

# Calculating $gg \rightarrow t\bar{t} + jets$ at Tree Level

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DESY Summer Students Program 2010

# Motivation

- We are interested in  $gg \rightarrow t\bar{t} + n \text{ gluons}$  at tree level.
- This is a partonic part of  $pp \rightarrow t\bar{t} + \text{jets}$ .
- This is a background for Higgs searches at the LHC.

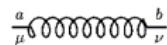
# QCD Feynman Rules

- Draw the Feynman diagrams of the process
- Label each line with a momentum
- Associate particular structures as follows:

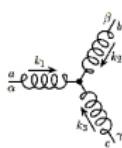
Fermion propagator :  $\frac{i(\not{k}+m)\delta_i^j\delta_{i'}^{j'}}{\not{k}^2-m^2+i\epsilon}$



Gluon propagator :  $\frac{-i\partial_\mu^b g_{\mu\nu}}{k^2+i\epsilon}$

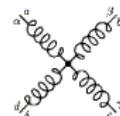


Three gluons vertex :  $-gf^{abc} [g^{\alpha\beta}(k_1-k_2)^\gamma + g^{\beta\gamma}(k_2-k_3)^\alpha + g^{\gamma\alpha}(k_3-k_1)^\beta]$

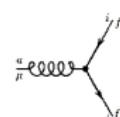


Four gluons vertex :

$$-ig^2 [f^{abe} f^{cde} (g^{\alpha\gamma} g^{\beta\delta} - g^{\alpha\delta} g^{\beta\gamma}) + f^{ace} f^{bde} (g^{\alpha\beta} g^{\gamma\delta} - g^{\alpha\delta} g^{\beta\gamma}) + f^{ade} f^{bce} (g^{\alpha\beta} g^{\delta\gamma} - g^{\alpha\gamma} g^{\delta\beta})]$$



Fermion vertex :  $i g \gamma^\mu \delta_{f'}^f (t^a)_i^j$



# QCD Feynman Rules

- Associate the external particles as follows:

*Initial external fermion :  $u(k, s)$*

*Final external fermion :  $\bar{u}(k, s)$*

*Initial external antifermion :  $\bar{v}(k, s)$*

*Final external antifermion :  $v(k, s)$*

*Initial external gluon :  $\epsilon_\mu(k)$*

*Final external gluon :  $\epsilon_\mu^*(k)$*

- Sum every term of diagrams and sum over colour, polarization, and spin

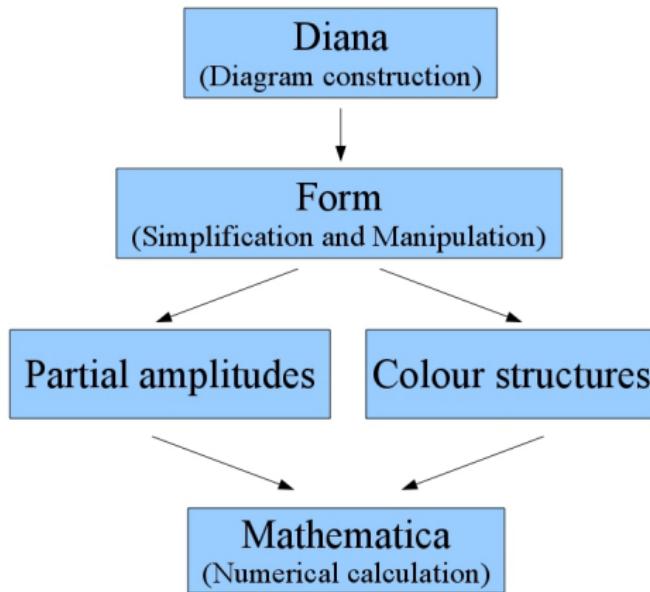
$gg \rightarrow t\bar{t} + n \text{ gluons}$

- The number of diagrams is increased rapidly when the number of outgoing gluons is increased

Process	$gg \rightarrow t\bar{t}$	$gg \rightarrow t\bar{t} + g$	$gg \rightarrow t\bar{t} + gg$
Number of diagrams	3	16	123

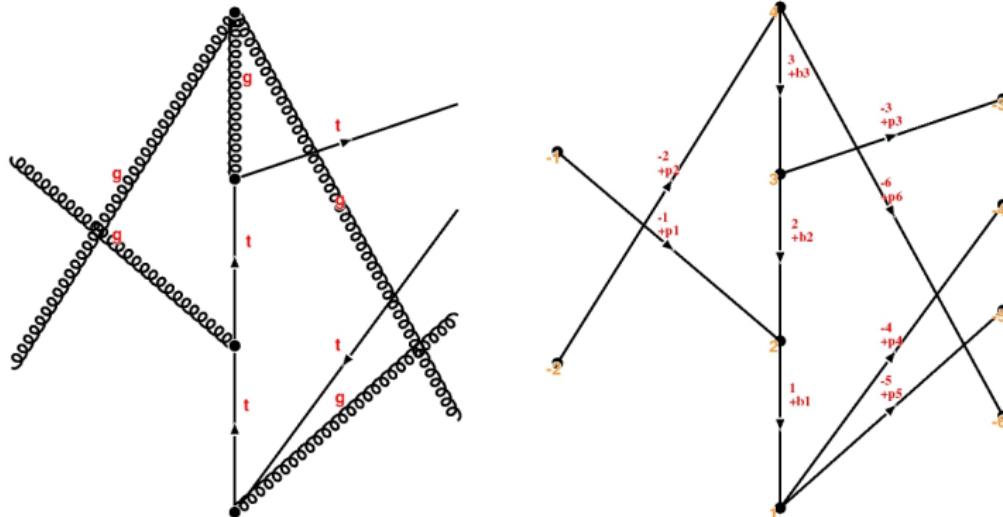
- The complete calculation can be done by using several programs in combination.

# The Method



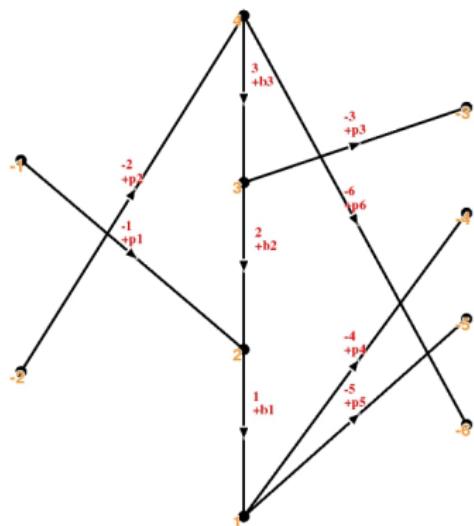
# Diagram Construction

- We ask Diana for the diagram construction.



# Example of a Diagram Contribution

- Feynman rules are defined as a set of functions in order to be used by Form.



```
#define LINE "3"
#define FERMIONLINE "1"
#define TOPOLOGY "i4_i082"
*****
#define b1 "(+p4+p5)"
#define b2 "(-p1+p4+p5)"
#define b3 "(-p1+p3+p4+p5)"
#define NM "3"
*****
I Rq =
1*F(3,1,li3,1,0,1)*i_ *gs*GM(ai3,fi3,fi1)*
FF(2,1,-b2,mt)*i_ *fdelta(fi1,fi)*
F(2,1,li1,1,0,1)*i_ *gs*GM(ai1,fi,fi2)*
FF(1,1,-b1,mt)*i_ *fdelta(fi2,fi5)*
F(1,1,li5,1,0,1)*i_ *gs*GM(ai5,fi5,fi4)*
VV(3,li,li3,+b3,3)*i_ *adelta(ai,ai3)*
V(4,li6,li2,li,-p6,+p2,-b3,3)*gs*Fabc(ai6,ai2,ai);
```

# Colour Structure and Partial Amplitude

- The amplitude can be divided in two parts:

$$M = \sum_i c_i M_i^{\text{partial}}$$

- Colour structures and partial amplitudes are simplified and manipulated separately by Form.

# Colour Structure

- Form reads out the different colour structures, for example,

```
#define colfactor1 "T(ai1,ai2,fi3,fi4)"  
  
#define colfactor2 "T(ai2,ai1,fi3,fi4)"
```

where  $T(ai_1, ai_2, fi_3, fi_4) = (t^{ai_1} t^{ai_2})_{fi_3 fi_4}$

- By squaring the amplitude and sum over colour, we get
- $$\sum_{colour} |M|^2 = \sum_{colour} \sum_{i,j} c_i^* M_i^* c_j M_j = \sum_{i,j} \left( \sum_{colour} c_i^* c_j \right) M_i^* M_j$$
- The matrix  $\sum c_i^* c_j$  can be simplified by using SU(3) algebra and kept in Mathematica format.

```
matrix[1,1]:={NF^-1*a^2 - 2*NF*a^2 + NF^3*a^2}  
matrix[1,2]:={NF^-1*a^2 - NF*a^2}  
matrix[2,1]:={NF^-1*a^2 - NF*a^2}  
  
matrix[2,2]:={NF^-1*a^2 - 2*NF*a^2 + NF^3*a^2}
```

# Partial Amplitude

- The partial amplitude is simplified by using properties of Dirac matrices and Dirac equation, and the output is in Mathematica format.

```
re1e2Sum11=
+SpinorUBar(p3,mt)*GS(p2)*SpinorV(p4,mt)*I*p1dp2^-1*e1de2
+SpinorUBar(p3,mt)*GS(e1)*SpinorV(p4,mt)*I*p1dp2^-1*p1de2
-SpinorUBar(p3,mt)*GS(e2)*SpinorV(p4,mt)*I*p1dp2^-1*p2de1;
re1e2Sum12=
-SpinorUBar(p3,mt)*GS(p2)*SpinorV(p4,mt)*I*p1dp2^-1*e1de2
-SpinorUBar(p3,mt)*GS(e1)*SpinorV(p4,mt)*I*p1dp2^-1*p1de2

+SpinorUBar(p3,mt)*GS(e2)*SpinorV(p4,mt)*I*p1dp2^-1*p2de1 ;
```

# Numerical Calculation

- We use Mathematica for numerical calculation.
- The appropriate phase space point (the set of momenta,  $p_1, \dots, p_n$ ) is set at the beginning.
- The gluon polarization vector basis ( $\varepsilon_1, \dots, \varepsilon_m$ ) are chosen.
- The representation of Dirac matrices is defined.

# 4 Point Process: $gg \rightarrow t\bar{t}$

- The partial amplitudes are compared with previous calculation [6] (R. K. Ellis, W. T. Giele, Z. Kunszt, K. Melnikov. Nucl.Phys.B822:270-282, 2009)
- The phase space point is  
 $p_1 = E(1, -\sin\theta, 0, -\cos\theta)$ ,  $p_2 = E(1, \sin\theta, 0, \cos\theta)$ ,  $p_3 = E(1, 0, 0, \beta)$ ,  $p_4 = E(1, 0, 0, -\beta)$ , where  
 $m_t = 1.75$ ,  $E = 10$ ,  $\beta = \sqrt{1 - m_t^2/E^2}$ , and  $\theta = \pi/3$ .

Helicities	Partial amplitude (colour structure 1)	Primitive amplitude [6]
$+\bar{t}, +1, +2, +t$	0.0009048290295650407i	0.000905i

- The squared matrix element is compared with previous calculation[7] (W. Bernreuther, A. Brandenburg, Z. G. Si, P. Uwer. Nucl.Phys. B690 (2004) 81-137)
- The phase space point is  
 $p_1 = \frac{s}{2}(1, 0, 0, 1)$ ,  $p_2 = \frac{s}{2}(1, 0, 0, -1)$ ,  $p_3 = \frac{s}{2}(1, \beta \sin\theta \cos\phi, \beta \sin\theta \sin\phi, \beta \cos\theta)$ ,  $p_4 = p_1 + p_2 - p_3$ ,  
where  $\beta = \sqrt{1 - 4m_t^2/s}$ ,  $\cos\theta$  is the angle between incoming particle and outgoing particle,  $s$  is center of mass energy.

$s$	$\cos\theta$	Numerical result of the method	Analytical result of [7]
20	0.842497	413.9748159148358	413.9748159148358
200	0.90523	772.2986631763597	772.2986631763583

# 5 Point Process: $gg \rightarrow t\bar{t} + g$

- The partial amplitudes are compared with previous calculation [6] (R. K. Ellis, W. T. Giele, Z. Kunszt, K. Melnikov. Nucl.Phys.B822:270-282, 2009)
- The phase space point is  
 $p_1 = E\xi(-1, 1, 0, 0)$ ,  $p_2 = E\xi(-\sqrt{2}, 0, 1, 1)$ ,  $p_3 = E(1, 0, 0, \beta)$ ,  $p_4 = E(1, 0, 0, -\beta)$ ,  $p_5 = p_1 + p_2 - p_3 - p_4$ ,  
 where  $m_t = 1.75$ ,  $E = 10$ ,  $\beta = \sqrt{1 - m_t^2/E}$ , and  $\xi = 2/(1 + \sqrt{2} + \sqrt{3})$

Helicities	Partial amplitude (colour structure 1)	Primitive amplitude [6]
$+_{\bar{t}}, +_{t+1}, +_2, +_5$	$-0.0005332686176129279 - 0.00013689856022906747i$	$-0.000533-0.000137i$

- The squared matrix element is compared with previous calculation [8] (S. Dittmaier, P. Uwer, S. Weinzierl. arXiv:0810.0452, hep-ph)
- The phase space point is  $p_1 = (500, 0, 0, 500)$ ,  $p_2 = (500, 0, 0, -500)$ ,  
 $p_3 = (458.53317553852783, 207.0255169909440, 0, 370.2932732896167)$ ,  
 $p_4 = (206.6000026080000, -10.65693677252589, 42.52372780926147, -102.39982104210421085)$ ,  
 $p_5 = (334.8668220067217, -196.3685802184181, -42.52372780926147, -267.8934522475083)$ .

	Squared matrix element ( $10^{-3} \text{ GeV}^{-2}$ )
Result of [8] Version 1	0.6566843362709776
Numerical result of the method	0.6566843357688175

# 6 Point Process: $gg \rightarrow t\bar{t} + gg$

- $gg \rightarrow t\bar{t} + gg$  compared with Madgraph.
- We use the benchmark phase space point [8] (S. Dittmaier, P. Uwer, S. Weinzierl. arXiv:0810.0452, hep-ph):
- $p_1 = (2100, 0, 0, 2100)$ ,  $p_2 = (2800, 0, 0, -2800)$ ,  
 $p_3 = (1581.118367308447, 1254.462316247655, -766.9360998604944, -554.7905976902205)$ ,  
 $p_4 = (1460.449317799282, -975.9731477430979, -466.5314749495881, 965.6402060944737)$ ,  
 $p_5 = (545.4084744819, 218.7220720302516, 472.0439121434804, -163.7241712507502)$ ,  
 $p_6 = (1313.023840410371, -497.2112405348086, 761.423662666602, -947.1254371535031)$ ,  
where  $m_t = 174$ .

	Squared matrix element [8] ( $10^{-10} \text{ GeV}^{-4}$ )
Numerical result of the method	2.34651551922455
MadGraph	2.34651551922455

# Summary

- By using several programs in combination, the complete method of calculation is given.
- The example results of  $gg \rightarrow t\bar{t} + n \text{ gluons}$  at tree level agree well with previous calculations.
- The advantage of the method is that we can compute the different processes by the same method with minimal changes.

# References I

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# Thank you