_1)\gamma_\mu\Psi(p_2)$  *)
  val fuse3 : flavor -> flavor -> flavor ->
    (flavor * constant Coupling.t) list (*  $\phi(p_{123}) \leftarrow g\phi(p_1)\phi(p_2)\phi(p_3)$  *)
  val fuse : flavor list -> (flavor * constant Coupling.t) list
end
```

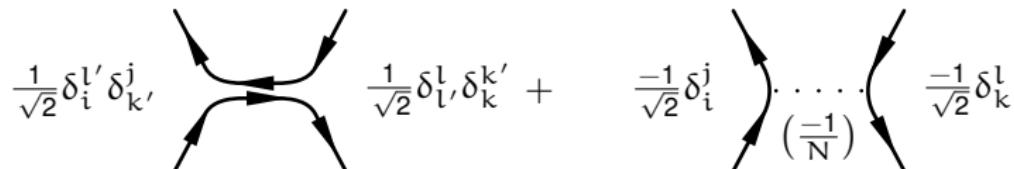
- For interfacing to parton shower and hadronization programs, we need the amplitudes for all possible color flows or color connections.

E.g. in $q\bar{q} \rightarrow q\bar{q}$



$$\sum_a (T^a)_i^j (T^a)_k^l = \frac{1}{2} \left(\delta_i^l \delta_j^k - \frac{1}{N} \delta_i^j \delta_k^l \right)$$

- This can be expressed by two diagrams, one for gluon and one for phantom exchange



$$\frac{1}{\sqrt{2}} \delta_i^{l'} \delta_{k'}^j + \frac{1}{\sqrt{2}} \delta_l^l \delta_{k'}^{k'} + \left(\frac{-1}{\sqrt{2}} \delta_i^j \right) \cdot \dots \cdot \left(\frac{-1}{\sqrt{2}} \delta_k^l \right)$$

- the sum over all colors can be written as a sum over all color flows

$$\sum_{IJ} N_C^{\lambda(J,J)} A_I A_J^*$$

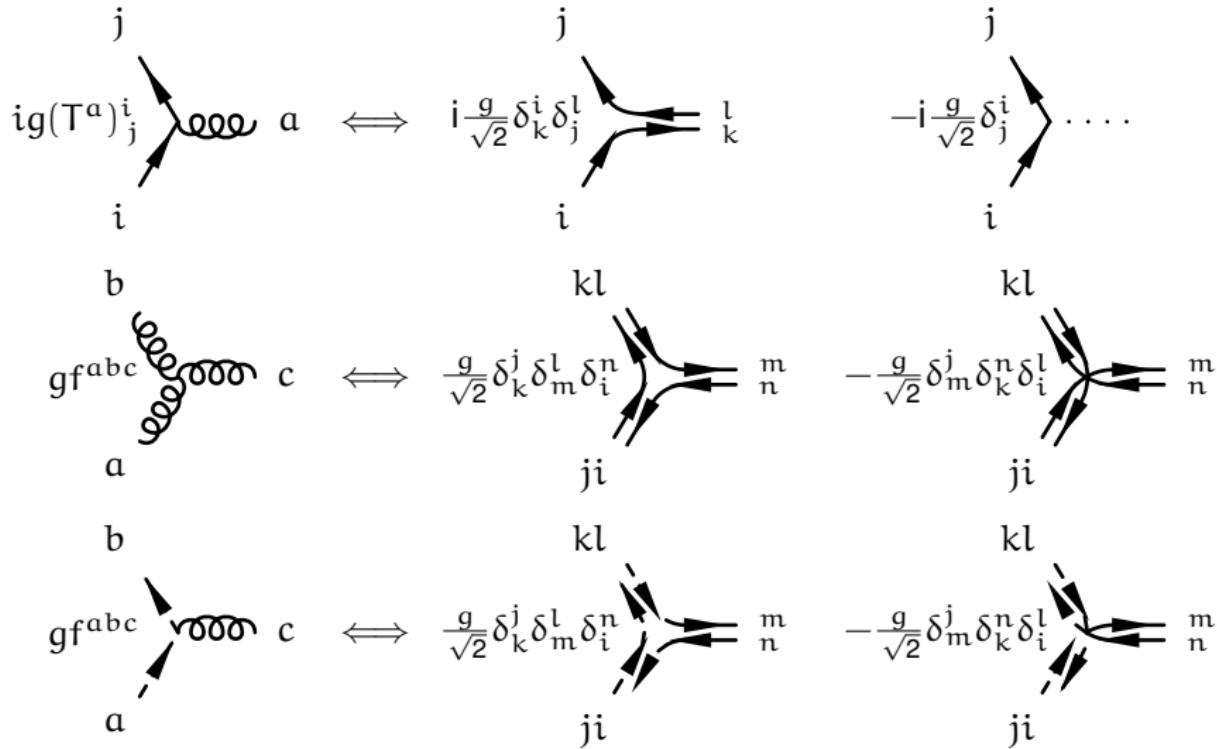
with $\lambda(J,J)$ the number of closed color loops in $A_I A_J^*$.

- one can use the completeness relation $2T^a \otimes T^a = \delta \otimes \delta - \delta \otimes \delta$ **repeatedly** to compute all the color flow amplitudes
- instead, one can formulate **equivalent** Feynman rules that give the color flow amplitudes **directly**
- Propagators**

$$\begin{array}{c}
 a \text{ } \overbrace{\hspace{1cm}}_{-i\delta^{ab}} \text{ } b \quad \Longleftrightarrow \quad j \text{ } \overbrace{\hspace{1cm}}_{-i\delta_k^j \delta_l^l} \text{ } k \quad + \quad \dots \dots \dots \\
 \\
 a \text{ } \overbrace{\hspace{1cm}}_{-i\delta^{ab}} \text{ } b \quad \Longleftrightarrow \quad j \text{ } \overbrace{\hspace{1cm}}_{-i\delta_k^j \delta_l^l} \text{ } k \\
 \\
 i \text{ } \overrightarrow{\hspace{1cm}}_{i\delta_j^i} \text{ } j \quad \Longleftrightarrow \quad j \text{ } \overrightarrow{\hspace{1cm}}_{i\delta_j^i} \text{ } i
 \end{array}$$

- The price to pay in the introduction of **phantom** particles that subtract the trace part of the gluons
- NB:** they're **not** required for the **Faddeev-Popov ghosts**, because the trace of the gluons behaves as if it was abelian.

► Cubic vertices:

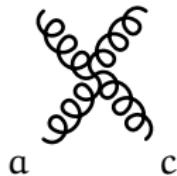


► Quartic vertices:

b

d

$$g^2(f^{abe}f^{cde} + f^{ace}f^{dbe} + f^{ade}f^{bce})$$



kl

mn

ji

op



$$\frac{g^2}{2} \delta_k^j \delta_m^l \delta_p^n \delta_i^o$$



$$\frac{g^2}{2} \delta_k^j \delta_p^l \delta_i^n \delta_m^o$$



$$\frac{g^2}{2} \delta_m^j \delta_p^l \delta_k^n \delta_i^o$$



$$\frac{g^2}{2} \delta_p^j \delta_m^l \delta_i^n \delta_k^o$$

kl

mn

ji

op

kl

mn

ji

op

kl

mn

ji

op



$$\frac{g^2}{2} \delta_p^j \delta_i^l \delta_k^n \delta_m^o$$



$$\frac{g^2}{2} \delta_m^j \delta_i^l \delta_p^n \delta_k^o$$



$$\frac{g^2}{2} \delta_p^j \delta_m^l \delta_i^n \delta_k^o$$



$$\frac{g^2}{2} \delta_i^j \delta_m^l \delta_k^n \delta_p^o$$

Example from **supersymmetry**: electroproduction of chargino pairs with bremsstrahlung, i.e. $e^+ e^- \rightarrow \tilde{\chi}_1^+ \tilde{\chi}_1^- \gamma$:

```
pure function l1b11cp1cm1a (k, s) result (amp)
  real(kind=omega_prec), dimension(0::), intent(in) :: k
  integer, dimension(:, :), intent(in) :: s
  complex(kind=omega_prec) :: amp
  type(momentum) :: p1, p2, p3, p4, p5
  type(bispinor) :: cp1_4, l1_2
  type(bispinor) :: cm1_3, l1b_1
  type(vector) :: a_5
  complex(kind=omega_prec) :: snl_24, snc1_13
  type(bispinor) :: cp1_45, l1_25
  type(bispinor) :: cm1_35, l1b_15
  type(vector) :: a_34, a_12, z_34, z_12
  type(momentum) :: p12, p13, p15, p24, p25, p34, p35, p45
  p1 = - k(:, 1) ! incoming e+
  p2 = - k(:, 2) ! incoming e-
  p3 = k(:, 3) ! outgoing ch1+
  p4 = k(:, 4) ! outgoing ch1-
  p5 = k(:, 5) ! outgoing A
  l1b_1 = u (mass(11), - p1, s(1))
  l1_2 = u (mass(11), - p2, s(2))
  cm1_3 = v (mass(69), p3, s(3))
  cp1_4 = v (mass(69), p4, s(4))
  a_5 = conjg (eps (mass(22), p5, s(5)))
  p12 = p1 + p2
  a_12 = pr_feynman(p12, + v_ff(qlep,l1b_1,l1_2))
  z_12 = pr_unitarity(p12,mass(23),wd_t1(p12,width(23)), &
    + va_ff(gncllep(1),gncllep(2),l1b_1,l1_2))
  p13 = p1 + p3
  snc1_13 = pr_phi(p13,mass(54),wd_t1(p13,width(54)), &
    + sr_ff(g_yuk_ch1_snl_1_c,l1b_1,cm1_3))
  p24 = p2 + p4
```

```
sn1_24 = pr_phi(p24,mass(54),wd_tl(p24,width(54)), &
+ sl_ff(g_yuk_ch1_sn1_1,11_2,cp1_4))
p34 = p3 + p4
a_34 = pr_feynman(p34, + v_ff(qchar,cm1_3,cp1_4))
z_34 = pr_unitarity(p34,mass(23),wd_tl(p34,width(23)), &
+ va_ff(-gczc_1_1(1),-gczc_1_1(2),cm1_3,cp1_4))
p15 = p1 + p5
l1b_15 = pr_psi(p15,mass(11),wd_tl(p15,width(11)), + f_vf(-qlep,a_5,l1b_1))
p25 = p2 + p5
l1_25 = pr_psi(p25,mass(11),wd_tl(p25,width(11)), + f_vf(qlep,a_5,l1_2))
p35 = p3 + p5
cm1_35 = pr_psi(p35,mass(69),wd_tl(p35,width(69)), &
+ f_vf(-qchar,a_5,cm1_3))
p45 = p4 + p5
cp1_45 = pr_psi(p45,mass(69),wd_tl(p45,width(69)), + f_vf(qchar,a_5,cp1_4))
amp = 0
amp = amp + sn1_24*( + sr_ff(g_yuk_ch1_sn1_1_c,cm1_35,l1b_1))
amp = amp + snc1_13*(- sl_ff(g_yuk_ch1_sn1_1,11_25,cp1_4) &
+ sl_ff(g_yuk_ch1_sn1_1,cp1_45,l1_2))
amp = amp + l1_25*(- f_vf(-qlep,a_34,l1b_1) &
- f_vaf(-(gncalep(1)),gncalep(2),z_34,l1b_1))
amp = amp + l1b_15*(- f_srf(g_yuk_ch1_sn1_1_c,sn1_24,cm1_3) &
+ f_vf(qlep,a_34,l1_2) + f_vaf(gncalep(1),gncalep(2),z_34,l1_2))
amp = amp + z_12*(- va_ff(-(-gczc_1_1(1)), -gczc_1_1(2),cp1_45,cm1_3) &
+ va_ff(-gczc_1_1(1),-gczc_1_1(2),cm1_35,cp1_4))
amp = amp + a_12*(- v_ff(-qchar,cp1_45,cm1_3) + v_ff(qchar,cm1_35,cp1_4))
end function l1b1cp1cm1a
```

28 fusions, 10 propagators, 12 diagrams

- ▶ readable code, can be edited for **exotic models** or **NLO vertex functions**

Remaining problem:

$$I(f) = \int_M d\mu(p) f(p)$$

1. non-trivial geometry of multi particle phase space

$$d\mu(p) = \delta^4(\sum_n k_n - P) \prod_n d^4 k_n \delta(k_n^2 - m_n^2)$$

2. ill-behaved function, i. e. squared matrix element w/ kinematical cuts

$$f(p) = |T(k_1, k_2, \dots)|^2 \cdot C(k_1, k_2, \dots)$$

Choose a (pseudo-)random sequence $p_g = \{p_1, p_2, \dots, p_N\}$ distributed according to $d\mu_g(p) = g(p)d\mu(p)$, then an estimator of $I(f)$ is

$$E(f) = \left\langle \frac{f}{g} \right\rangle_g = \frac{1}{N} \sum_{i=1}^N \frac{f(p_i)}{g(p_i)}$$

The sampling error is estimated by the square root of the variance

$$V(f, g) = \frac{1}{N-1} \left(\left\langle \left(\frac{f}{g} \right)^2 \right\rangle_g - \left\langle \frac{f}{g} \right\rangle_g^2 \right)$$

which depends on g , even after averaging over p .

Conflicting goals for g

1. make $d\mu_g(p)$ simple enough, so that p_g can be generated w/ reasonable effort
 2. choose g to minimize $V(f, g)$: importance sampling or stratified sampling
- ▶ for multi particle phase space, $d\mu(p)$ is very intricate and the generation of p_g is not trivial even for $g(p) = 1/\text{Vol}(M)$.
- ∴ RAMBO: elegant trick only for $m_n = 0$ and g constant
- ∴ parametrizations $]0, 1[^{\otimes \dim(M)} \rightarrow M$: require compensation of bad Jacobians

Practical considerations for particle physics:

∴ only a small number of different manifolds M :

- number of particles 2, 3, 4, 5, 6, 7, ...
- massless vs. massive particles

∴ it makes sense to invest manpower into an optimal treatment of the geometry, i. e. $d\mu$

∴ f changes w/

- physics model du jour
- physical parameters in the model
- changing external cuts that can affect singular regions

∴ automatic and computer aided adaptive approaches, i. e. numerical optimizations, are appropriate

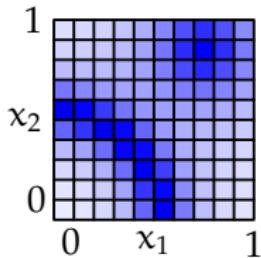
For over a quarter century, Peter Lepage's VEGAS has been the workhorse for adaptive Monte Carlo in particle physics.

For simplicity

$$x \in M =]0, 1[^{\otimes n}, \quad d\mu(x) = dx_1 \wedge dx_2 \wedge \dots \wedge dx_n$$

How can we implement efficiently a variable weight g in $d\mu_g(x) = g(x)d\mu(x)$?

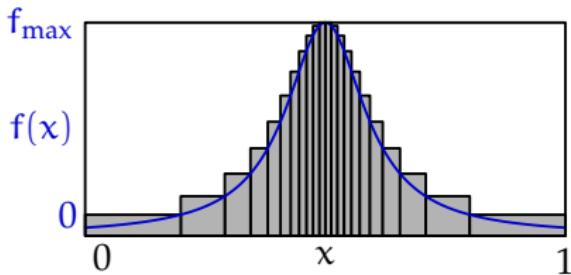
- ▶ optimization of expansion coefficients α in $g(x) = \sum_l \alpha_l g_l(x)$ popular, but not exciting for **generator generation**
 - ∴ selection of g_l requires expert human input
 - ∴ can't deal very efficiently w/ cuts
- ▶ fixed grid w/ variable weights



x (i. e. characteristic functions as g_l)
not useful at all

∴ locally fixed resolution can **not** adapt to the typical power law singularities over orders of magnitude

- ▶ alternative in **one** dimension: instead of adjusting weights of fixed bins, adjust density of equal weight bins
 - ∴ globally fixed resolution can nevertheless adjust locally over many orders of magnitude:



iteratively adjust grid, use estimates to either

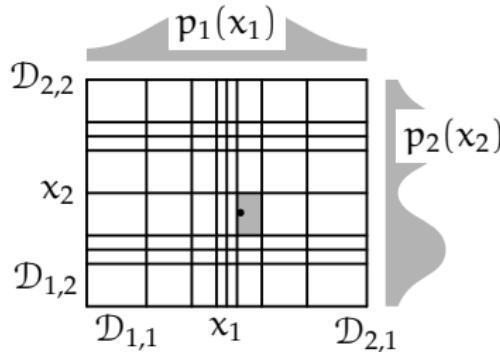
- ▶ approximate **f** locally (importance sampling \Rightarrow event generation)
- ▶ or equidistribute the variance (stratified sampling \Rightarrow high precision integration).

Factorized ansatz

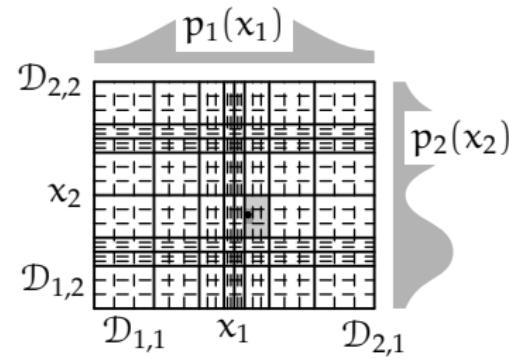
$$g(x) = g_1(x_1)g_2(x_2) \cdots g_n(x_n)$$

- guarantees hypercubic properties and simple one-dimensional formulae (w/ averaging over other dimensions)
- computational costs grow only **linearly** w/ the number of dimensions

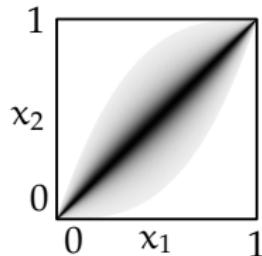
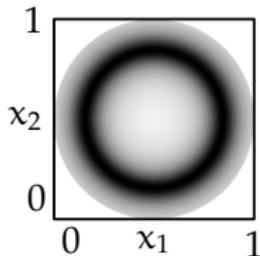
VEGAS grid structure for **importance sampling**:



for genuinely **stratified** sampling, used in low dimensions:

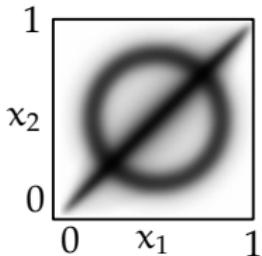


VEGAS' factorized ansatz handles



separately after mappings.

- ▶ fails for overlapping singularities



which is the common case
(for more than one diagram)

∴ adaptive multichannel approach

$$I(f) = \int_M d\mu(p) f(p)$$

$$I(f) = \sum_{i=1}^{N_c} \alpha_i \int_0^1 g_i(x) d^n x \frac{f(\phi_i(x))}{g(\phi_i(x))}$$

with

$$g = \sum_{i=1}^{N_c} \alpha_i \cdot (g_i \circ \phi_i^{-1}) \left| \frac{\partial \phi_i^{-1}}{\partial p} \right|$$

- ▶ works with **factorized** g_i adapted by **VEGAS** and α_i adapted by variance reduction.

- ▶ in general, $g \circ \phi_i$ does not factorize, even if all g_i factorize.
- ▶ $\pi_{ij} = \phi_j^{-1} \circ \phi_i$: coordinate transformations among coordinate systems in which different singularities factorize.
- ▶ pure geometry: economical studies of dependence on cuts and parameters
 - :: π_{ij} universal and are calculated automatically by WHIZARD
 - :: VEGAS can optimize the g_i for each set of parameters and cuts

However:

- :: singularity structure determined by Feynman diagrams
- ▶ naive application brings the combinatorial explosion in through the back door!
- ∴ WHIZARD uses heuristics to select the important channels
 - ▶ s-channel resonances
 - ▶ 1/t-poles for massless particlesand permutation symmetries to eliminate equivalent channels

- In Monte Carlo **integration** of integrals of observables f on phasespace weighted with the cross section

$$\Sigma(f) = \int d\sigma(\Phi) f(\Phi) = \int d\Phi \frac{d\sigma}{d\Phi}(\Phi) f(\Phi)$$

it suffices to generate **weighted** phase space configurations

$$\mathcal{W} = \{(\phi_i, w_i)\}_{i \in \mathbb{N}}$$

such that the integral is approximated by the **weighted sum**

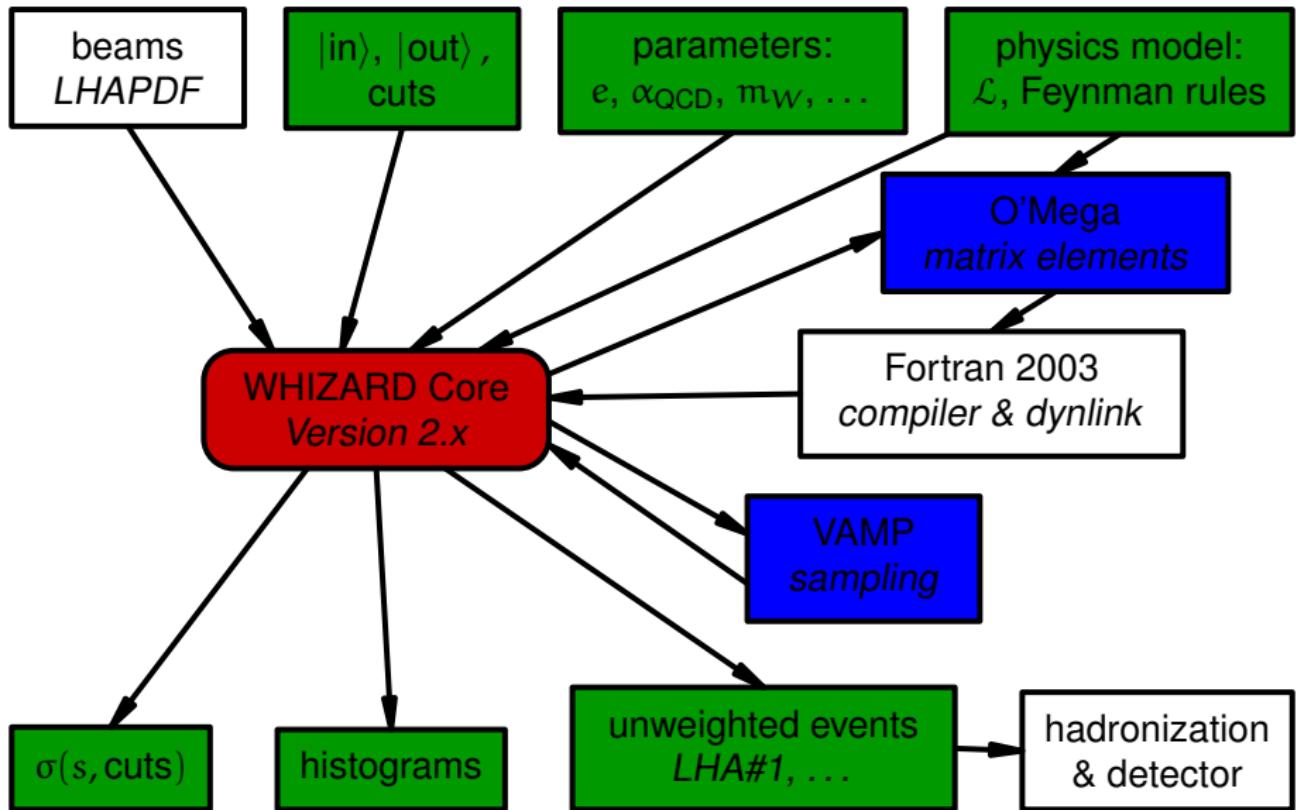
$$\Sigma(f) \approx \sum_{i \in \mathbb{N}} w_i f(\phi_i).$$

- In the case of wildly varying cross sections, this is often **much simpler** than generating **unweighted** phase space configurations

$\mathcal{U} = \{\phi_i\}_{i \in \mathbb{N}}$ with

$$\Sigma(f) \approx \sum_{i \in \mathbb{N}} f(\phi_i),$$

- ... by they **are** required for realistic **detector simulations**.





- ▶ the **names** of particles and couplings depend on the model and can be looked up in the share/whizard/models directory, e.g. in

```
/home/HEP/toolbox-1.1.7/whizard/share/whizard/models/SM.mdl  
for the standard model
```

- ▶ **input parameters**

```
parameter GF      = 1.16639E-5    # Fermi constant  
parameter mZ      = 91.1882       # Z-boson mass  
parameter mW      = 80.419        # W-boson mass  
parameter mH      = 125           # Higgs mass  
parameter alphas = 0.1178        # Strong coupling constant (Z point)  
parameter me      = 0.000510997   # electron mass
```

```
...
```

- ▶ **derived parameters**

```
derived v      = 1 / sqrt (sqrt (2.) * GF)      # v (Higgs vev)  
derived cw     = mW / mZ                         # cos(theta-W)
```

```
...
```

- ▶ **particles**

```
particle D_QUARK 1 parton  
  spin 1/2 charge -1/3  isospin -1/2  color 3  
  name d down  
  anti dbar D "d~"  
  tex_anti "\bar{d}"
```

► WHIZARD input `xsect_eemm.sin`

```
model = SM
process mumu = e1, E1 => e2, E2
compile
sqrtS = 90 GeV
beams = e1, E1
integrate (mumu) { iterations = 2:1000, 3:5000 }
```

► run WHIZARD

```
$ /home/HEP/toolbox-1.1.7/whizard/bin/whizard xsect_eemm.sin
```

► console output

```
| Integrating process 'mumu':
|=====
| It    Calls  Integral[fb]  Error[fb]   Err[%]   Acc  Eff[%]  Chi2 N[It] |
|=====
| 1      1000  1.0605313E+06  8.45E+02   0.08   0.03*  60.96
| 2      1000  1.0597762E+06  5.35E+02   0.05   0.02*  61.25
| -----
| 2      2000  1.0599922E+06  4.52E+02   0.04   0.02   61.25   0.57  2
| -----
| 3      5000  1.0601102E+06  1.34E+02   0.01   0.01*  61.12
| 4      5000  1.0599916E+06  9.36E+01   0.01   0.01*  78.78
| 5      5000  1.0598832E+06  7.58E+01   0.01   0.01*  67.98
| -----
| 5      15000 1.0599559E+06  5.39E+01   0.01   0.01   67.98   1.19  3
| -----
|=====
```

5 15000 1.0599559E+06 5.39E+01 0.01 0.01 67.98 1.19 3

► WHIZARD input

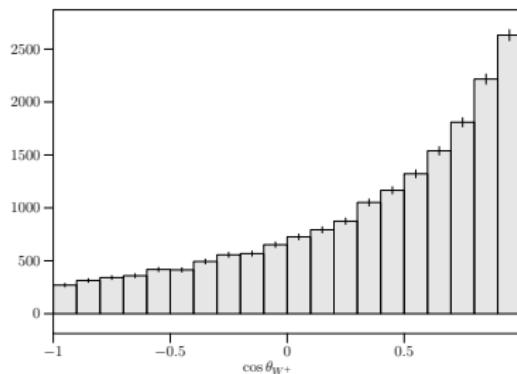
```
model = SM
alias parton = u:U:d:D:g
process tt = parton, parton => b, Wp, B, Wm
compile
sqrtS = 8 TeV
beams = p, p => pdf_builtin
cuts = all Pt > 10 GeV [b:Wp:B:Wm]
integrate (tt) { iterations = 2:10000, 5:20000 }
```

► console output

```
[1/5] gl gl -> b W+ bbar W- ... allowed.
[2/5] d dbar -> b W+ bbar W- ... allowed.
[3/5] dbar d -> b W+ bbar W- ... allowed.
[4/5] u ubar -> b W+ bbar W- ... allowed.
[5/5] ubar u -> b W+ bbar W- ... allowed.
=====
| It      Calls  Integral[fb]  Error[fb]   Err[%]    Acc  Eff[%]  Chi2 N[It] |
|=====|
 1      10000  9.9611255E+04  2.89E+04  29.06*   29.06*  2.02
 2      10000  1.1805759E+05  1.65E+04  13.95*   13.95*  2.63
|-----|
 2      20000  1.1354494E+05  1.43E+04  12.61    17.83    2.63   0.31  2
|-----|
 3      20000  1.1757383E+05  1.64E+04  13.91    19.67    1.45
 4      20000  1.2032373E+05  8.96E+03  7.45     10.53*   2.32
 5      20000  1.1948028E+05  5.69E+03  4.76     6.73*   3.09
 6      20000  1.2279117E+05  4.20E+03  3.42     4.84*   3.86
 7      20000  1.2270293E+05  3.64E+03  2.97     4.20*   3.42
|-----|
 7      100000  1.2190186E+05  2.36E+03  1.94     6.13     3.42   0.09  5
|-----|
 7      100000  1.2190186E+05  2.36E+03  1.94     6.13     3.42   0.09  5
|=====|
 | time estimate for generating 10000 unweighted events: 14m:18s
```

```
▶ model = SM
process ww = e1, E1 => Wp, Wm
compile
$x_label = "$\cos\theta_{W^+}$"
?draw_errors = true
histogram costh (-1, 1, 0.1)
sqrtS = 180 GeV
beams = e1, E1
luminosity = 1000 / 1 pbarn
integrate (ww) { iterations = 2:1000, 5:50000 }
simulate (ww) { analysis = record costh (eval cos (Theta) [Wm]) }
compile_analysis { $out_file = "ww.dat" }
```

▶



$pp \rightarrow gg \rightarrow \mu^-\bar{\nu}_e u \bar{d}$ as a **Les Houches Event File (LHEF)**:

```
model = SM
process hWW = g, g => e2, N2, u, D
compile
sqrtS = 14 TeV
beams = p, p => pdf_builtin
sample_format = lhef
simulate (hWW) { n_events = 1 }
```

```
<LesHouchesEvents version="1.0">
<header>
  <generator_name>WHIZARD</generator_name>
  <generator_version>2.1.1</generator_version>
</header>
<init>
  2212    2212  7000.0000000000000000  7000.0000000000000000      -1          -1          -1          -1
  420.46515665414375  249.25905165127301  1.0000000000000000
</init>
<event>
  6      1  1.0000000000000000  136.31643354892731      -1.0000000000000000  0.11780000000000000
  21     -1      0      0      501      502  0.0000000000000000  0.0000000000000000  64.59551
  21     -1      0      0      503      501  0.0000000000000000  0.0000000000000000  -71.91741
  13      1      1      2      0      0  16.068546922717953  -15.231242240310792  28.41895
  -14     1      1      2      0      0  -35.486359371670360  14.938719772529632  -38.73337
   2      1      1      2      503      0  14.687109514917077  13.162760335459950  -13.89070
  -1      1      1      2      0      502  4.7307029340353255  -12.870237867678787  16.88323
</event>
</LesHouchesEvents>
```

$pp \rightarrow gg \rightarrow \mu^-\bar{\nu}_e u\bar{d}$ as a **HepMC** file:

```
model = SM
process hWW = g, g => e2, N2, u, D
compile
sqrtS = 14 TeV
beams = p, p => pdf_builtin
sample_format = hepmc
simulate (hWW) { n_events = 1 }
```

```
HepMC::Version 2.06.09
HepMC::IO_GenEvent-START_EVENT_LISTING
E 1 -1 1.3631643354892731e+02 1.1780000000000000e-01 -1.0000000000000000e+00 1 0 3 10001 10002 0 4 1.0000000000000000e+00 7.08284436471
N 4 "0" "1" "2" "3"
U GEV MM
C 4.2046515665414375e+02 2.4925905165127301e+02
V -1 0 0 0 0 1 2 0
P 10001 2212 0 0 7.0000000000000000e+03 7.0000000000000000e+03 0 4 0 0 -1 0
P 10003 21 0 0 6.4595518178633455e+01 6.4595518178633455e+01 0 3 0 0 -3 2 1 1 2 2
P 10005 93 0 0 6.9354044818213670e+03 6.9354044818213670e+03 0 3 0 0 0 2 1 2 2 1
V -2 0 0 0 0 1 2 0
P 10002 2212 0 0 -7.0000000000000000e+03 7.0000000000000000e+03 0 4 0 0 -2 0
P 10004 21 0 0 -7.1917412304487158e+01 7.1917412304487158e+01 0 3 0 0 -3 2 1 3 2 1
P 10006 93 0 0 -6.9280825876955132e+03 6.9280825876955132e+03 0 3 0 0 0 2 1 1 2 3
V -3 0 0 0 0 0 4 0
P 10007 13 1.6068546922717953e+01 -1.5231242240310792e+01 2.8418956865688791e+01 3.6025507816263499e+01 1.0565838899979969e-01 1 0 0 0
P 10008 -14 -3.5486359371670360e+01 1.4938719772529632e+01 -3.8733375441153328e+01 5.4614296873280757e+01 0 1 0 0 0 0
P 10009 2 1.4687109514917077e+01 1.3162760335459950e+01 -1.3890707675246947e+01 2.4123043035054472e+01 4.7683715820312500e-07 1 0 0 0 1
P 10010 -1 4.7307029340353255e+00 -1.2870237867678787e+01 1.6883232124857773e+01 2.1750082758521934e+01 4.1295309247228556e-07 1 0 0 0
HepMC::IO_GenEvent-END_EVENT_LISTING
```

```

model = SM
process hWW = g, g => e2, N2, u, D
compile
sqrtS = 14 TeV
beams = p, p => lhapdf
sample_format = hepmc, lhef
integrate (hWW) { iterations = 1:1000 }
?ps_fsr_active = true
?ps_isr_active = true
?hadronization_active = true
?ps_use_PYTHIA_shower = true
simulate (hWW) { n_events = 1 }

<event>

```

310	1	1.0000000000000000	103.17887552082534	-1.0000000000000000	0.11780000000000000	
21	-1	0	0	501	502	0.0000000000000000
21	-1	0	0	502	503	0.0000000000000000
93	3	0	0	502	501	0.0000000000000000
93	3	0	0	503	502	0.0000000000000000
-14	1	1	2	0	0	-27.571846966921186
13	1	1	2	0	0	8.2564499233431601
22	1	1	2	0	0	-0.60053536107337491
211	1	1	2	0	0	0.18025914014147792
-211	1	1	2	0	0	-0.83369544321947398
2112	1	1	2	0	0	-0.34678766523372612
-2212	1	1	2	0	0	-7.7893814158812913E-002
2212	1	1	2	0	0	0.46265670377713941
-2212	1	1	2	0	0	0.71417103301686868
211	1	1	2	0	0	2.7762381422163100
-211	1	1	2	0	0	0.6556572660978855
211	1	1	2	0	0	1.1650307289886463
-211	1	1	2	0	0	0.10194216370164123
-321	1	1	2	0	0	1.2505517529312573



- ▶ Compute the cross section for **single top production**

$$pp \rightarrow tb$$

in the standard model at the Tevatron and at the 14 TeV LHC.

- ▶ Plot the rapidity distribution of the produced top quark.
- ▶ Include the two-particle decay of the top quark.
- ▶ Include the irreducible backgrounds.
- ▶ Estimate the influence of bottom-quark parton distribution functions.