

aMC@NLO and top pair production at LC

Olivier Mattelaer

Universite Catholique de Louvain

for the MadGraph/aMC@NLO team

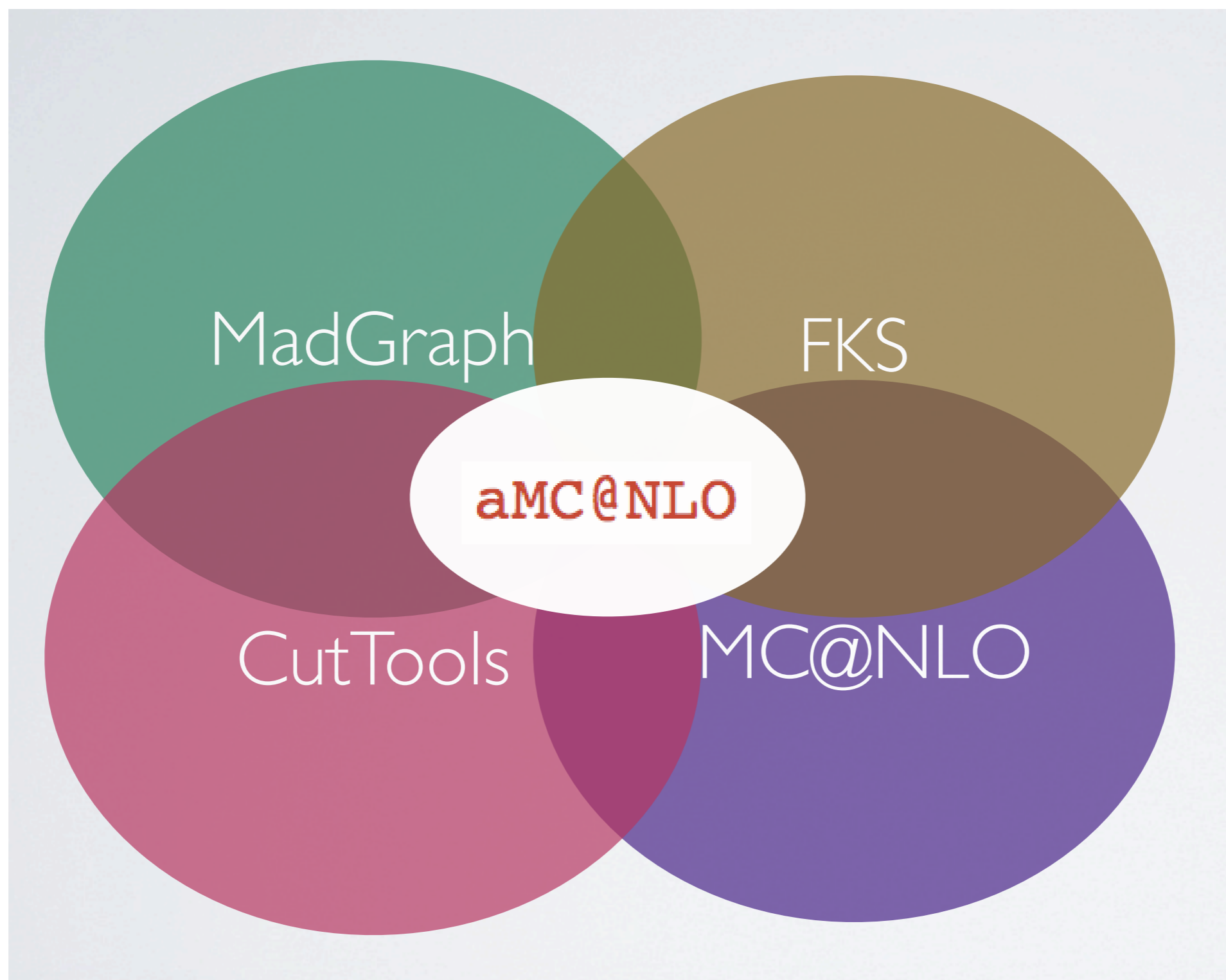
Full list of contributors:

<http://amcatnlo.web.cern.ch/amcatnlo/people.htm>

Plan of the Talk

- aMC@NLO
 - ➔ MadLoop
 - ➔ MadFKS
 - ➔ NLO+PS
- MadSpin
- DEMO
- top pair production at LC
- Conclusion

aMC@NLO: A Joint Venture



aMC@NLO

- Why automation?
 - ➔ Time: Less tools, means more time for physics
 - ➔ Robust: Easier to test, to trust
 - ➔ Easy: One framework/tool to learn

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 - ➔ Reliable prediction of the total rate
 - ➔ Reduction of the theoretical uncertainty

aMC@NLO

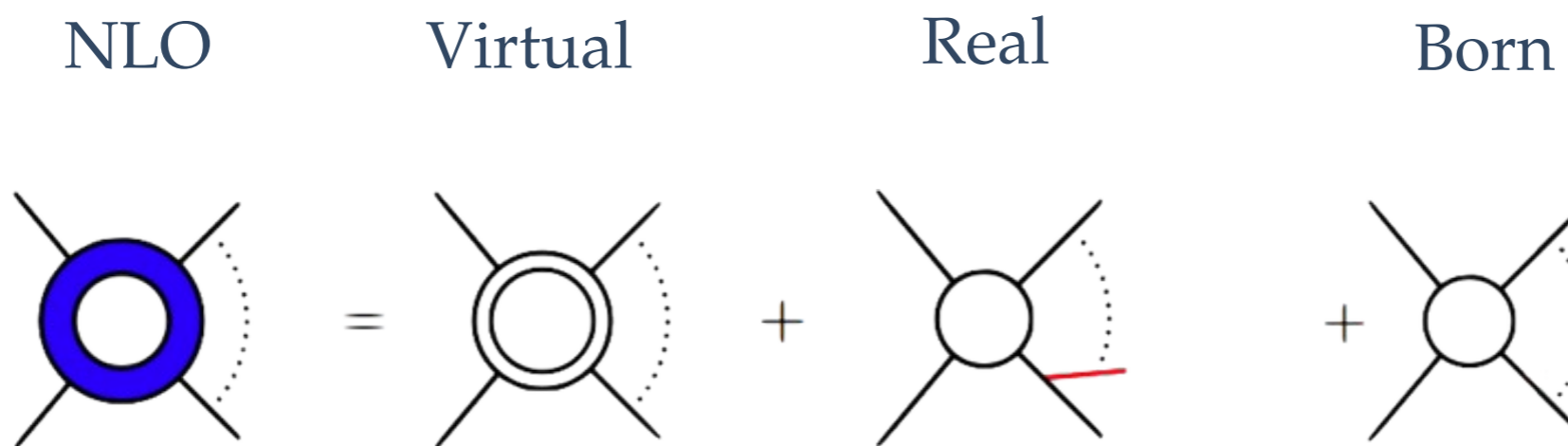
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- Why **matched to the PS**?
 - ➔ Parton are not an detector observables
 - ➔ Matching cure some fix-order ill behaved observables

NLO Basics

NLO	Virtual	Real	Born
=			
+			
+			

$$\sigma^{NLO} = \int_m d^{(d)} \sigma^V + \int_{m+1} d^{(d)} \sigma^R + \int_m d^{(4)} \sigma^B$$

NLO Basics

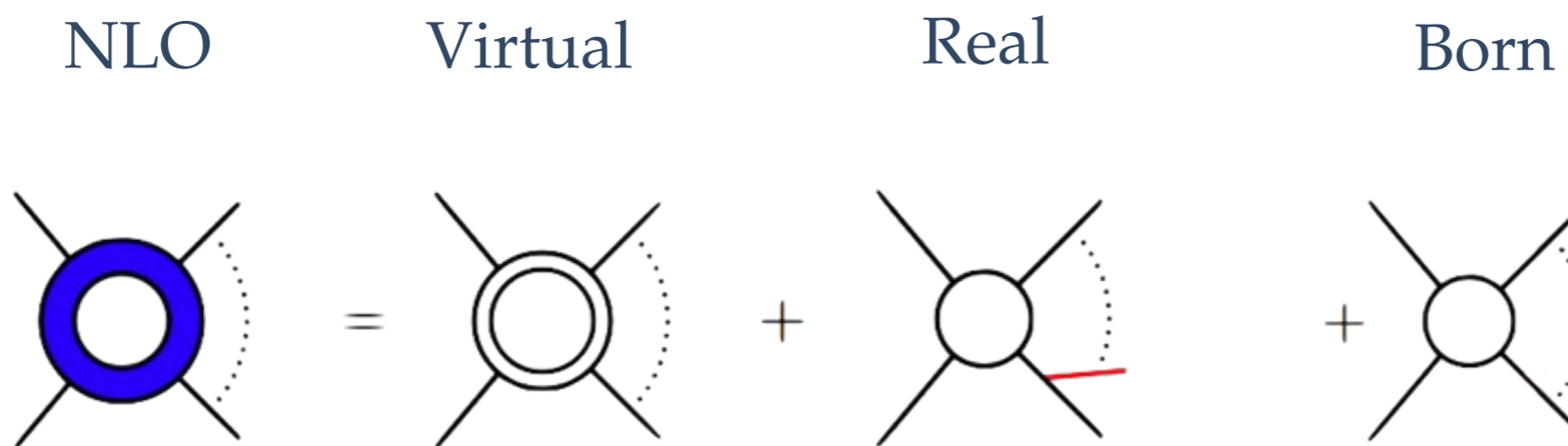


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Need to deal with singularities

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MadLoop

MadFKS

MadGraph

MADLOOP

The virtual

OPP Reduction

- decomposition to scalar integrals works at the level of the **integrals**

$$\begin{aligned}\mathcal{M}^{1\text{-loop}} &= \sum_{i_0 < i_1 < i_2 < i_3} d_{i_0 i_1 i_2 i_3} \text{Box}_{i_0 i_1 i_2 i_3} \\ &+ \sum_{i_0 < i_1 < i_2} c_{i_0 i_1 i_2} \text{Triangle}_{i_0 i_1 i_2} \\ &+ \sum_{i_0 < i_1} b_{i_0 i_1} \text{Bubble}_{i_0 i_1} \\ &+ \sum_{i_0} a_{i_0} \text{Tadpole}_{i_0} \\ &+ R + \mathcal{O}(\epsilon)\end{aligned}$$

[Ossola, Papadopoulos, Pittau 2006]

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$$\begin{aligned}
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Spurious term

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Spurious term

- feed cut tools with numerator value and it returns the coefficients

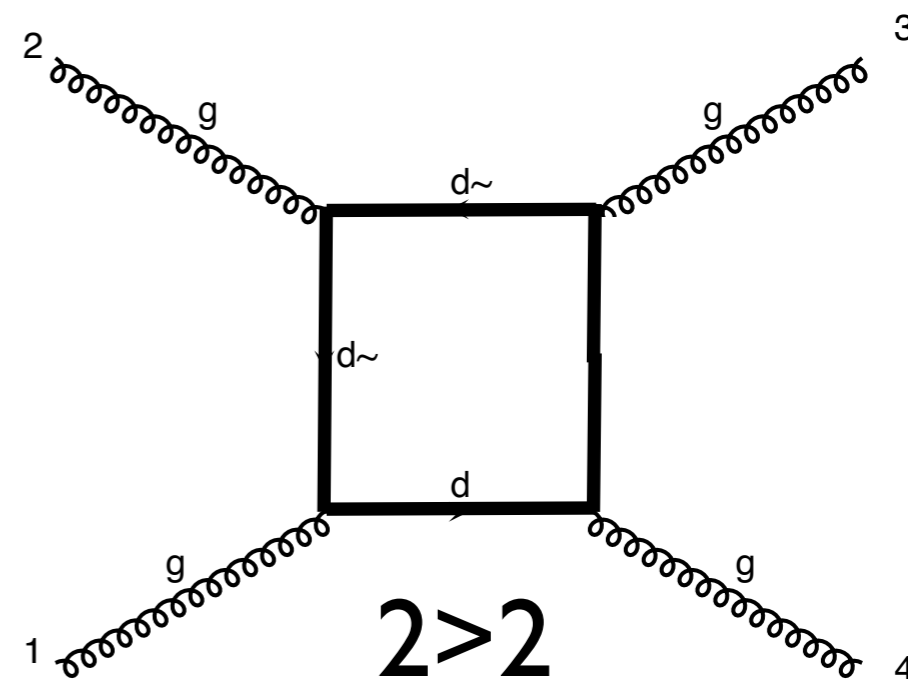
[Ossola, Papadopoulos, Pittau 2006]

OPP in a nutshell

- In OPP reduction we reduce the system at the integrand level.
- We can solve the system numerically: we only need a numerical function of the (numerator of) integrand. We can set-up a system of linear equations by choosing specific values for the loop momentum l , depending on the kinematics of the event
- OPP reduction is implemented in CutTools (publicly available). Given the integrand, CutTools provides all the coefficients in front of the scalar integrals and the R1 term
- The OPP reduction leads to numerical instabilities whose origins are not well under control. Require quadruple precision.
- Analytic information is needed for the R2 term, but can be compute once and for all for a given model

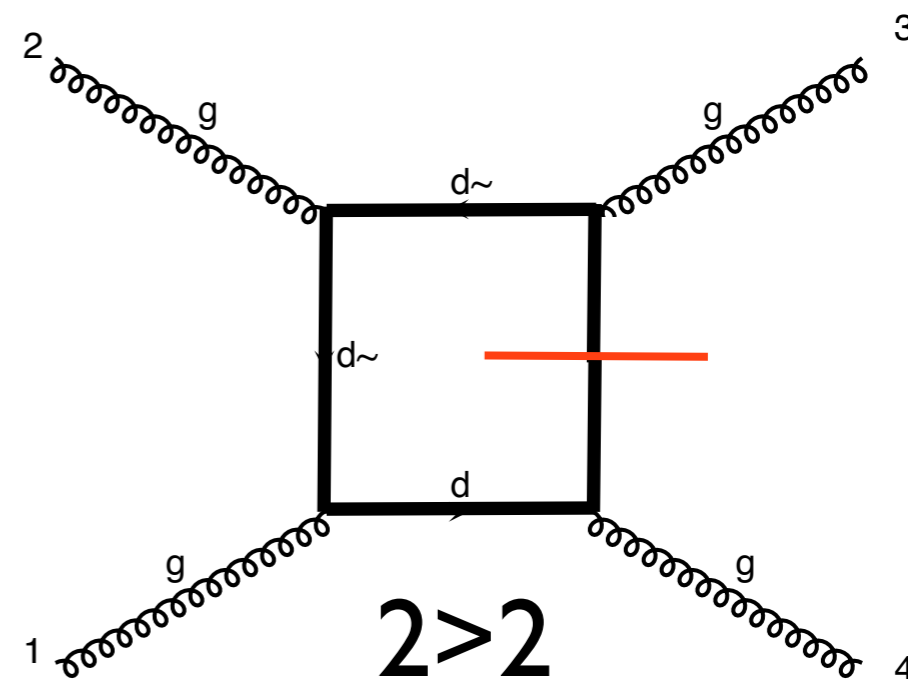
MADLOOP

- Diagram Generation



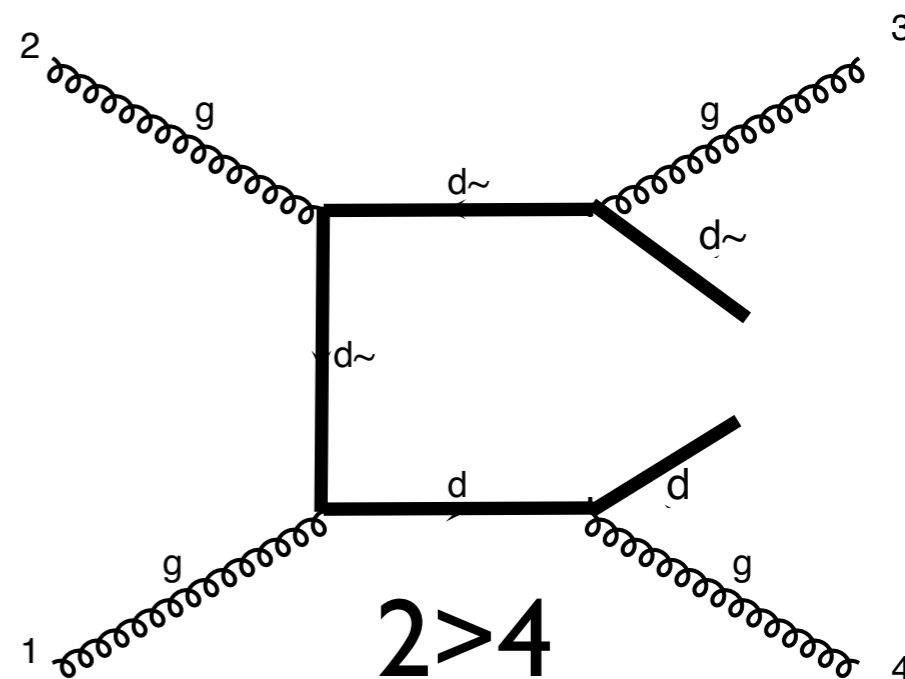
MADLOOP

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MADLOOP

- Diagram Generation
 - ➔ Generate diagrams with 2 extra particles
 - ➔ Need to filter result



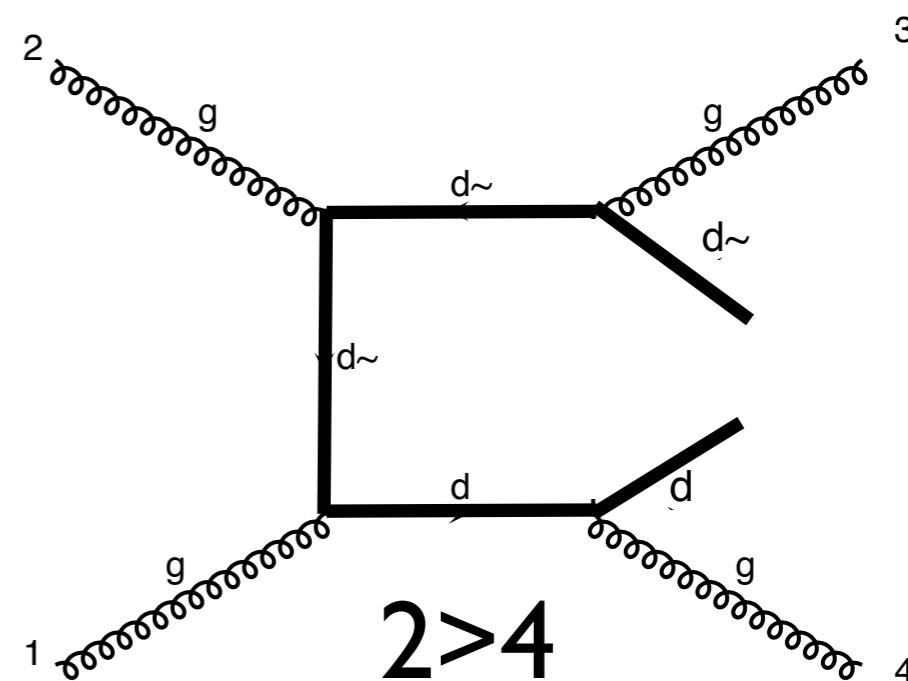
MADLOOP

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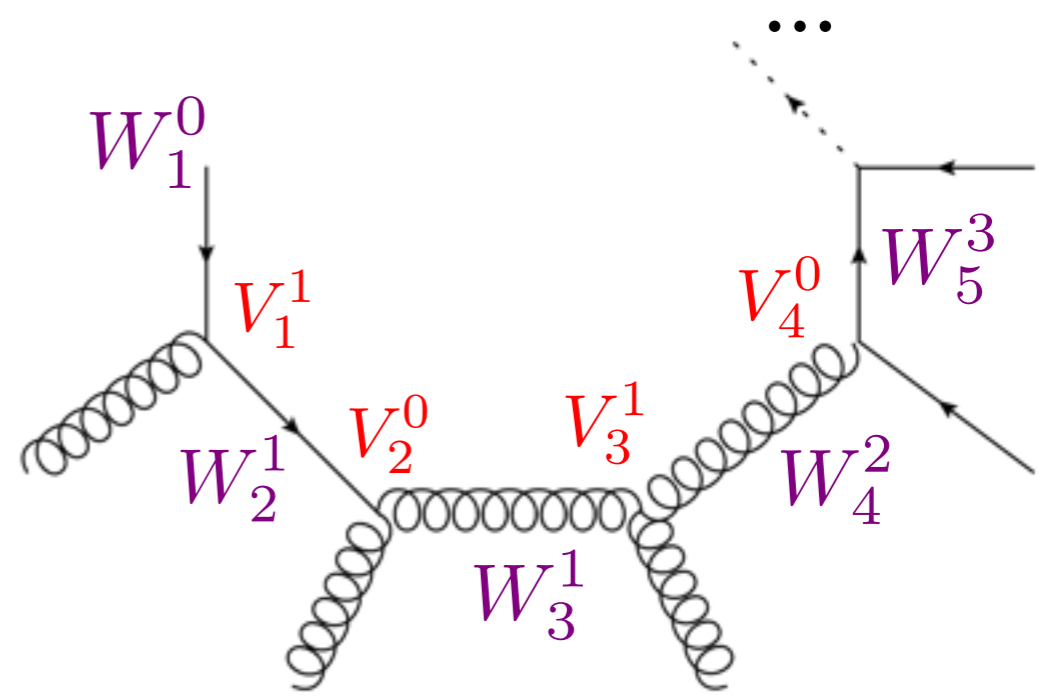
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- Evaluation of the Numerator:

- ➔ OpenLoops technique [S. Pozzorini & al.(2011)]



$$\mathcal{N}(l^\mu) = \sum_{r=0}^{r_{max}} C_{\mu_0 \mu_1 \dots \mu_r}^{(r)} l^{\mu_0} l^{\mu_1} \dots l^{\mu_r}$$



MADFKS

The real

FKS subtraction

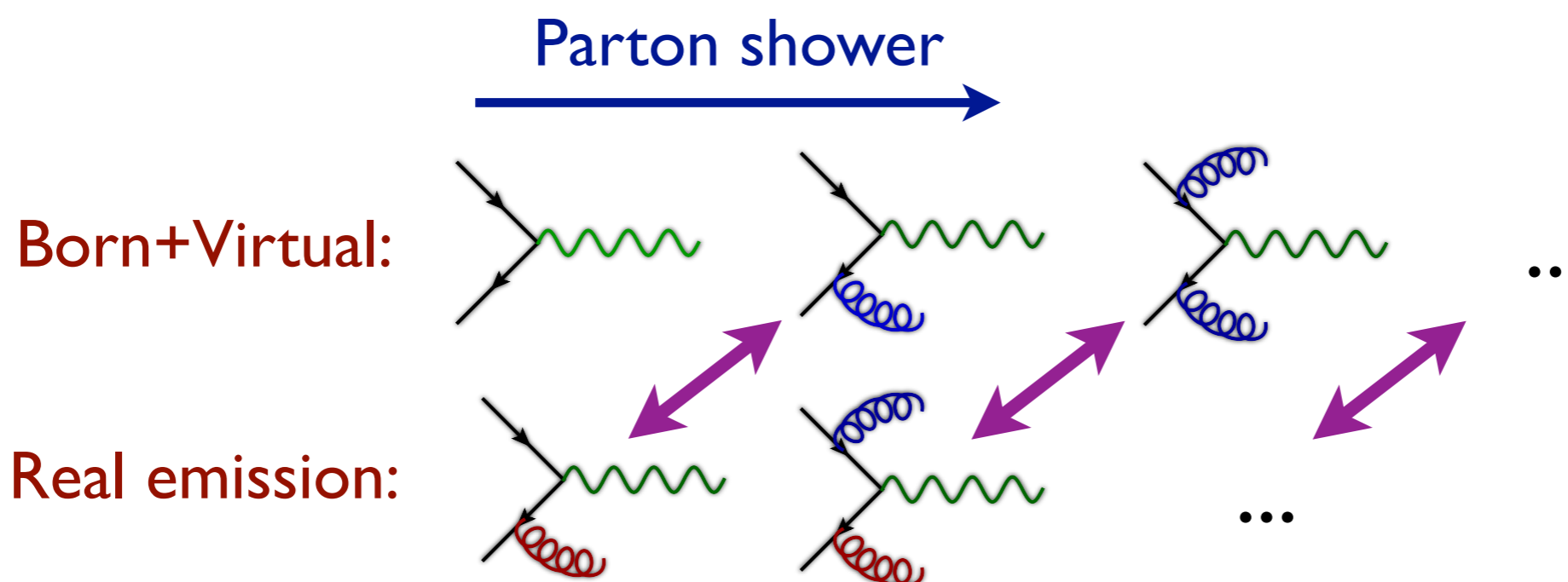
- Find parton pairs i, j that can give collinear singularities
- Split the phase space into regions with one collinear singularities
- Integrate them independently
 - ➔ with an adhoc PS parameterization
 - ➔ can be run in parallel
- # of contributions $\sim n^2$

[S. Frixione, Z Kunst, A Signer (1995)]

MC@NLO

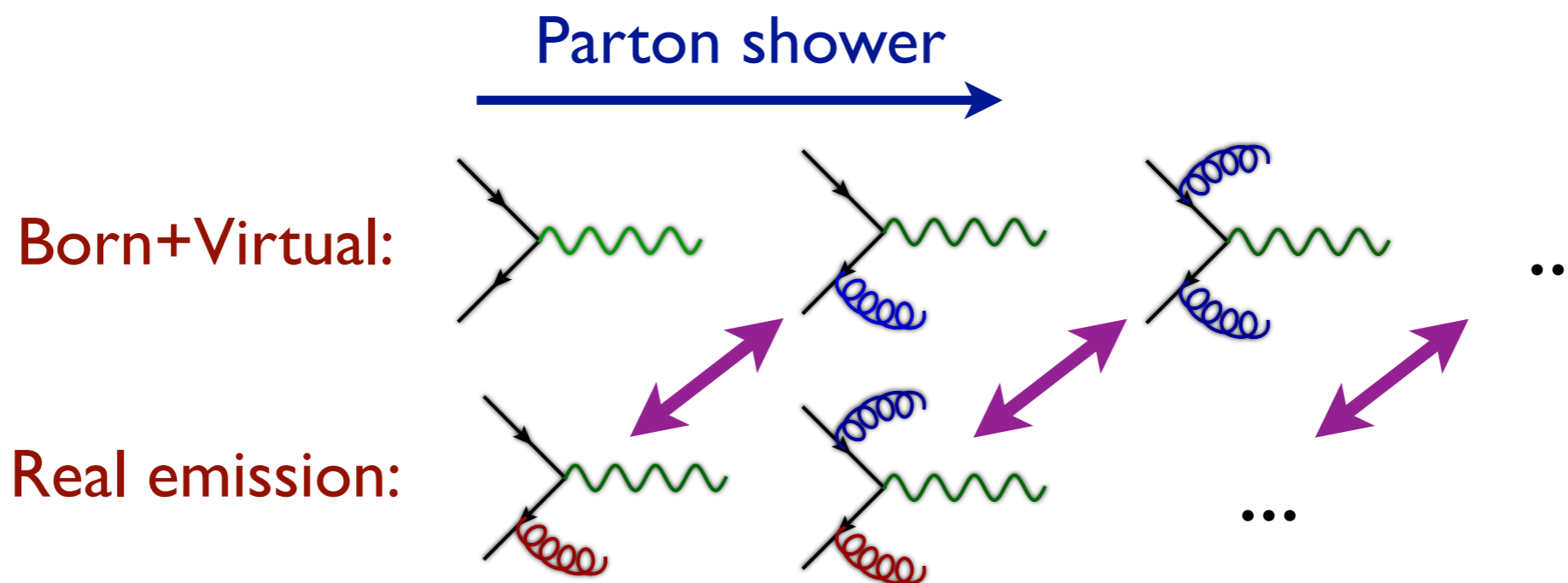
Matching to the shower

Sources of double counting



- There is double counting between the real emission matrix elements and the parton shower: the extra radiation can come from the matrix elements or the parton shower
- There is also an overlap between the virtual corrections and the Sudakov suppression in the zero-emission probability

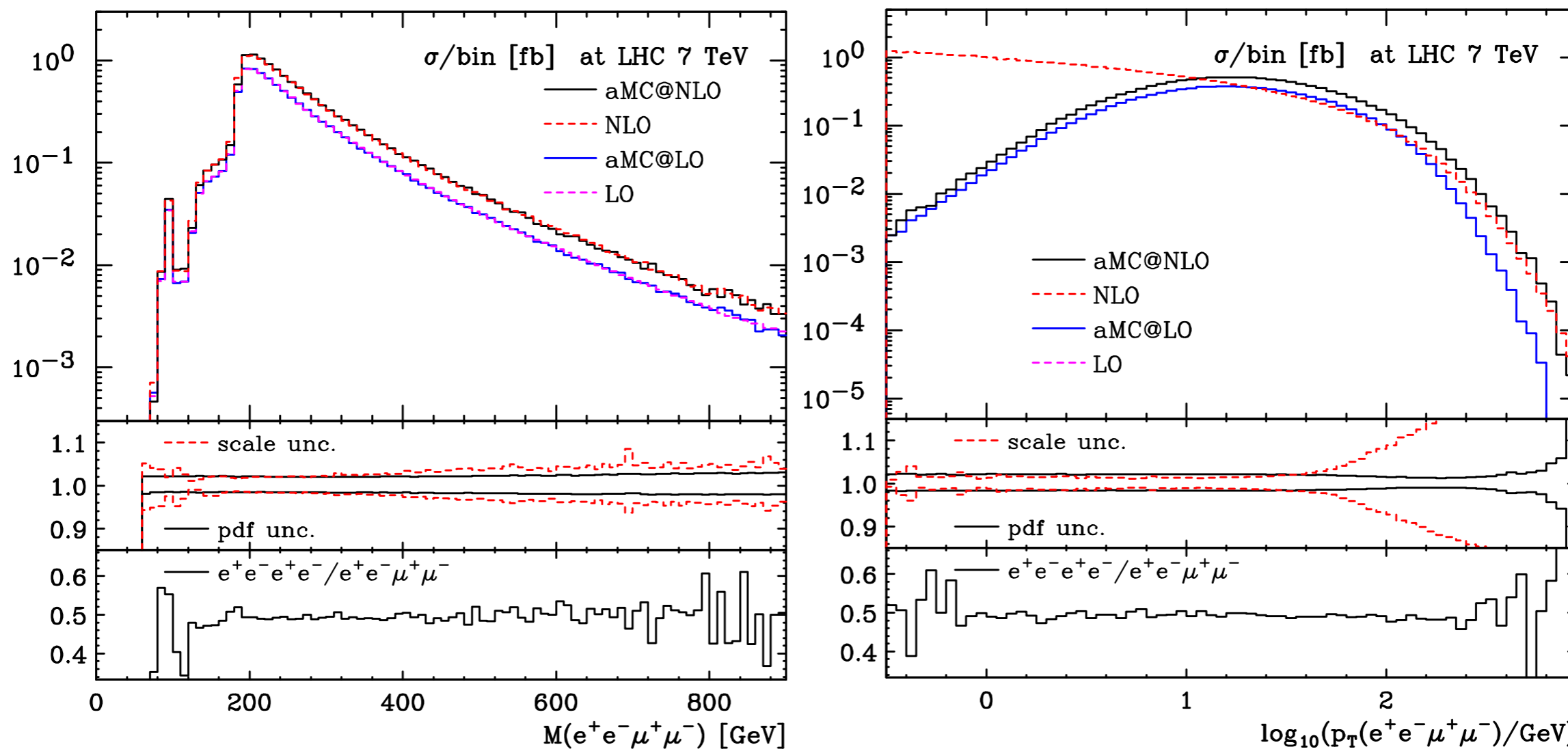
MC@NLO procedure



$$\frac{d\sigma_{\text{NLOwPS}}}{dO} = \left[d\Phi_m (B + \int_{\text{loop}} V + \int d\Phi_1 MC) \right] I_{\text{MC}}^{(m)}(O) + \left[d\Phi_{m+1} (R - MC) \right] I_{\text{MC}}^{(m+1)}(O)$$

- Double counting is explicitly removed by including the “shower subtraction terms”

Four-lepton production



- 4-lepton invariant mass is almost insensitive to parton shower effects.
- 4-lepton transverse momenta is extremely sensitive

[Frederix, Frixione, Hirschi, Maltoni, Pittau & Torrielli (2011)]

results

- Errors are the MC integration uncertainty only
- Cuts on jets, γ^*/Z decay products and photons, but **no cuts on b quarks** (their mass regulates the IR singularities)
- Efficient handling of **exceptional phase-space points**: their uncertainty always at least two orders of magnitude smaller than the integration uncertainty
- Running time: **two weeks on ~150 node cluster** leading to rather small integration uncertainties

	Process	μ	n_{lf}	Cross section (pb)	
				LO	NLO
a.1	$pp \rightarrow t\bar{t}$	m_{top}	5	123.76 ± 0.05	162.08 ± 0.12
a.2	$pp \rightarrow tj$	m_{top}	5	34.78 ± 0.03	41.03 ± 0.07
a.3	$pp \rightarrow tjj$	m_{top}	5	11.851 ± 0.006	13.71 ± 0.02
a.4	$pp \rightarrow t\bar{b}j$	$m_{top}/4$	4	25.62 ± 0.01	30.96 ± 0.06
a.5	$pp \rightarrow t\bar{b}jj$	$m_{top}/4$	4	8.195 ± 0.002	8.91 ± 0.01
b.1	$pp \rightarrow (W^+ \rightarrow)e^+\nu_e$	m_W	5	5072.5 ± 2.9	6146.2 ± 9.8
b.2	$pp \rightarrow (W^+ \rightarrow)e^+\nu_e j$	m_W	5	828.4 ± 0.8	1065.3 ± 1.8
b.3	$pp \rightarrow (W^+ \rightarrow)e^+\nu_e jj$	m_W	5	298.8 ± 0.4	300.3 ± 0.6
b.4	$pp \rightarrow (\gamma^*/Z \rightarrow)e^+e^-$	m_Z	5	1007.0 ± 0.1	1170.0 ± 2.4
b.5	$pp \rightarrow (\gamma^*/Z \rightarrow)e^+e^- j$	m_Z	5	156.11 ± 0.03	203.0 ± 0.2
b.6	$pp \rightarrow (\gamma^*/Z \rightarrow)e^+e^- jj$	m_Z	5	54.24 ± 0.02	56.69 ± 0.07
c.1	$pp \rightarrow (W^+ \rightarrow)e^+\nu_e b\bar{b}$	$m_W + 2m_b$	4	11.557 ± 0.005	22.95 ± 0.07
c.2	$pp \rightarrow (W^+ \rightarrow)e^+\nu_e t\bar{t}$	$m_W + 2m_{top}$	5	0.009415 ± 0.000003	0.01159 ± 0.00001
c.3	$pp \rightarrow (\gamma^*/Z \rightarrow)e^+e^- b\bar{b}$	$m_Z + 2m_b$	4	9.459 ± 0.004	15.31 ± 0.03
c.4	$pp \rightarrow (\gamma^*/Z \rightarrow)e^+e^- t\bar{t}$	$m_Z + 2m_{top}$	5	0.0035131 ± 0.0000004	0.004876 ± 0.000002
c.5	$pp \rightarrow \gamma t\bar{t}$	$2m_{top}$	5	0.2906 ± 0.0001	0.4169 ± 0.0003
d.1	$pp \rightarrow W^+W^-$	$2m_W$	4	29.976 ± 0.004	43.92 ± 0.03
d.2	$pp \rightarrow W^+W^- j$	$2m_W$	4	11.613 ± 0.002	15.174 ± 0.008
d.3	$pp \rightarrow W^+W^+ jj$	$2m_W$	4	0.07048 ± 0.00004	0.1377 ± 0.0005
e.1	$pp \rightarrow HW^+$	$m_W + m_H$	5	0.3428 ± 0.0003	0.4455 ± 0.0003
e.2	$pp \rightarrow HW^+ j$	$m_W + m_H$	5	0.1223 ± 0.0001	0.1501 ± 0.0002
e.3	$pp \rightarrow HZ$	$m_Z + m_H$	5	0.2781 ± 0.0001	0.3659 ± 0.0002
e.4	$pp \rightarrow HZ j$	$m_Z + m_H$	5	0.0988 ± 0.0001	0.1237 ± 0.0001
e.5	$pp \rightarrow Ht\bar{t}$	$m_{top} + m_H$	5	0.08896 ± 0.00001	0.09869 ± 0.00003
e.6	$pp \rightarrow Hb\bar{b}$	$m_b + m_H$	4	0.16510 ± 0.00009	0.2099 ± 0.0006
e.7	$pp \rightarrow Hjj$	m_H	5	1.104 ± 0.002	1.036 ± 0.002

MadSpin

Decay with Full Spin correlation

[P.Artoisenet, R. Frederix, OM, R. RietKerk (2012)]

MadSpin

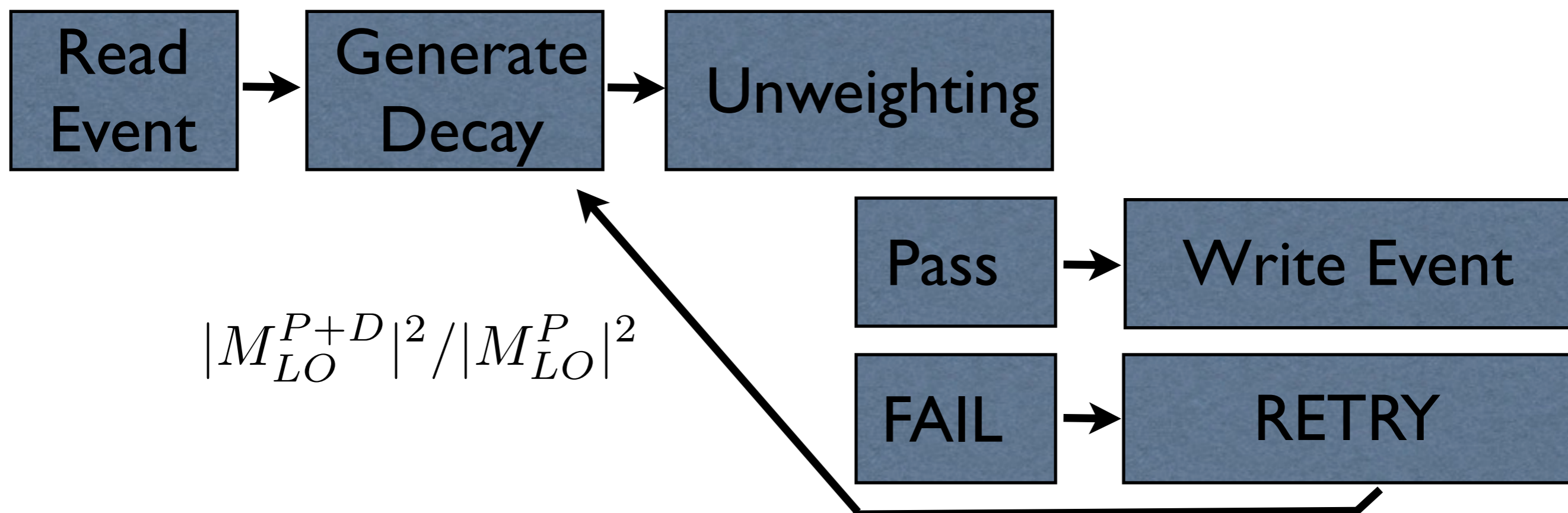
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 - ➔ For a sample of events include the decay of unstable final states particles.
 - ➔ Keep full spin correlations and finite width effect
 - ➔ Keep unweighted events

MadSpin

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- Solution:

[Frixione, Leenen, Motylinski, Webber (2007)]



MadSpin

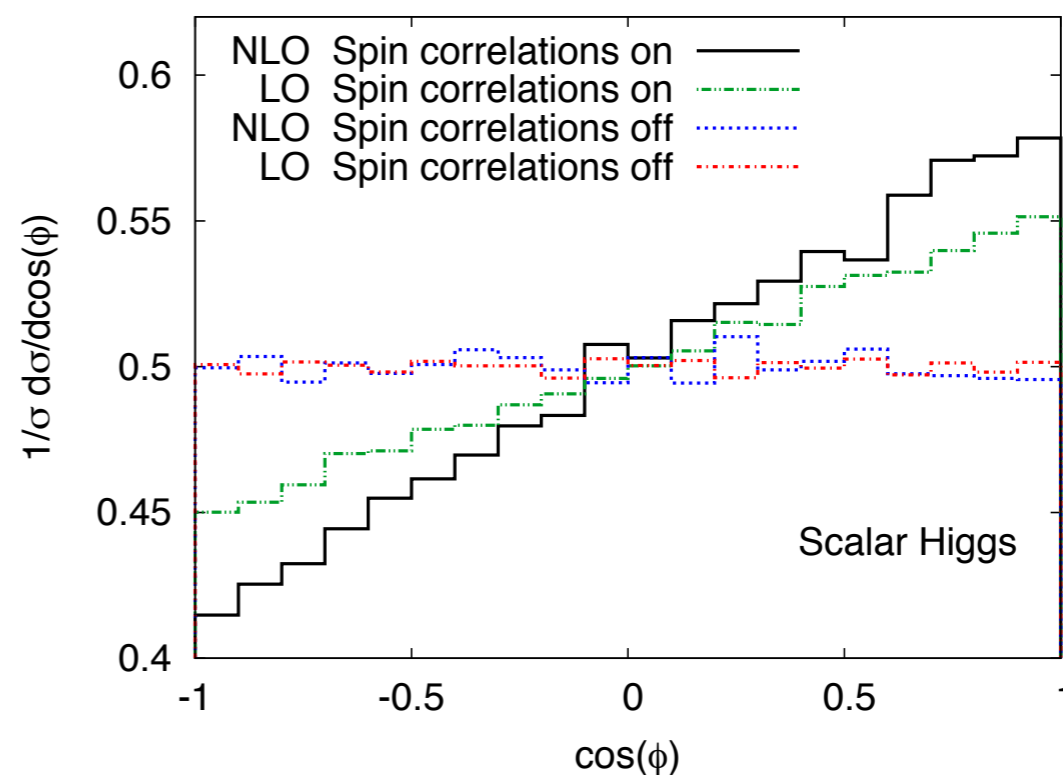
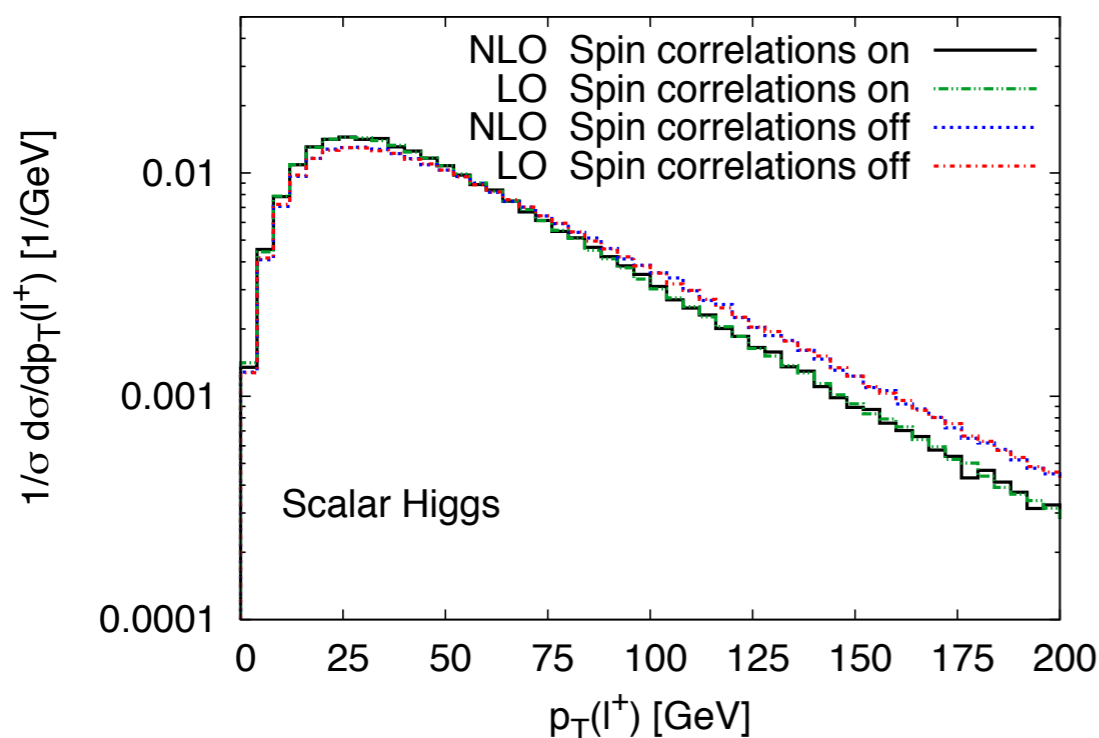
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- Example $t \bar{t} h$:



DEMO

Is it really automatic?

DEMO

- 1) Download the code

The MadGraph Matrix Element Generator version 5

Registered 2009-09-15 by [Michel Herquet](#)

The version 5 of the MadGraph Matrix Element Generator for the simulation of parton-level events for decay and collision processes at high energy colliders. Allows matrix element generation and event generation for any model that can be written as a Lagrangian, using the output of the FeynRules Feynman rule calculator. Provides output in multiple formats and languages, including Fortran MadEvent, Fortran Standalone matrix elements, C++ matrix elements, and Pythia 8 process libraries.

Note that process generation can also be done directly online at <http://madgraph.phys.ucl.ac.be> or <http://madgraph.hep.uiuc.edu>.
If you use MadGraph 5, please cite JHEP 1106(2011)128, arXiv:1106.0522 [hep-ph].

Installation:
MadGraph 5 needs Python version 2.6 or 2.7. The latest stable release is in the trunk, which can be branched using the Bazaar versioning system:
`bzr branch lp:madgraph5`
or be downloaded as a tar.gz package to the right. This release contains everything needed for process generation in multiple models, as well as event generation through MadEvent, and standalone matrix element evaluation for Fortran or C++ output.
In order to use the process library output for Pythia 8, you need Pythia 8.150 or later installed.

Getting started:
Run `bin/mg5` and type "help" to learn how to run MadGraph 5 using the command interface, or run the interactive quick-start tutorial by typing "tutorial".
Or copy the Template, edit the `Cards/proc_card_mg5.dat` and run `bin/newprocess_mg5`.

Examples of process generation syntax:
`pp > w+jj`
`pp > tt, t > bjj, t- > b-l-vl-`
`e+e- > z > n2 n2, (n2 > x1+w, x1 > l+v-l-), n2 > jj n1`

To output model files for MadGraph 5 with FeynRules, use version 1.6 or later, and use the WriteUFO command.

[Change branding](#)

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Downloads

Latest version is 1.5.0

[MadGraph5_v1.5.9.tar.gz](#)

[MadGraph5_v...eta3.tar.gz](#)

Released on 2012-09-29

[All downloads](#)

Announcements

DEMO

- launch the code [./bin/mg5]
 - ➔ Exactly like MG5 !!!

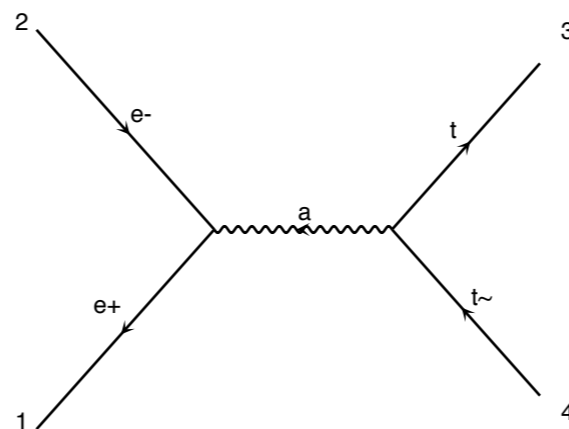
```
Terminal — Python — 201x58
bash
madevent
madgraph@server02:~
omatt@in...madgraph5 ...
omatt@in...5_interface
bash
bash
Python

[@oliviers-MacBook-Pro ~]$ cd MadGraph5_v2
MadGraph5_v2.0.0.beta3.tar.gz MadGraph5_v2.0.0.beta3.tar.gz.1 MadGraph5_v2_0_0_beta3/
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[@oliviers-MacBook-Pro ~]$ cd MadGraph5_v2_0_0_beta3/
[@oliviers-MacBook-Pro MadGraph5_v2_0_0_beta3]$ ./bin/mg5
*****
*
*     W E L C O M E to M A D G R A P H  5
*
*
*     * * * * * 5 * * * * *
*
*
*
*     VERSION 2.0.0.beta3           2013-02-14
*
*
* The MadGraph Development Team - Please visit us at
* https://server06.fynu.ucl.ac.be/projects/madgraph
*
*
*     Type 'help' for in-line help.
*     Type 'tutorial' to learn how MGS works
*     Type 'tutorial aMCatNLO' to learn how aMC@NLO works
*     Type 'tutorial MadLoop' to learn how MadLoop works
*
*
*****
load MGS configuration from /Users/omatt/.mg5/mg5_configuration.txt
load MGS configuration from input/mg5_configuration.txt
set lhpdf to lhpdf-config
set fastjet to /Users/omatt/programme/ma5_v1.1.2/tools/fastjet/bin/fastjet-config
Loading default model: sm
INFO: load particles
INFO: load vertices
INFO: Restrict model sm with file models/sm/restrict_default.dat .
INFO: Run "set stdout_level DEBUG" before import for more information.
INFO: Change particles name to pass to MGS convention
Defined multiparticle p = g u c d s u~ c~ d~ s~
Defined multiparticle j = g u c d s u~ c~ d~ s~
Defined multiparticle l+ = e+ mu+
Defined multiparticle l- = e- mu-
Defined multiparticle vl = ve vm vt
Defined multiparticle vl~ = ve~ vm~ vt~
Defined multiparticle all = g u c d s u~ c~ d~ s~ a ve vm vt e- mu- ve~ vm~ vt~ e+ mu+ t b t~ b~ z w+ h w- ta- ta+
MGS>
```

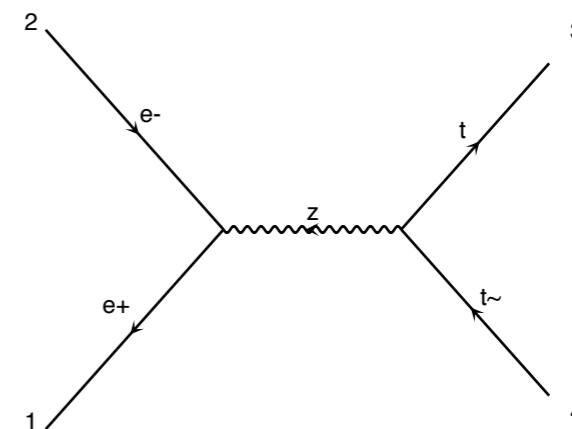
- You can enter **ANY** process!
 - ➔ add [QCD] for NLO functionalities
 - ✦ generate $p p \rightarrow t t^{\sim}$ [QCD]
 - ✦ generate $p p \rightarrow e^+ e^- \mu^+ \mu^-$ [QCD]
 - ✦ generate $e^+ e^- \rightarrow t t^{\sim}$ [QCD]

```
mg5>generate e+ e- > t t~ [QCD]
Switching from interface MG5 to aMC@NLO
The default sm model does not allow to generate loop processes. MG5 now loads 'loop_sm' instead.
import model loop_sm
INFO: load particles
INFO: load vertices
INFO: Restrict model loop_sm with file models/loop_sm/restrict_default.dat .
INFO: Run "set stdout_level DEBUG" before import for more information.
INFO: Change particles name to pass to MG5 convention
Kept definitions of multiparticles l- / j / vl / l+ / p / vl~ unchanged
Defined multiparticle all = g gh gh~ d u s c d~ u~ s~ c~ a ve vm vt e- mu- ve~ vm~ vt~ e+ mu+ b t b~ t~ z w+ h
INFO: Generating FKS-subtracted matrix elements for born process: e+ e- > t t~ [ QCD ]
INFO: Generating virtual matrix elements using MadLoop:
INFO: Generating virtual matrix element with MadLoop for process: e+ e- > t t~ [ QCD ]
INFO: Generated 1 subprocesses with 4 real emission diagrams, 2 born diagrams and 2 virtual diagrams
aMC@NLO>
```

- Born

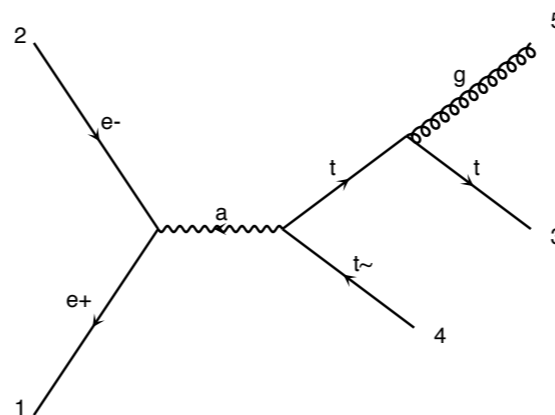


born diagram 1 QCD=0, QED=2

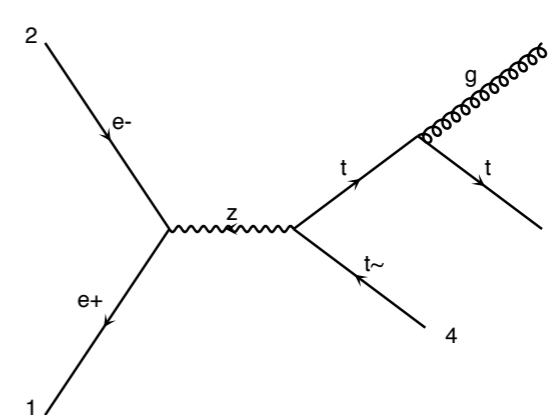


born diagram 2 QCD=0, QED=2

- real



real diagram 1 QCD=1, QED=2



real diagram 2 QCD=1, QED=2

- Virtual

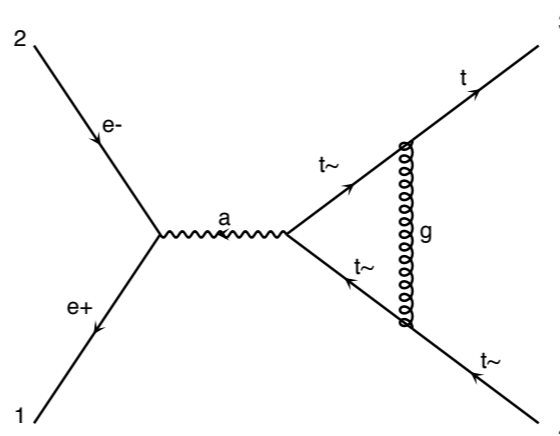


diagram 1 QCD=2, QED=2

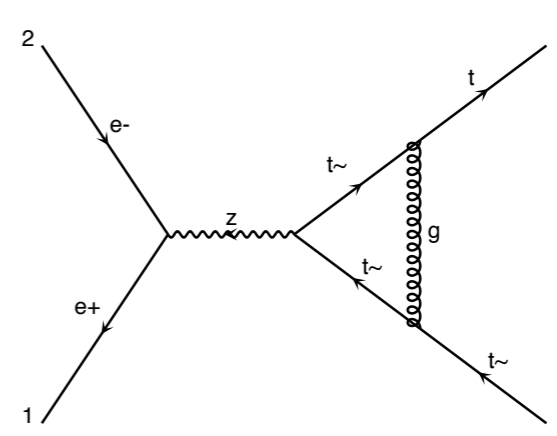


diagram 2 QCD=2, QED=2

- Create your aMC@NLO code
 - ➔ output PATH

- Create your aMC@NLO code

- ➔ output PATH

```
INFO: Writing out the aMC@NLO code, using optimized Loops
INFO: initialize a new directory: PROD_TT2
INFO: remove old information in PROD_TT2
INFO: Generating real emission matrix-elements...
INFO: Generating Helas calls for FKS process: e+ e- > t t~ [ QCD ]
ALOHA: aloha creates FFV1 routines
ALOHA: aloha creates FFV1 routines
ALOHA: aloha creates FFV1 routines
ALOHA: aloha creates FFV2 routines
ALOHA: aloha creates FFV5 routines
INFO: Processing color information for process: e+ e- > t t~ [ QCD ]
INFO: Writing files in P0_epem_ttx
INFO: Creating files in directory V0_epem_ttx
INFO: Computing diagram color coefficients
INFO: Drawing loop Feynman diagrams for Process: e+ e- > t t~ [ QCD ]
INFO: Generating born Feynman diagrams for Process: e+ e- > t t~ [ QCD ]
History written to /Users/omatt/Documents/eclipse/2.0.0beta4/PROD_TT2/Cards/proc_card_mg5.dat
Export UFO model to MG4 format
ALOHA: aloha creates FFV2 routines
ALOHA: aloha creates FFV1 routines
ALOHA: aloha creates FFV4 routines
ALOHA: aloha creates FFV5 routines
ALOHA: aloha creates FFV2_5 routines
ALOHA: aloha creates FFV2_4 routines
INFO: Use Fortran compiler gfortran
INFO: Generate jpeg diagrams
INFO: Generate web pages
The option group_subprocesses is modified [Auto] but will not be written in the configuration files.
If you want to make this value the default for future session, you can run 'save options --all'
The option complex_mass_scheme is modified [False] but will not be written in the configuration files.
If you want to make this value the default for future session, you can run 'save options --all'
save configuration file to /Users/omatt/Documents/eclipse/2.0.0beta4/PROD_TT2/Cards/amcatnlo_configurat
Type "launch" to generate events from this process, or see
/Users/omatt/Documents/eclipse/2.0.0beta4/PROD_TT2/README
Run "open index.html" to see more information about this process.
```


- Create your aMC@NLO code
 - ➔ output PATH
- Run it:
 - ➔ launch [PATH]

```
INFO: *****
*
*      WELCOME to MADGRAPH 5      *
*      aMC@NLO                    *
*
*      *          *                *
*      *    **    *                *
*      *   ** 5 **   *            *
*      *    **    *                *
*      *          *                *
*
*      VERSION 2.0.0.beta4        2013-06-22
*
*      The MadGraph Development Team - Please visit us at
*      http://amcatnlo.cern.ch
*
*      Type 'help' for in-line help.
*
*****
INFO: load configuration from /Users/omatt/.mg5/mg5_configuration.txt
INFO: load configuration from /Users/omatt/Documents/eclipse/2.0.0beta4/PROCNLO_loop_sm_
INFO: load configuration from /Users/omatt/Documents/eclipse/2.0.0beta4/input/mg5_config
INFO: load configuration from /Users/omatt/Documents/eclipse/2.0.0beta4/PROCNLO_loop_sm_
set group_subprocesses Auto
set ignore_six_quark_processes False
set loop_optimized_output True
set gauge unitary
set complex_mass_scheme False
launch auto
The following switches determine which operations are executed:
 1 Perturbative order of the calculation:          order=NLO
 2 Fixed order (no event generation and no MC@N]LO matching):  fixed_order=OFF
 3 Shower the generated events:                    shower=ON
 4 Decay particles with the MadSpin module:        madspin=OFF
  Either type the switch number (1 to 4) to change its default setting,
  or set any switch explicitly (e.g. type 'order=L0' at the prompt)
  Type '0', 'auto', 'done' or just press enter when you are done.
 [0, 1, 2, 3, 4, auto, done, order=L0, order=NLO, ... ][60s to answer]
>
```

- Create your aMC@NLO code
 - ➔ output PATH
- Run it:
 - ➔ launch [PATH]

First Question:

The following switches determine which operations are executed:

- | | |
|---------------------------------------------------------------|-----------------|
| 1 Perturbative order of the calculation: | order=NLO |
| 2 Fixed order (no event generation and no MC@[N]LO matching): | fixed_order=OFF |
| 3 Shower the generated events: | shower=ON |
| 4 Decay particles with the MadSpin module: | madspin=OFF |

Either type the switch number (1 to 4) to change its default setting,
or set any switch explicitly (e.g. type 'order=L0' at the prompt)

Type '0', 'auto', 'done' or just press enter when you are done.

[0, 1, 2, 3, 4, auto, done, order=L0, order=NLO, ...][60s to answer]

>[timer stopped]



- Create your aMC@NLO code
 - ➔ output PATH
- Run it:
 - ➔ launch [PATH]

Second Question:

```
INFO: will run in mode: aMC@NLO
Do you want to edit a card (press enter to bypass editing)?
 1 / param      : param_card.dat
 2 / run        : run_card.dat
 3 / madspin    : madspin_card.dat
 4 / shower     : shower_card.dat
you can also
- enter the path to a valid card or banner.
- use the 'set' command to modify a parameter directly.
  The set option works only for param_card and run_card.
  Type 'help set' for more information on this command.
[0, done, 1, param, 2, run, 3, madspin, 4, enter path, ... ][60s to answer]
>
```

- each beam at 250 GeV

- The code runs:

```
INFO: For gauge cancellation, the width of 't' has been set to zero.
```

```
INFO: Compiling source...
```

```
INFO: ...done, continuing with P* directories
```

```
INFO: Compiling directories...
```

```
INFO: Compiling on 8 cores
```

```
INFO: Compiling P0_epem_ttx...
```

```
INFO: P0_epem_ttx done.
```

```
INFO: Checking test output:
```

```
INFO: P0_epem_ttx
```

```
INFO: Result for test_ME:
```

```
INFO: Passed.
```

```
INFO: Result for test_MC:
```

```
INFO: Passed.
```

```
INFO: Result for check_poles:
```

```
INFO: Poles successfully cancel for 20 points over 20 (tolerance=1.0e-05)
```

```
INFO: Starting run
```

```
INFO: Using 8 cores
```

```
INFO: Cleaning previous results
```

```
INFO: Doing NLO matched to parton shower
```

```
INFO: Setting up grid
```

```
INFO: Idle: 0, Running: 2, Completed: 2 [ current time: 17h19 ]
```

```
INFO: Idle: 0, Running: 1, Completed: 3 [ 1.2s ]
```

```
INFO: Idle: 0, Running: 0, Completed: 4 [ 1.3s ]
```

```
INFO: Determining the number of unweighted events per channel
```

```
Intermediate results:
```

```
Random seed: 36
```

```
Total cross-section: 6.232e-01 +- 4.2e-03 pb
```

```
Total abs(cross-section): 7.010e-01 +- 2.5e-03 pb
```

Compilation

Check Poles cancelation

Integration

```
INFO: Computing upper envelope
INFO: Idle: 0, Running: 2, Completed: 2 [ current time: 17h19 ]
INFO: Idle: 0, Running: 1, Completed: 3 [ 1.3s ]
INFO: Idle: 0, Running: 0, Completed: 4 [ 1.3s ]
INFO: Updating the number of unweighted events per channel

Intermediate results:
Random seed: 36
Total cross-section:      6.183e-01 +- 3.7e-03 pb
Total abs(cross-section): 6.986e-01 +- 1.9e-03 pb
```

```
INFO: Generating events
INFO: Idle: 0, Running: 2, Completed: 2 [ current time: 17h19 ]
INFO: Idle: 0, Running: 1, Completed: 3 [ 0.6s ]
INFO: Idle: 0, Running: 0, Completed: 4 [ 1.3s ]
INFO: Doing reweight
INFO: Idle: 0, Running: 1, Completed: 3 [ current time: 17h19 ]
INFO: Idle: 0, Running: 0, Completed: 4 [ 0.74s ]
INFO: Collecting events
INFO:
Summary:
Process e+ e- > t t~ [QCD]
Run at l-l collider (250 + 250 GeV)
Total cross-section: 6.183e-01 +- 3.7e-03 pb
Ren. and fac. scale uncertainty: +1.0% -0.8%
Number of events generated: 10000
Parton shower to be used: HERWIG6
```

```
decay_events -from_cards
INFO: Running MadSpin
INFO: This functionality allows for the decay of resonances
INFO: in a .lhe file, keeping track of the spin correlation effects.
INFO: BE AWARE OF THE CURRENT LIMITATIONS:
INFO: (1) Only a succession of 2 body decay are currently allowed
*****
*
*      W E L C O M E  t o  M A D S P I N
*
*
*****
```

Integration

Events Generation

Top Decay

```
INFO: Estimating the maximum weight
INFO: *****
INFO: Probing the first 75 events
INFO: with 400 phase space points
INFO:
INFO: Event 1/75 : 0.18s
INFO: Event 6/75 : 1s
INFO: Event 11/75 : 1.9s
INFO: Event 16/75 : 2.8s
INFO: Event 21/75 : 3.8s
INFO: Event 26/75 : 4.9s
INFO: Event 31/75 : 5.8s
INFO: Event 36/75 : 6.7s
INFO: Event 41/75 : 8s
INFO: Event 46/75 : 8.9s
INFO: Event 51/75 : 9.7s
INFO: Event 56/75 : 10.8s
INFO: Event 61/75 : 11.7s
INFO: Event 66/75 : 12.6s
INFO: Event 71/75 : 13.6s
```

estimation of
the maximum
weight

```
INFO:
INFO: Decaying the events...
INFO: Event nb 1000 2.1s
INFO: Event nb 2000 3.8s
INFO: Event nb 3000 5.6s
INFO: Event nb 4000 7.3s
INFO: Event nb 5000 9.1s
INFO: Event nb 6000 10.8s
INFO: Event nb 7000 12.5s
INFO: Event nb 8000 14.4s
INFO: Event nb 9000 16.1s
INFO: Event nb 10000 17.9s
INFO: Total number of events written: 10000/10000
INFO: Average number of trial points per production event: 10.9322
INFO: Branching ratio to allowed decays: 1
INFO: Number of events with weights larger than max_weight: 0
INFO: Number of subprocesses 8
INFO: Number of failures when restoring the Monte Carlo masses: 1
INFO: Decayed events have been written in /Users/omatt/Documents/eclipse/2.0.
INFO: The decayed event file has been moved to the following location:
INFO: /Users/omatt/Documents/eclipse/2.0.0beta4/PROCNLO_loop_sm_16/Events/rur
INFO: MadSpin Done
```

adding decay
event by event

```
INFO: Preparing MCatNLO run
INFO: Compiling MCatNLO for HERWIG6...
INFO: ... done
INFO: Running MCatNLO in /Users/omatt/Documents/eclipse/2.0.0beta4/PROCNLO_lc
INFO: The file /Users/omatt/Documents/eclipse/2.0.0beta4/PROCNLO_loop_sm_16/E
It contains showered and hadronized events in the StdHEP format obtained show
cayed_1/events.lhe.gz with HERWIG6
quit
```

Shower

DEMO

Is it really automatic?

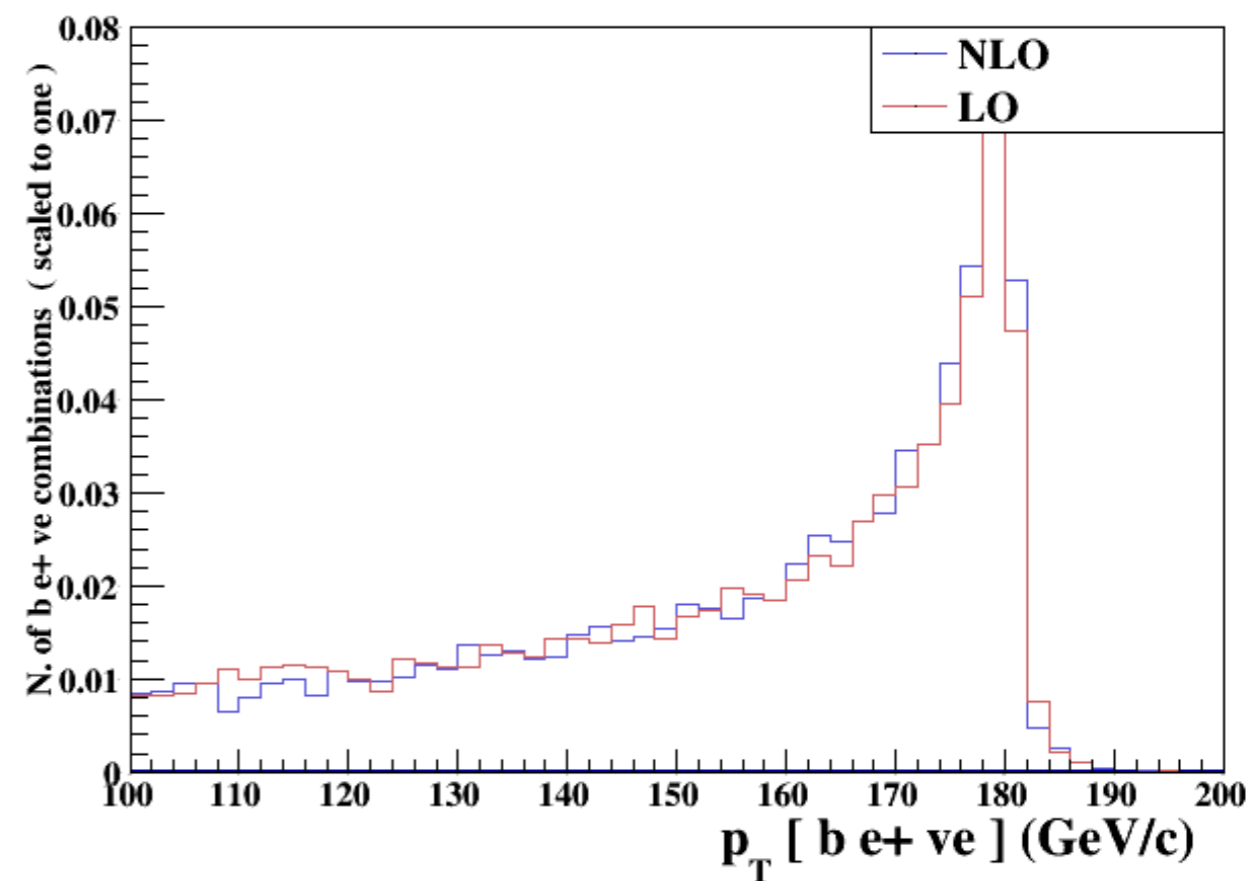
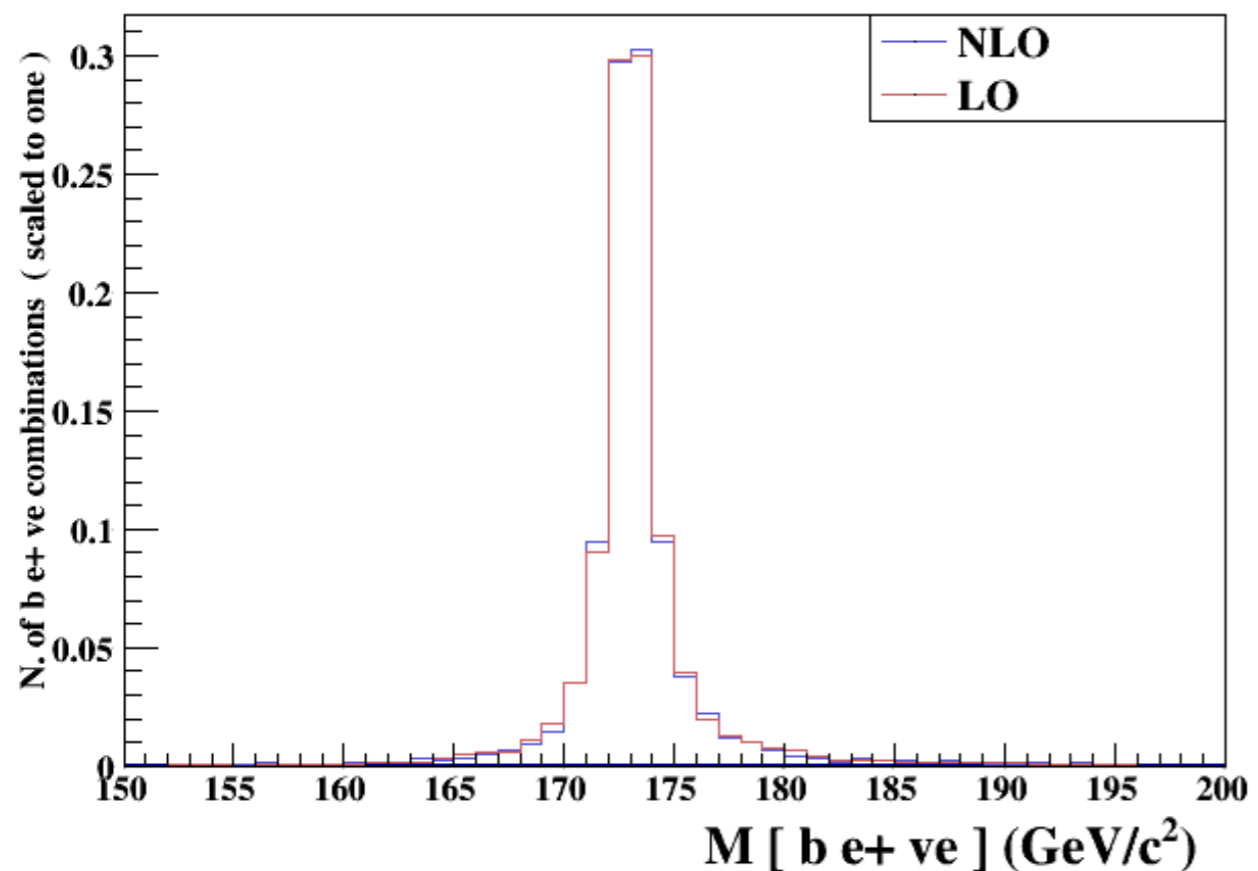
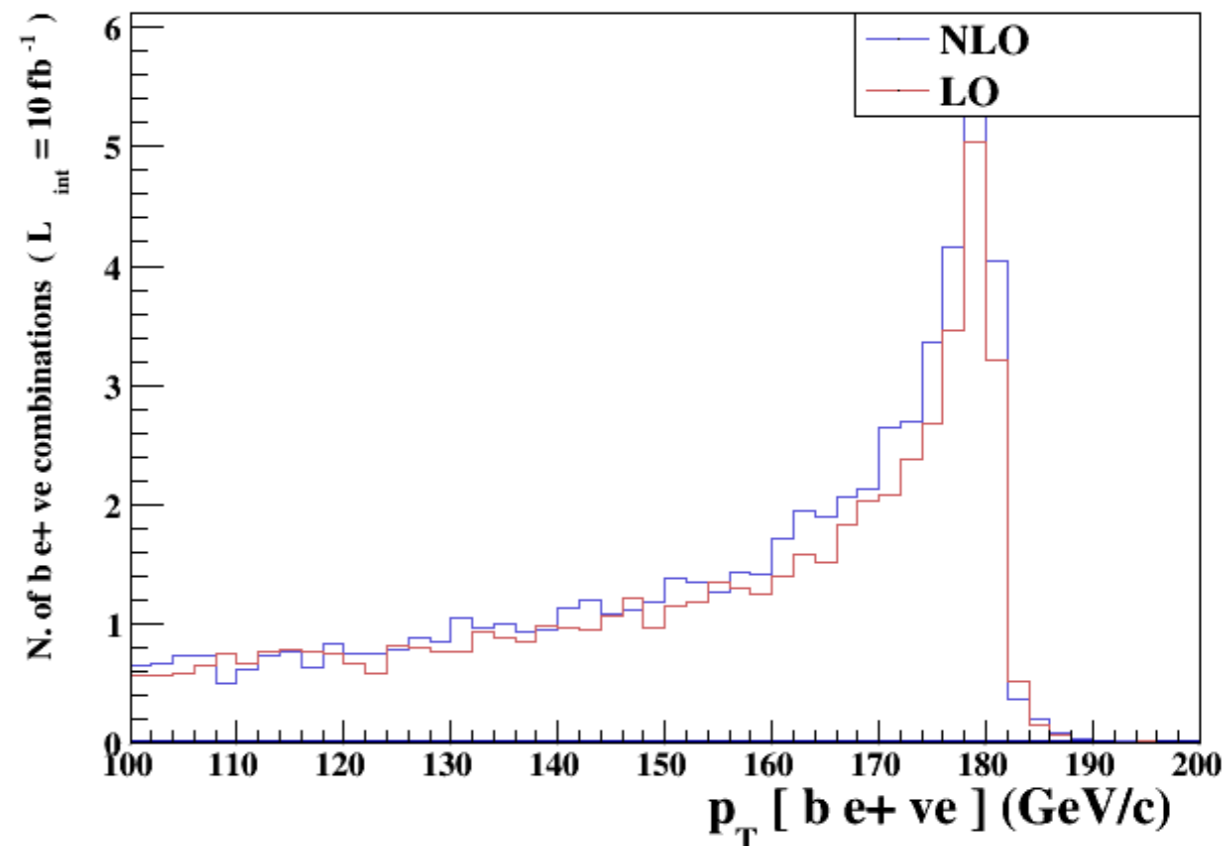
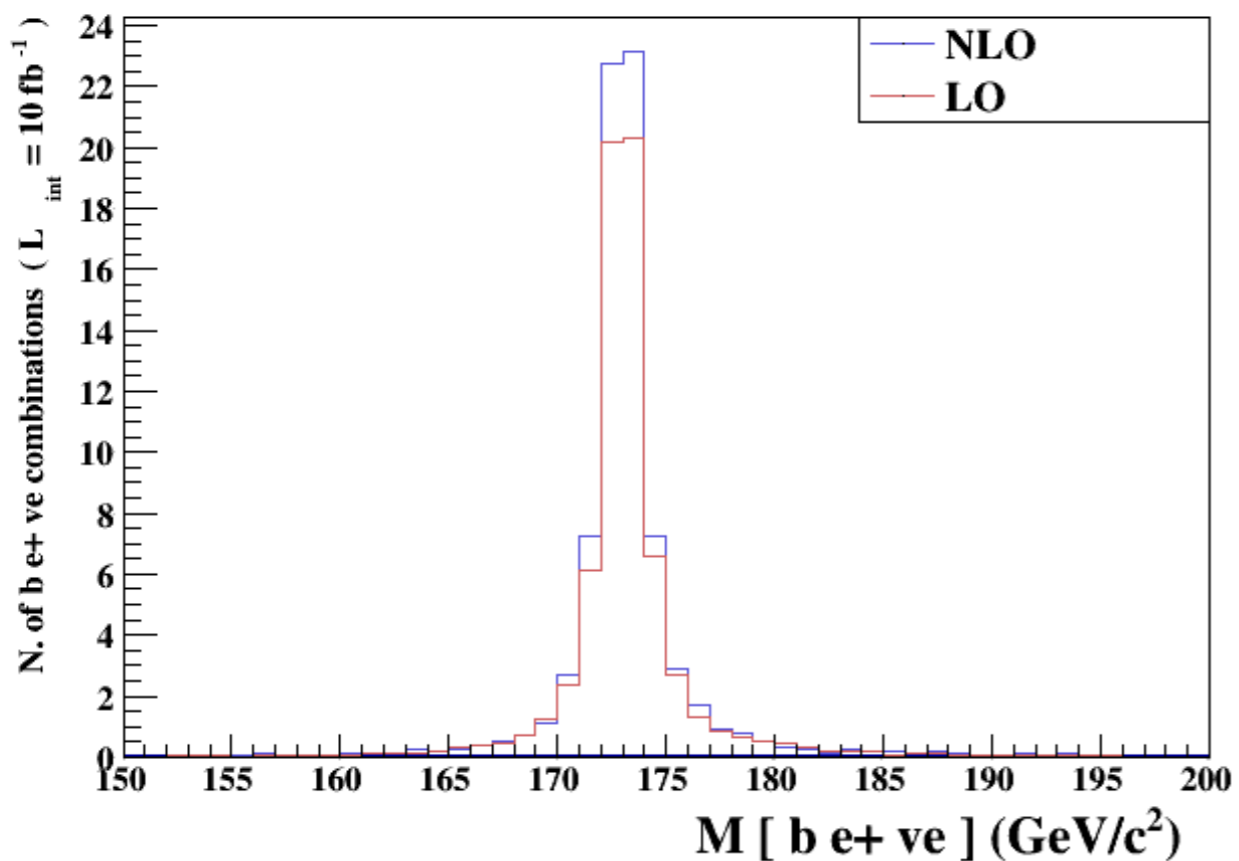
DEMO

Is it really automatic?

As much as LO!

Top-quark pair production at ILC

Preliminary



Offshell effect at NLO

- Diagrams with unstable particles present in general an imaginary part in the Dyson-resummed propagator:

$$P(p) = [p^2 - m_0^2 + Pi(p^2)]^{-1}$$

- Mixing of different perturbative orders breaks gauge invariance. Fine cancellations spoiled, leading to enhanced violation of unitarity
- **No pole cancelation** at NLO for fix-width scheme
- Solution: Complex Mass-Scheme: $M \rightarrow \sqrt{M^2 - iM\Gamma}$,

$$c_W^2 = \frac{M_w^2 + iM_W\Gamma_W}{M_Z^2 + iM_Z\Gamma_Z}$$

Gauge dependence at LO

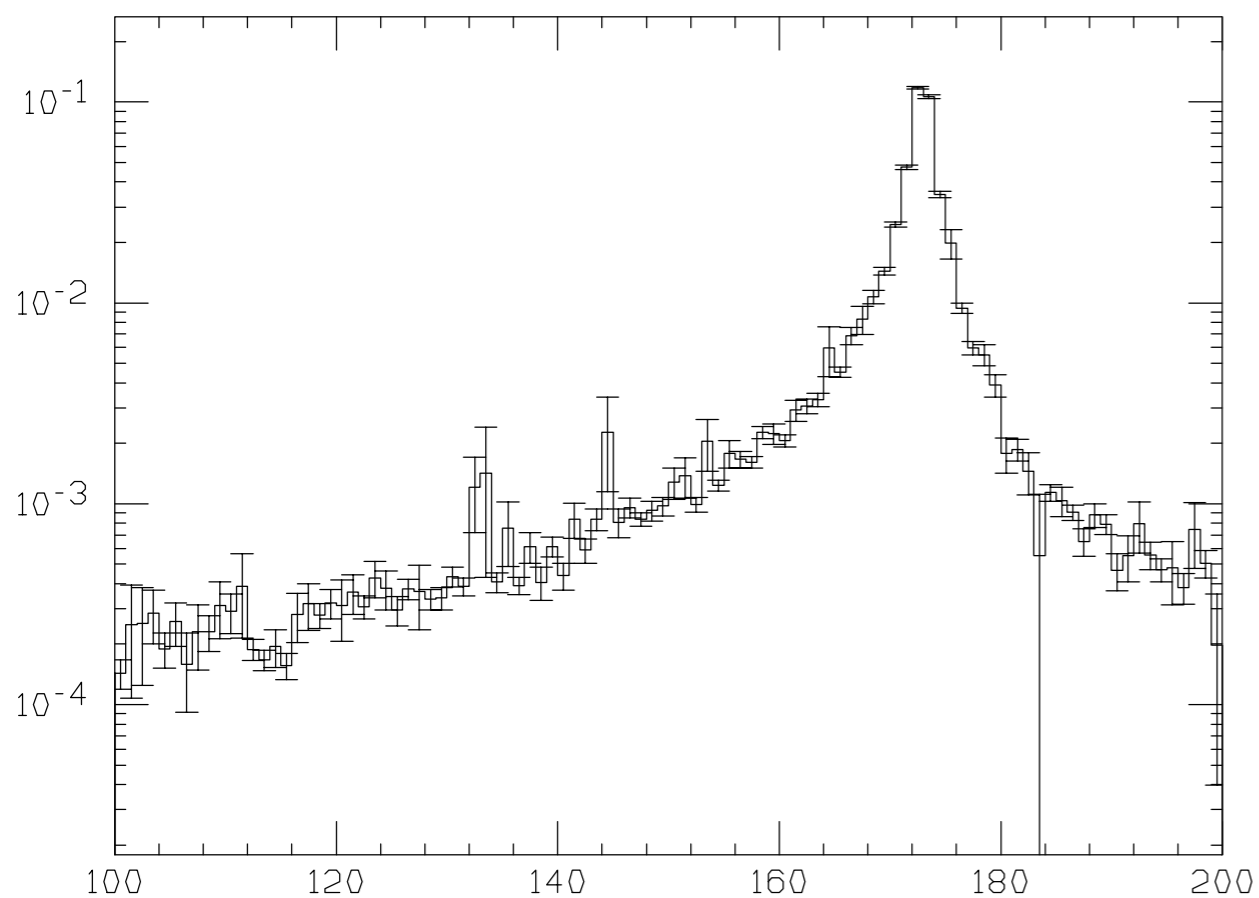
$ A ^2 - \text{Feynman-unitary} /\text{unitary}$	complex mass	fixed width
$e^+ e^- \rightarrow u\bar{u}d\bar{d}$	1.5334067678e-15	1.2312200197e-09
$u\bar{u} \rightarrow u\bar{u}d\bar{d}$	2.0862057616e-16	2.7696013365e-10
$u\bar{u} \rightarrow b\bar{b}e^+ \nu_e \mu^- \nu_\mu$ (real Yuk)	1.7934842084e-06	2.2832833007e-05
" (complex Yuk)	8.5986902303e-16	2.2832833007e-05

$\sigma(pb)$ for $gg \rightarrow b\bar{b}e^+ \nu_e \mu^- \nu_\mu$			
gauge - scheme	complex-mass	fix width	no width
feynman	$1.796e-05 \pm 2.3e-08$	$1.787e-05 \pm 2.5e-08$	
unitary	$1.792e-05 \pm 2.1e-08$	$1.778e-05 \pm 2.4e-08$	$1.810e-05 \pm 2.4e-08$

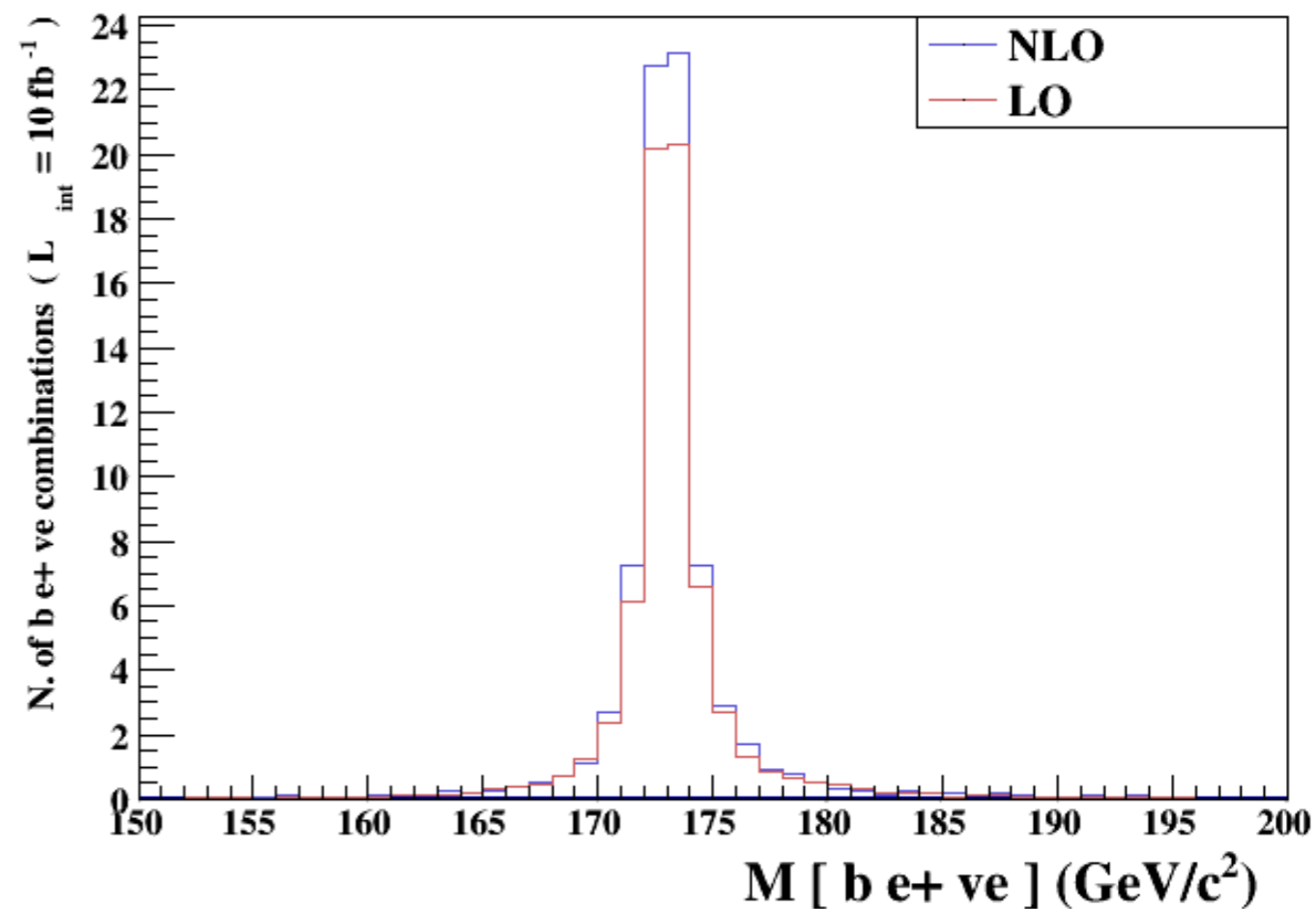
Offshell effect at NLO

$$e^+ e^- \rightarrow w^+ w^- b \bar{b}$$

M t NLO



$$e^+ e^- \rightarrow t \bar{t}$$



Conclusion

- aMC@NLO is
 - ➔ public
 - ➔ automatic
 - ➔ flexible
- MadSpin
 - ➔ decay with full spin correlations
 - ➔ keep finite width effect
- complex-mass
- This is only the beginning of this Tool!

Process	μ	n_{lf}	Cross section (pb)	
			LO	NLO
a.1 $pp \rightarrow t\bar{t}$	m_{top}	5	123.76 ± 0.05	162.08 ± 0.12
a.2 $pp \rightarrow tj$	m_{top}	5	34.78 ± 0.03	41.03 ± 0.07
a.3 $pp \rightarrow tjj$	m_{top}	5	11.851 ± 0.006	13.71 ± 0.02
a.4 $pp \rightarrow t\bar{b}j$	$m_{top}/4$	4	25.62 ± 0.01	30.96 ± 0.06
a.5 $pp \rightarrow t\bar{b}jj$	$m_{top}/4$	4	8.195 ± 0.002	8.91 ± 0.01
b.1 $pp \rightarrow (W^+ \rightarrow)e^+\nu_e$	m_W	5	5072.5 ± 2.9	6146.2 ± 9.8
b.2 $pp \rightarrow (W^+ \rightarrow)e^+\nu_e j$	m_W	5	828.4 ± 0.8	1065.3 ± 1.8
b.3 $pp \rightarrow (W^+ \rightarrow)e^+\nu_e jj$	m_W	5	298.8 ± 0.4	300.3 ± 0.6
b.4 $pp \rightarrow (\gamma^*/Z \rightarrow)e^+e^-$	m_Z	5	1007.0 ± 0.1	1170.0 ± 2.4
b.5 $pp \rightarrow (\gamma^*/Z \rightarrow)e^+e^- j$	m_Z	5	156.11 ± 0.03	203.0 ± 0.2
b.6 $pp \rightarrow (\gamma^*/Z \rightarrow)e^+e^- jj$	m_Z	5	54.24 ± 0.02	56.69 ± 0.07
c.1 $pp \rightarrow (W^+ \rightarrow)e^+\nu_e b\bar{b}$	$m_W + 2m_b$	4	11.557 ± 0.005	22.95 ± 0.07
c.2 $pp \rightarrow (W^+ \rightarrow)e^+\nu_e t\bar{t}$	$m_W + 2m_{top}$	5	0.009415 ± 0.000003	0.01159 ± 0.00001
c.3 $pp \rightarrow (\gamma^*/Z \rightarrow)e^+e^- b\bar{b}$	$m_Z + 2m_b$	4	9.459 ± 0.004	15.31 ± 0.03
c.4 $pp \rightarrow (\gamma^*/Z \rightarrow)e^+e^- t\bar{t}$	$m_Z + 2m_{top}$	5	0.0035131 ± 0.0000004	0.004876 ± 0.000002
c.5 $pp \rightarrow \gamma t\bar{t}$	$2m_{top}$	5	0.2906 ± 0.0001	0.4169 ± 0.0003
d.1 $pp \rightarrow W^+W^-$	$2m_W$	4	29.976 ± 0.004	43.92 ± 0.03
d.2 $pp \rightarrow W^+W^- j$	$2m_W$	4	11.613 ± 0.002	15.174 ± 0.008
d.3 $pp \rightarrow W^+W^+ jj$	$2m_W$	4	0.07048 ± 0.00004	0.1377 ± 0.0005
e.1 $pp \rightarrow HW^+$	$m_W + m_H$	5	0.3428 ± 0.0003	0.4455 ± 0.0003
e.2 $pp \rightarrow HW^+ j$	$m_W + m_H$	5	0.1223 ± 0.0001	0.1501 ± 0.0002
e.3 $pp \rightarrow HZ$	$m_Z + m_H$	5	0.2781 ± 0.0001	0.3659 ± 0.0002
e.4 $pp \rightarrow HZ j$	$m_Z + m_H$	5	0.0988 ± 0.0001	0.1237 ± 0.0001
e.5 $pp \rightarrow Ht\bar{t}$	$m_{top} + m_H$	5	0.08896 ± 0.00001	0.09869 ± 0.00003
e.6 $pp \rightarrow Hb\bar{b}$	$m_b + m_H$	4	0.16510 ± 0.00009	0.2099 ± 0.0006
e.7 $pp \rightarrow Hjj$	m_H	5	1.104 ± 0.002	1.036 ± 0.002

Work in Progress in aMC@NLO

What to expect in the future

Perspectives

Perspectives

- FeynRules@NLO:

Perspectives

- FeynRules@NLO:
 - ➔ NLO not only for the SM but for New Physics

Perspectives

- FeynRules@NLO:
 - ➔ NLO not only for the SM but for New Physics
- ElectroWeak corrections (matched to the shower)

Perspectives

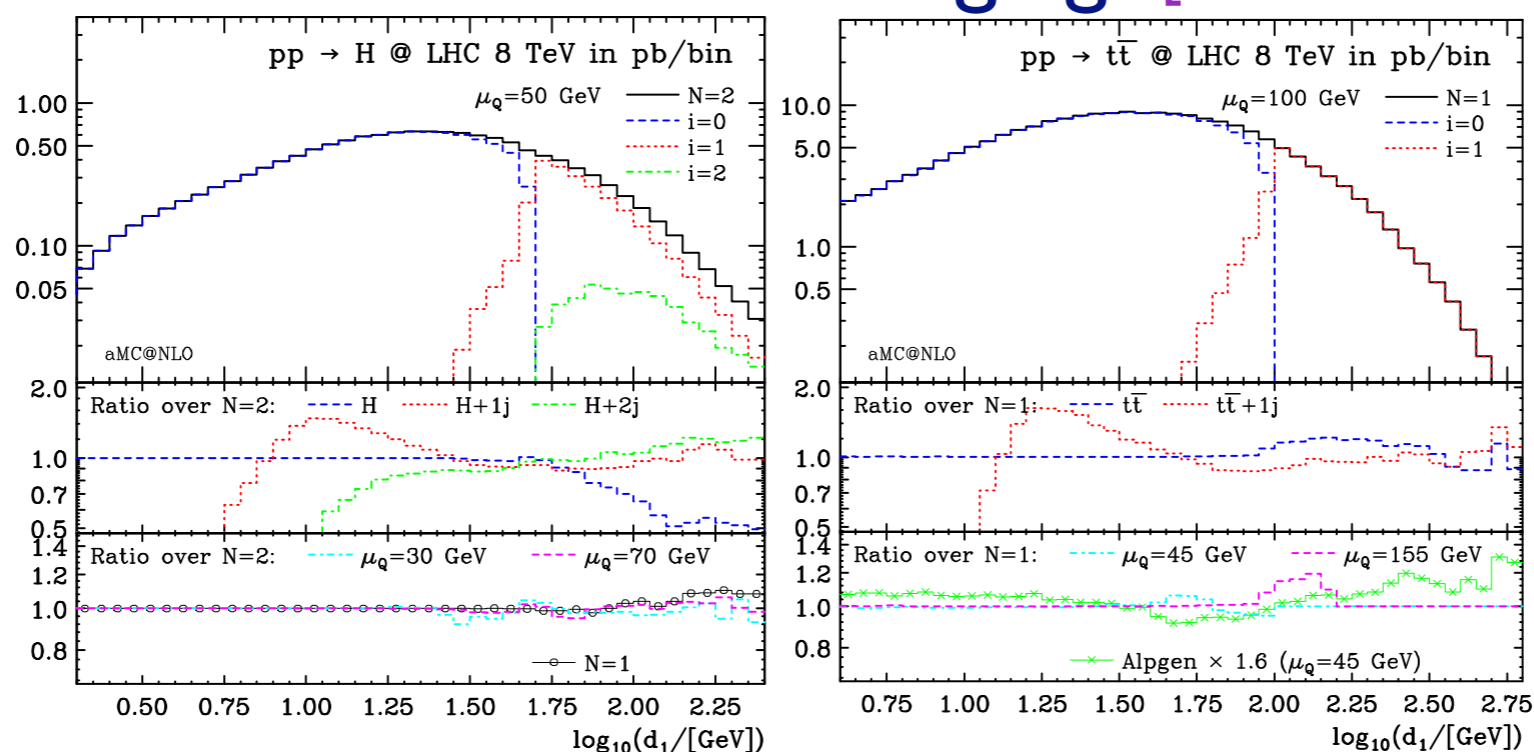
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 - ➔ MadLoop ready (currently in validation)

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- Full automation of FxFx merging [R. Frederix, S. Frixione (2012)]



$0 \rightarrow 1$ rates in H^0 and $t\bar{t}$ production

Perspectives

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- Full automation of FxFx merging [R. Frederix, S. Frixione (2012)]
- Automation of loop-induced processes
- Interface to Pythia8
- Complex mass scheme

MC@NLO properties

- Good features of including the subtraction counter terms
 1. **Double counting avoided:** The rate expanded at NLO coincides with the total NLO cross section
 2. **Smooth matching:** MC@NLO coincides (in shape) with the parton shower in the soft/collinear region, while it agrees with the NLO in the hard region
 3. **Stability:** weights associated to different multiplicities are separately finite. The **MC** term has the same infrared behavior as the real emission (there is a subtlety for the soft divergence)
- Not so nice feature (for the developer):
 1. **Parton shower dependence:** the form of the **MC** terms depends on what the parton shower does exactly. Need special subtraction terms for each parton shower to which we want to match