

MADDIPOLE: DIPOLE SUBTRACTION METHOD IN MADGRAPH

R I K K E R T F R E D E R I X
C E N T E R F O R P A R T I C L E P H Y S I C S
P H E N O M E N O L O G Y (U C L O U V A I N)
A N D
C E R N - T H

In collaboration with
Thomas Gehrmann & Nicolas Greiner (Univ. Zurich)
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OUTLINE

- ✿ Motivation
- ✿ What is available now in the MadDipole package
- ✿ What will be available soon

MOTIVATION...

$$\sigma^{\text{NLO}} = \int_{m+1} \left[d^{(4)} \sigma^R - d^{(4)} \sigma^A \right] + \int_m \left[\int_{\text{loop}} d^{(d)} \sigma^V + \int_1 d^{(d)} \sigma^A \right]_{\epsilon=0}$$

- ✿ “Most” of the contributions to a NLO cross section are tree-level like
- ✿ In particular, the Born, real and subtraction terms can be described by tree-level processes
- ✿ Many tree-level Monte Carlo event generators available!
- ✿ We use MadGraph/MadEvent
[Alwall, Demin, de Visscher, RF, Herquet, Maltoni, Plehn, Rainwater e Stelzer]

...AND GOAL

$$\sigma^{\text{NLO}} = \int_{m+1} \left[d^{(4)}\sigma^R - d^{(4)}\sigma^A \right] + \int_m \left[\int_{\text{loop}} d^{(d)}\sigma^V + \int_1 d^{(d)}\sigma^A \right]_{\epsilon=0}$$

- ✿ We want a **completely automatic** Monte Carlo program that calculates all the tree-level like contributions to any NLO computation in QCD
- ✿ Within the MadGraph/MadEvent framework
- ✿ Dipoles for the subtraction terms [*Catani & Seymour*]
- ✿ Massless and Massive final state particles [*Catani, Dittmaier, Seymour & Trocsanyi*]
- ✿ Open source: available to all

**WHAT IS
AVAILABLE NOW**

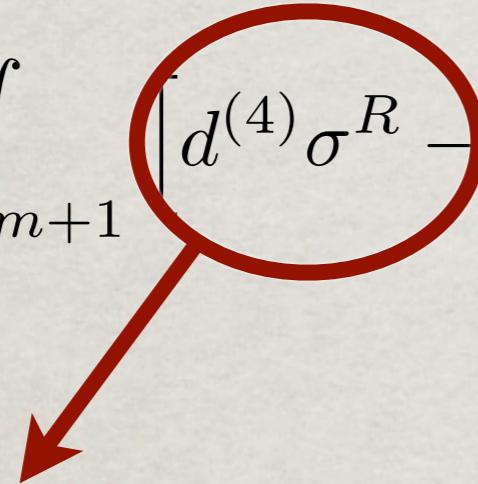
WHAT IS IN THE PACKAGE

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Normal
MadGraph



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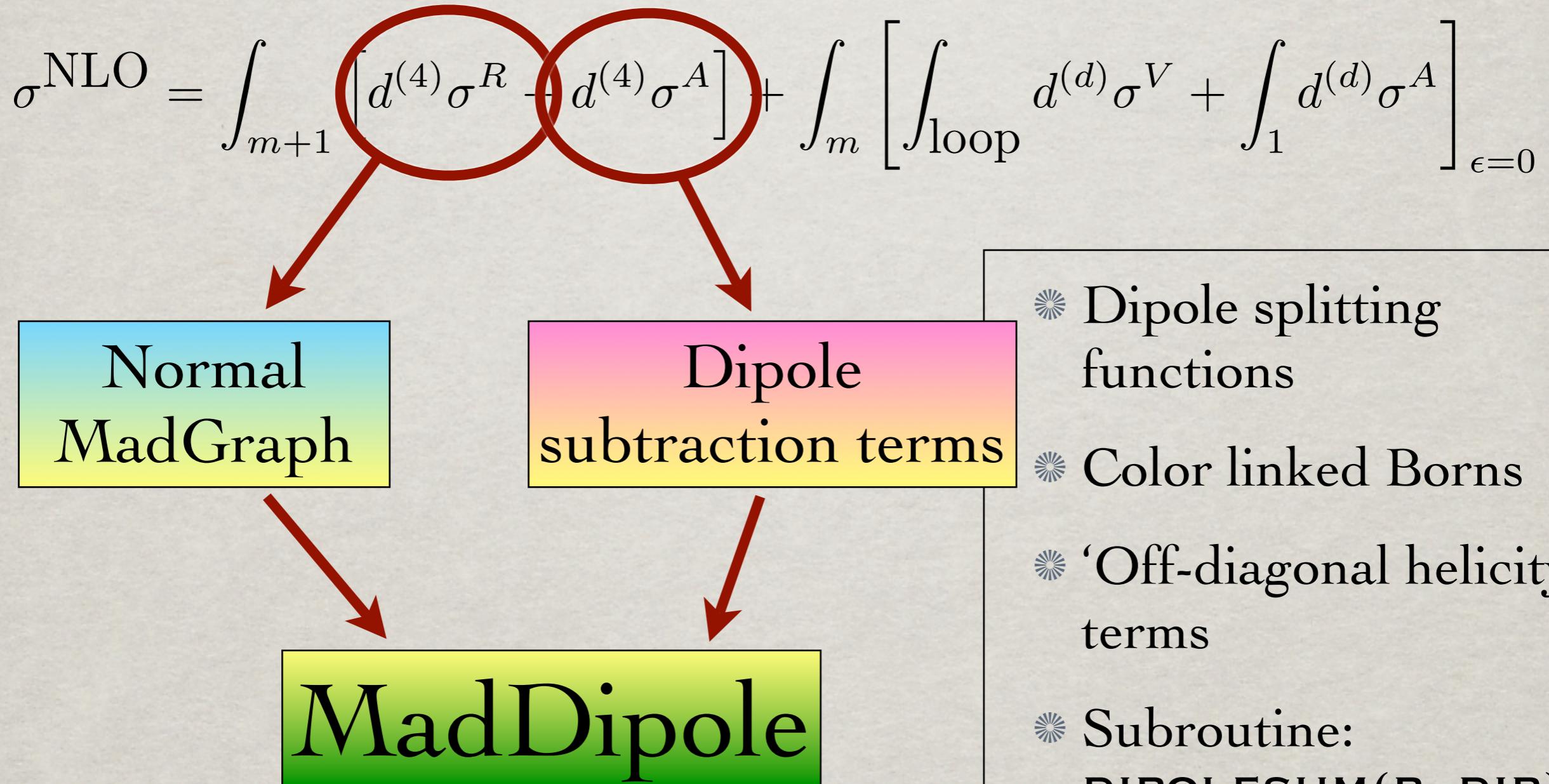
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Normal
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Dipole
subtraction terms

- ✿ Dipole splitting functions
- ✿ Color linked Borns
- ✿ ‘Off-diagonal helicity’ terms
- ✿ Subroutine:
DIPOLESUM(P, DIP)

WHAT IS IN THE PACKAGE



IN PRACTICE

- ✿ Download the MadDipole package from any of the MadGraph web pages, e.g., <http://madgraph.phys.ucl.ac.be>
- ✿ Edit the *./Cards/proc_card.dat* and give the real radiation ($n+1$) process, e.g., *gg>tt~bb~g*
- ✿ Run *./bin/newprocess* and all necessary files will be created in the *./SubProcesses/P1_gg_ttxbbxg/* directory
- ✿ Exactly the same syntax as running the normal MadGraph code!

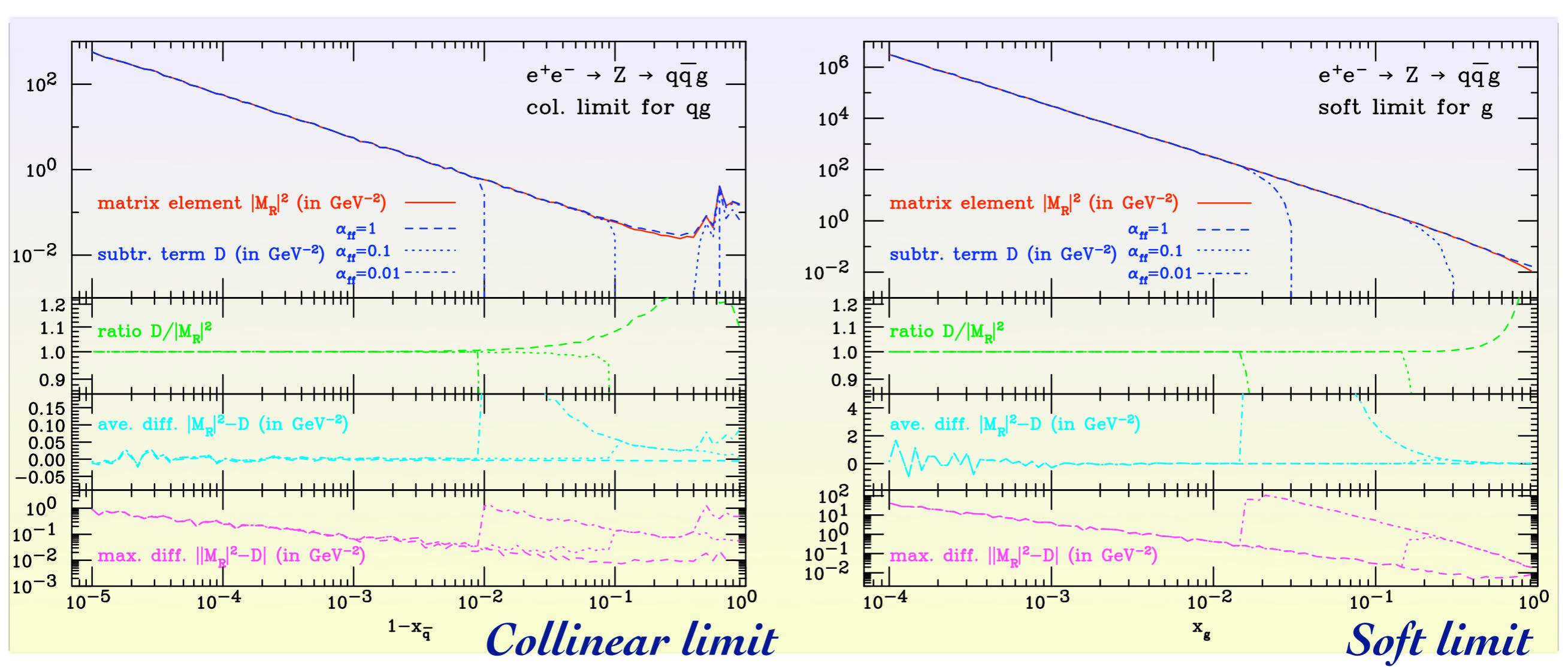
FINITE DIPOLES DUE TO HEAVY PARTICLES

- In processes with top or bottom quarks there are dipoles for which the unresolved parton is massive
- In this case the collinear singularities are regulated by the mass. However, it could develop potentially large logarithms
- Our code calculates all dipoles, but puts the non-divergent into a separate subroutine: **DIPOLESUMFINITE(...)**
- This subroutine can easily be included to cancel the potentially large logarithms
- In the limit of zero mass (or infinite energy), including the non-divergent dipoles is the same as generating the code with zero masses from the start

PHASE-SPACE RESTRICTION

- ✿ Subtraction is only needed when integrating close to the singular region
- ✿ Include the dipoles only when approaching the collinear or soft limits, using the “alpha parameter”
[Nagy & Trocsanyi]
- ✿ For the subtraction terms to the reals, a simple theta function is sufficient
- ✿ Four parameters: `alpha_ff`, `alpha_hi`, `alpha_if` & `alpha_ii` adjustable by the user

CHECKS



- ➊ Divergences are cancelled by the subtraction terms:
ratio approaches one
- ➋ Left-over rise in difference is an integrable singularity

FURTHER CHECKS AGAINST MCFM

process	subprocesses	(n+1)
Drell-Yan (W)	$q\bar{q}' \rightarrow W^+(\rightarrow e^+\nu_e)g$ $qg \rightarrow W^+(\rightarrow e^+\nu_e)q'$	
Drell-Yan (Z)	$q\bar{q} \rightarrow Z(\rightarrow e^+e^-)g$ $qg \rightarrow Z(\rightarrow e^+e^-)q$	
Drell-Yan (Z +jet)	$q\bar{q} \rightarrow Z(\rightarrow e^+e^-)q'\bar{q}'$ $q\bar{q} \rightarrow Z(\rightarrow e^+e^-)q\bar{q}$ $q\bar{q} \rightarrow Z(\rightarrow e^+e^-)gg$ $q\bar{g} \rightarrow Z(\rightarrow e^+e^-)qg$ $g\bar{g} \rightarrow Z(\rightarrow e^+e^-)q\bar{q}$	
top quark pair ($t\bar{t}$)	$q\bar{q} \rightarrow t(\rightarrow bl^+\nu_l)\bar{t}(\rightarrow \bar{b}l^-\bar{\nu}_l)g$ $qg \rightarrow t(\rightarrow bl^+\nu_l)\bar{t}(\rightarrow \bar{b}l^-\bar{\nu}_l)q$ $gg \rightarrow t(\rightarrow bl^+\nu_l)\bar{t}(\rightarrow \bar{b}l^-\bar{\nu}_l)g$	
t -channel single top with massive b -quark [63]	$gg \rightarrow t\bar{b}q\bar{q}'$ $qq' \rightarrow t\bar{b}q'q''$ $qq' \rightarrow t\bar{b}q'q''$ $qg \rightarrow t\bar{b}q'g$	

- ✿ Point-by-point in phase-space check against MCFM
[Campbell & Ellis]
- ✿ No discrepancies found
- ✿ This shows that all possible FI, IF and II dipoles (massless and massive) are okay

CHECK PROGRAM

- With the package a small code, called **CHECK_DIP**, is provided
- It calculates all the collinear and soft limits for the given process
- It prints the matrix element and subtraction terms for points closer and closer to the limit on the screen

CHECK PROGRAM

$$e^+ e^- \rightarrow Z \rightarrow u\bar{u}g$$

Collinear limit: $p(4).p(5)$ goes to zero

$p(4).p(5)/s(1,2)$,	$\sqrt{s(4,5)}$,	$ M ^2$,	$ Sub.term ^2$,	$ M ^2/ Sub.term ^2$
0.315364E+00 ,	0.561573E+03 ,	0.500612E-07 ,	0.863461E-07 ,	0.579774E+00
0.737460E-01 ,	0.271562E+03 ,	0.512878E-06 ,	0.511343E-06 ,	0.100300E+01
0.537588E-01 ,	0.231859E+03 ,	0.118507E-05 ,	0.120238E-05 ,	0.985603E+00
0.332833E-01 ,	0.182437E+03 ,	0.105116E-05 ,	0.109007E-05 ,	0.964305E+00
0.143302E-01 ,	0.119709E+03 ,	0.940775E-05 ,	0.922832E-05 ,	0.101944E+01
0.113982E-01 ,	0.106762E+03 ,	0.387269E-05 ,	0.390105E-05 ,	0.992730E+00
0.101978E-01 ,	0.100984E+03 ,	0.377575E-04 ,	0.378739E-04 ,	0.996928E+00
0.573922E-02 ,	0.757576E+02 ,	0.409948E-04 ,	0.407579E-04 ,	0.100581E+01
0.253423E-02 ,	0.503411E+02 ,	0.194822E-04 ,	0.196108E-04 ,	0.993440E+00
0.251508E-02 ,	0.501505E+02 ,	0.113841E-03 ,	0.113628E-03 ,	0.100188E+01
0.192366E-02 ,	0.438595E+02 ,	0.532619E-03 ,	0.530903E-03 ,	0.100323E+01
0.479893E-03 ,	0.219065E+02 ,	0.908399E-04 ,	0.903827E-04 ,	0.100506E+01
0.317967E-03 ,	0.178316E+02 ,	0.355093E-03 ,	0.354674E-03 ,	0.100118E+01
0.168300E-03 ,	0.129730E+02 ,	0.579438E-03 ,	0.581225E-03 ,	0.996925E+00
0.145372E-03 ,	0.120571E+02 ,	0.284682E-03 ,	0.285169E-03 ,	0.998294E+00
0.609018E-04 ,	0.780396E+01 ,	0.969336E-02 ,	0.969342E-02 ,	0.999994E+00
0.259173E-04 ,	0.509091E+01 ,	0.107919E-02 ,	0.107984E-02 ,	0.999399E+00
0.186237E-04 ,	0.431552E+01 ,	0.914959E-02 ,	0.914747E-02 ,	0.100023E+01
0.178298E-04 ,	0.422253E+01 ,	0.149607E-02 ,	0.149572E-02 ,	0.100023E+01
0.910116E-05 ,	0.301681E+01 ,	0.436849E-02 ,	0.437017E-02 ,	0.999616E+00
0.156505E-05 ,	0.125102E+01 ,	0.277892E-01 ,	0.277948E-01 ,	0.999800E+00
0.152407E-05 ,	0.123453E+01 ,	0.287201E-01 ,	0.287121E-01 ,	0.100028E+01
0.140303E-05 ,	0.118449E+01 ,	0.236653E-01 ,	0.236693E-01 ,	0.999831E+00
0.353217E-06 ,	0.594321E+00 ,	0.128877E+00 ,	0.128881E+00 ,	0.999970E+00

CHECK PROGRAM

$$e^+ e^- \rightarrow Z \rightarrow u\bar{u}g$$

$$\alpha_{ff} = 0.1$$

Limit: p(5) goes soft

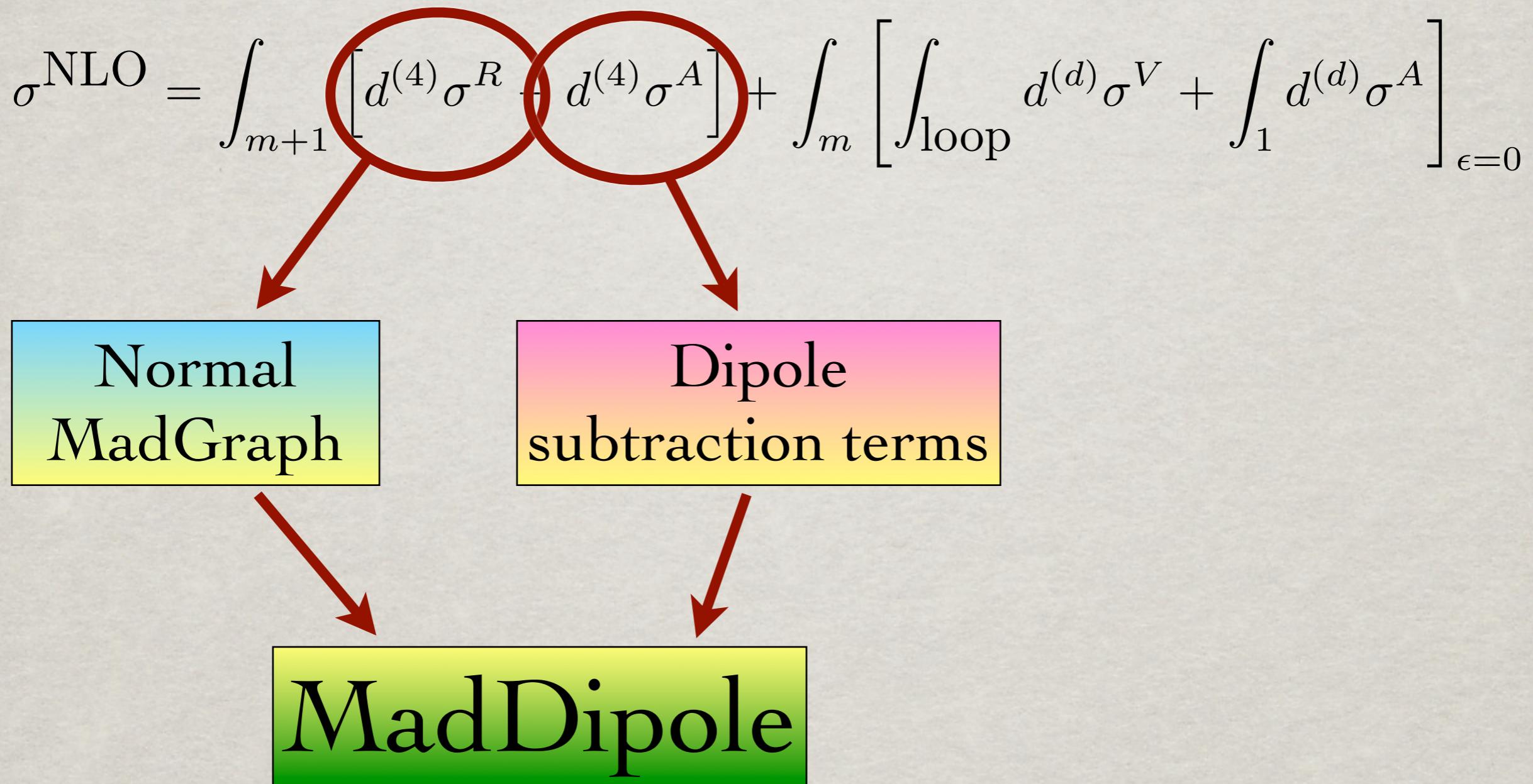
p(0, 5)^2/s(1,2) ,	M ^2 ,	Sub.term ^2 ,	M ^2/ Sub.term ^2
0.869211E-03 ,	0.109575E-06 ,	0.000000E+00 ,	INF
0.688787E-03 ,	0.193565E-06 ,	0.000000E+00 ,	INF
0.421426E-03 ,	0.699189E-06 ,	0.000000E+00 ,	INF
0.387974E-03 ,	0.659780E-06 ,	0.000000E+00 ,	INF
0.336094E-03 ,	0.154721E-05 ,	0.000000E+00 ,	INF
0.185406E-03 ,	0.376282E-05 ,	0.265364E-05 ,	0.141799E+01
0.137483E-03 ,	0.818395E-05 ,	0.816317E-05 ,	0.100255E+01
0.103854E-03 ,	0.269806E-04 ,	0.269632E-04 ,	0.100064E+01
0.637535E-04 ,	0.314766E-04 ,	0.314798E-04 ,	0.999898E+00
0.376601E-04 ,	0.122914E-03 ,	0.122984E-03 ,	0.999433E+00
0.191145E-04 ,	0.780061E-03 ,	0.778950E-03 ,	0.100143E+01
0.179479E-04 ,	0.565700E-03 ,	0.565041E-03 ,	0.100117E+01
0.129797E-04 ,	0.111247E-02 ,	0.111223E-02 ,	0.100021E+01
0.105788E-04 ,	0.203953E-02 ,	0.203945E-02 ,	0.100004E+01
0.774787E-05 ,	0.346362E-02 ,	0.346388E-02 ,	0.999927E+00
0.563720E-05 ,	0.610302E-02 ,	0.610181E-02 ,	0.100020E+01
0.343292E-05 ,	0.235165E-01 ,	0.235160E-01 ,	0.100002E+01
0.909122E-06 ,	0.347171E+00 ,	0.347161E+00 ,	0.100003E+01

**WHAT WILL BE
AVAILABLE SOON**

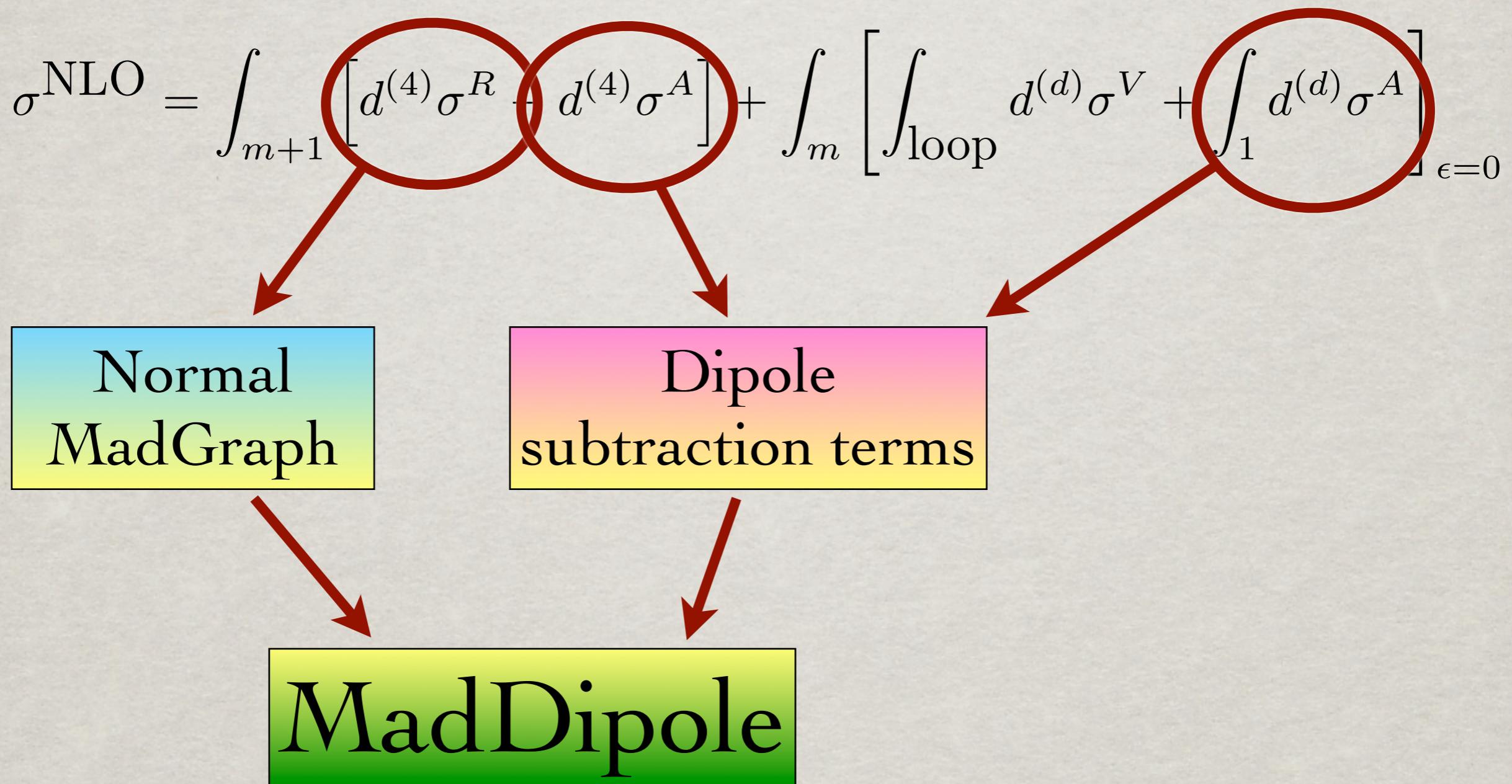
INTEGRATED DIPOLES

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INTEGRATED DIPOLES



INTEGRATED DIPOLES



INTEGRATED DIPOLES: SOME DETAILS

- ✿ For integrated splitting function with initial state particles contains plus-distributions and delta function: we need the Born's at x and also at $x=1$. (With x the parton energy fraction)
- ✿ Non-trivial dependence on the alpha parameter
 - New integrals were needed
- ✿ Results depend on regularization scheme: both 't Hooft-Veltman and Dimensional Reduction implemented
- ✿ Common factors in loop integrals can be excluded
 - Two implementations: with and without expanding this term:

$$\mathcal{V}_{ij,k} = \int [dp_i(\tilde{p}_{ij}, \tilde{p}_k)] \frac{1}{2p_i \cdot p_j} < \mathbf{V}_{ij,k} > \equiv \frac{\alpha_S}{2\pi} \underbrace{\frac{1}{\Gamma(1-\epsilon)} \left(\frac{4\pi\mu^2}{s_{ij,k}} \right)^\epsilon}_{=1+\left(\log\left(\frac{4\pi\mu}{s_{ij,k}}\right)-\gamma\right)\epsilon+O(\epsilon^2)} \mathcal{V}_{ij}(\epsilon)$$

INTEGRATED DIPOLES: IN PRACTICE

- ✿ One new subroutine: **INTDIPOLES(P,X, EPSSQ, EPS, FINITE)**
- ✿ Momenta and parton energy fractions as input, and returns three 5-dimensional vectors with
 - 1) $\delta(1-x)$ terms
 - 2) terms regular in x
 - 3) plus-distribution (singular in $x=1$)
- ✿ and for the massive case ($x_+=1-4m_f^2/Q^2$)
 - 4) $\delta(x_+-x)$ terms
 - 5) x_+ -distributions
- ✿ Implementation done → needs heavy testing

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Divergent terms okay
Alpha dependence okay
Overall normalization is
in progress...

PHASE SPACE INTEGRATION

- ✿ The phase space integration over subtracted NLO matrix elements is more complicated than at LO
- ✿ Not a problem for simple $2 \rightarrow 2$ (or $2 \rightarrow 3$) processes, but becomes an issue in more complicated cases
- ✿ We are working on a fully automatic phase space generator that can efficiently and in a parallel nature integrate the subtracted matrix elements

BEYOND THE SM

- ✿ MadGraph knows how to do BSM physics...
- ✿ However, in New Physics models with new colored particles different types of dipoles are needed
 - ✿ For example, SUSY squarks are heavy **color triplet scalars**
- ✿ The MadDipole code is modular enough to easily deal with these **new structures**, but the dipoles need to be implemented, tested, ...

CONCLUSIONS

- ⌘ **MadDipole** is a package that calculates any real matrix element, including all dipole subtraction terms
- ⌘ It calculates also the contributions to be added to the virtual corrections
- ⌘ The phase space integration of the subtracted matrix elements is still work in progress
- ⌘ An automatic NLO package that only needs the virtual corrections as external input is within reach

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wiki-page (accessible from any of the MadGraph web pages):
<http://cp3wks05.fynu.ucl.ac.be/twiki/bin/view/Software/MadDipole>