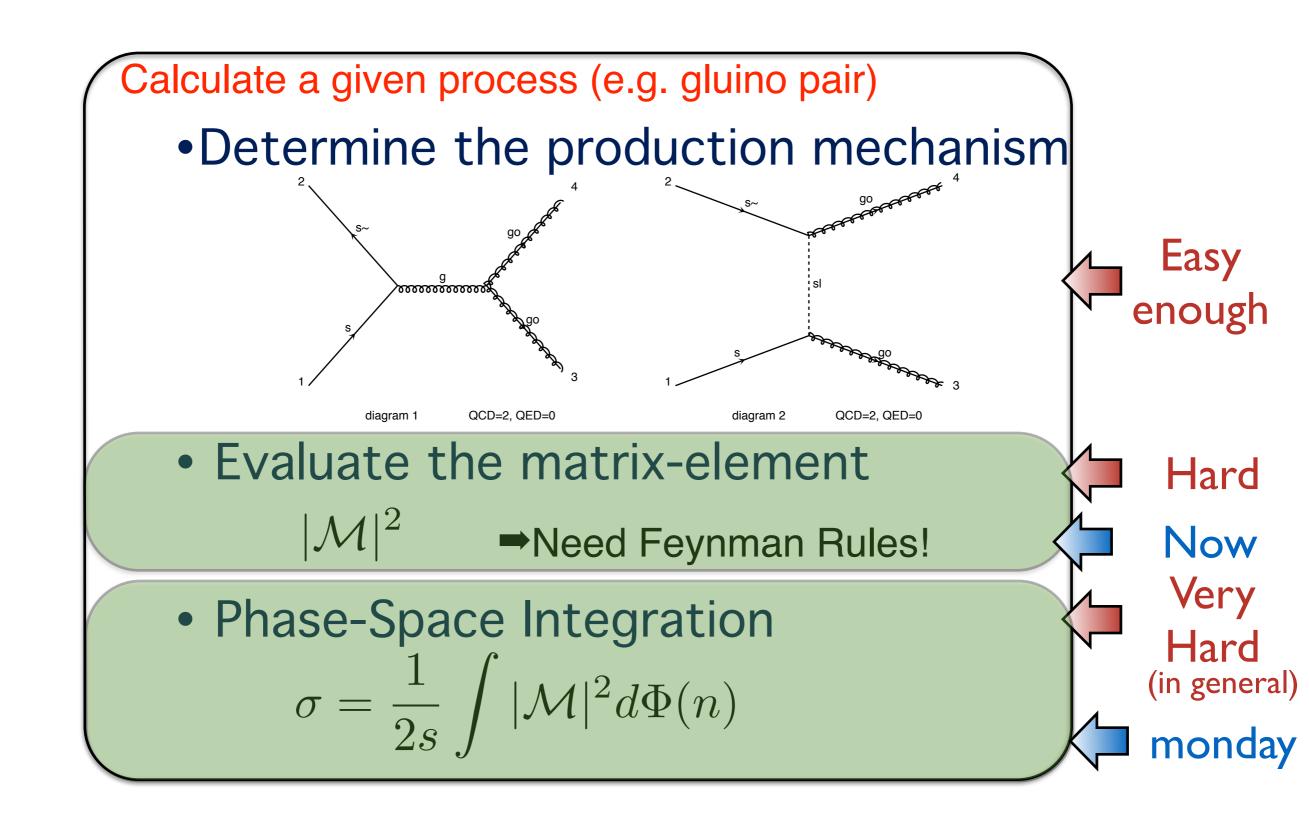
MadGraph5

Olivier Mattelaer IPPP/Durham



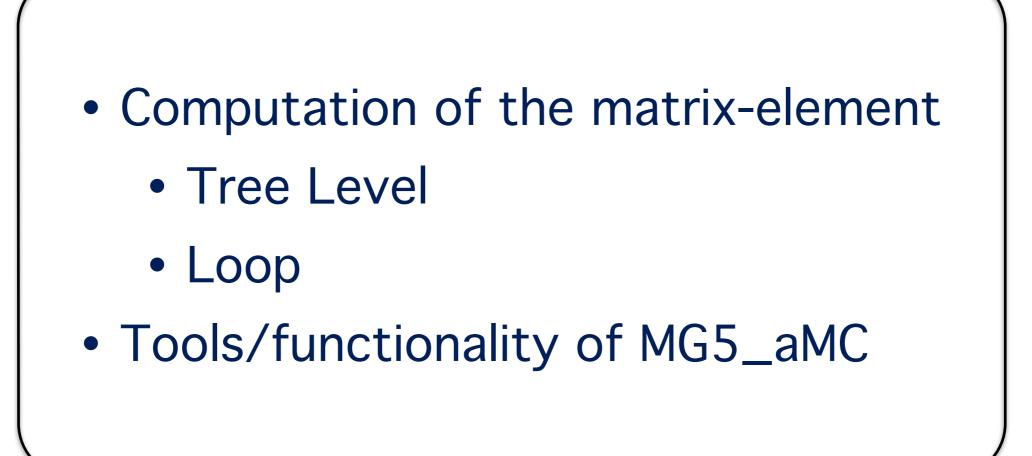
Matrix-Element







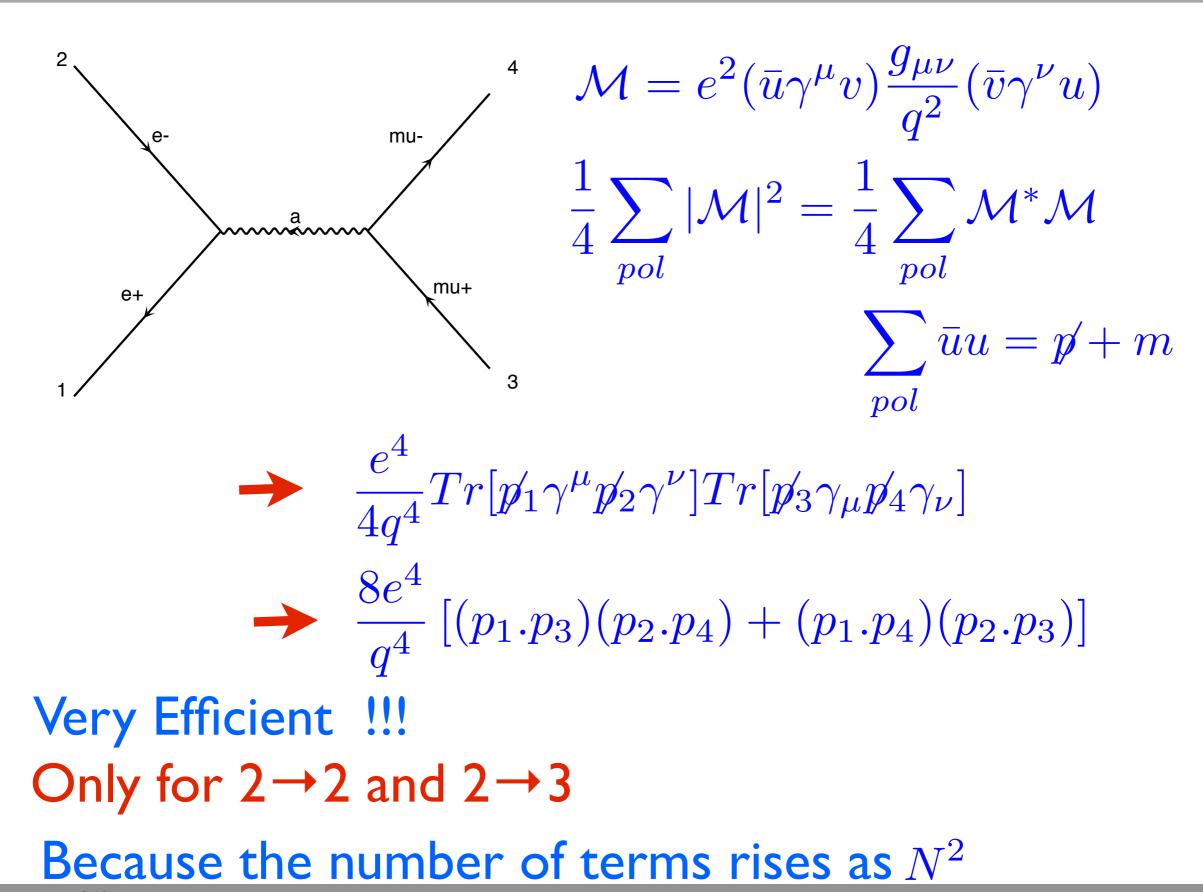






Matrix Element

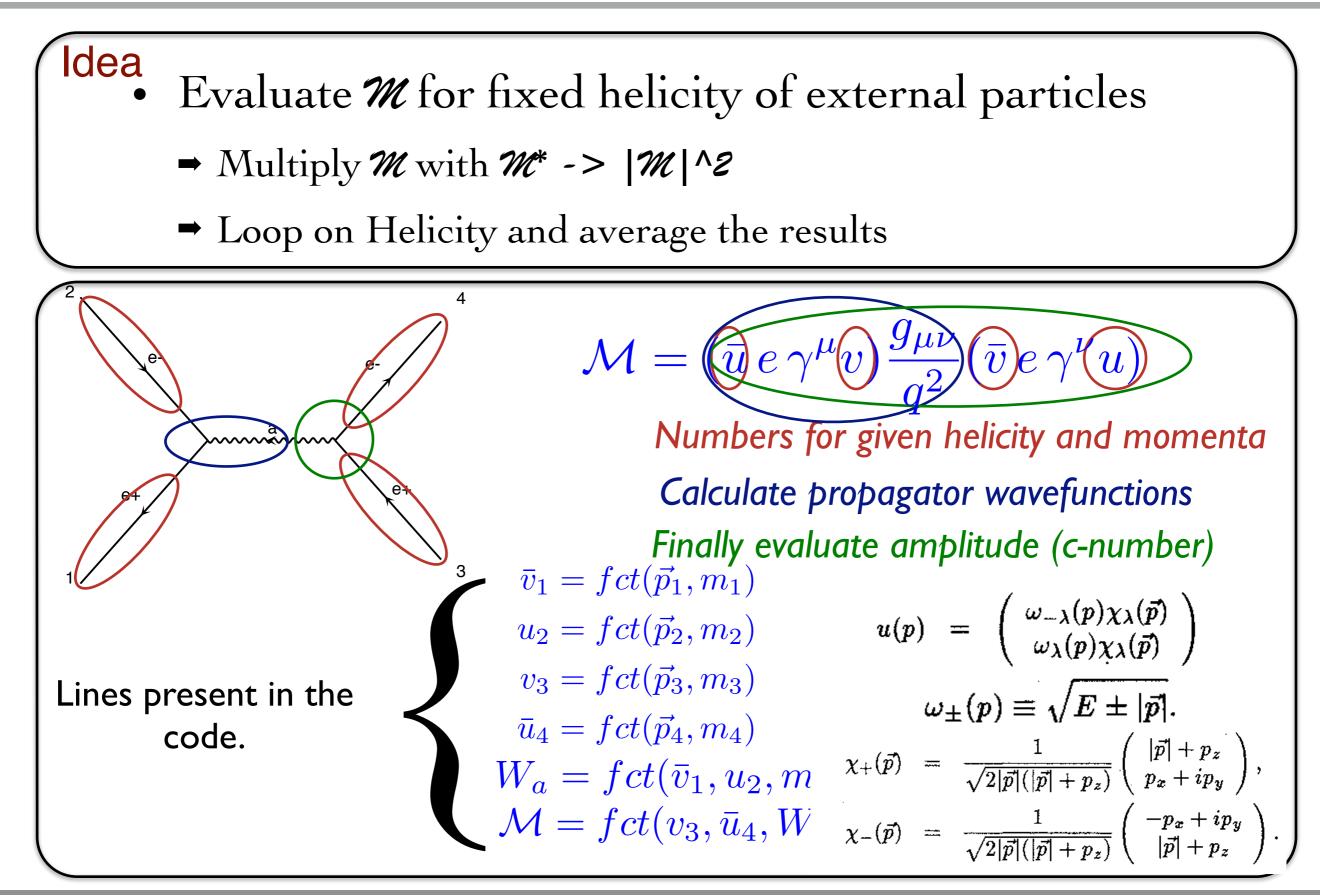




Mattelaer Olivier





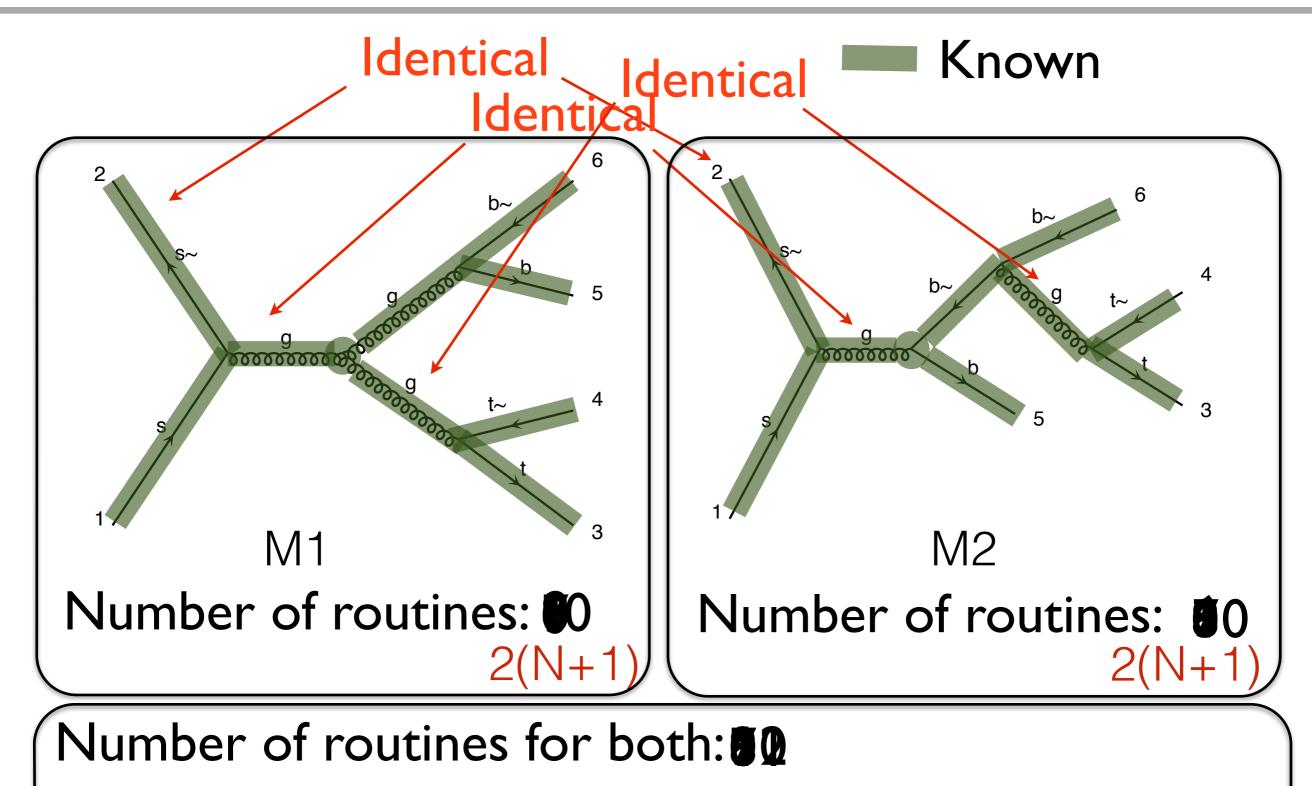


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Real case





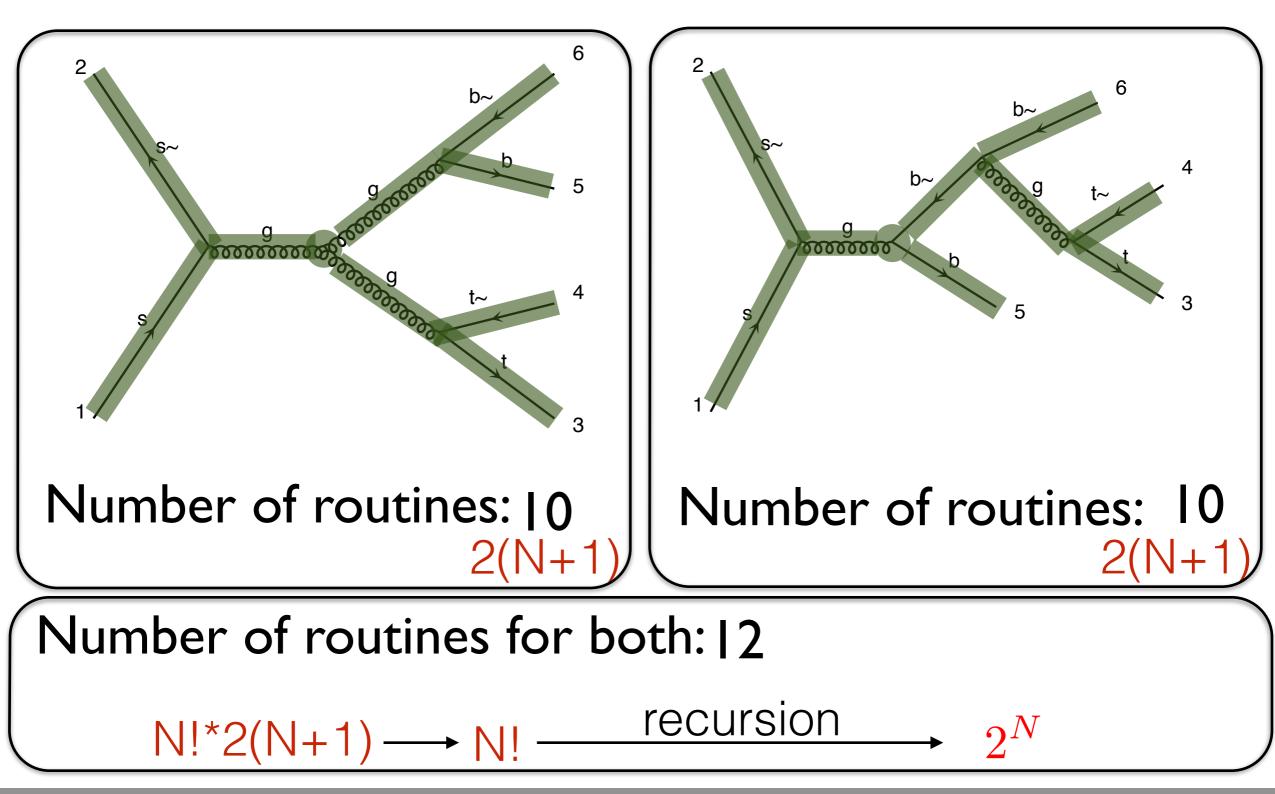
$$|M|^2 = |M_1 + M_2|^2$$



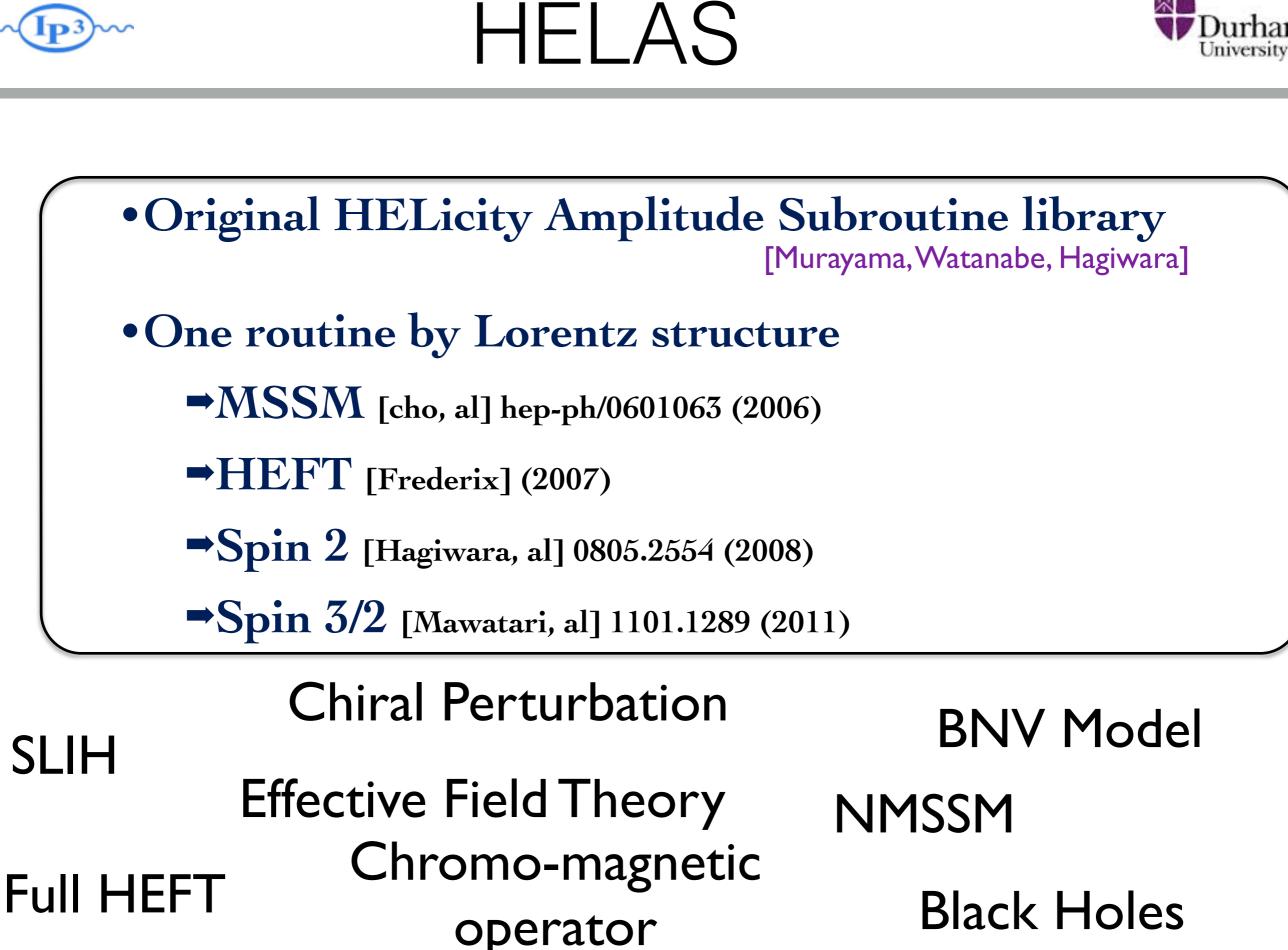
Real case















ALOHA



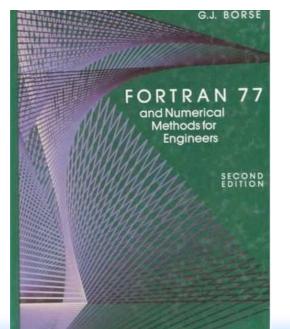
From: UFO 🔽 🔄

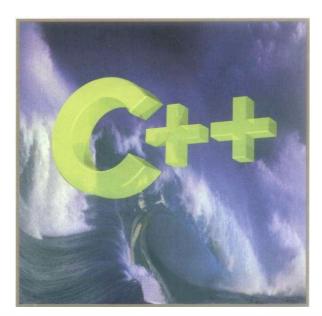
🗧 📷 Helicity

Translate

Basically, any new operator can be handle by MG5/Pythia8 out of the box!

Type text or a website address or translate a document.





PYTHON programming





WESLEY J. CHUN

Brussels October 2010





To Remember





- We are able to compute matrix-element
 - for large number of final state
 - for any BSM theory



Loop Computation

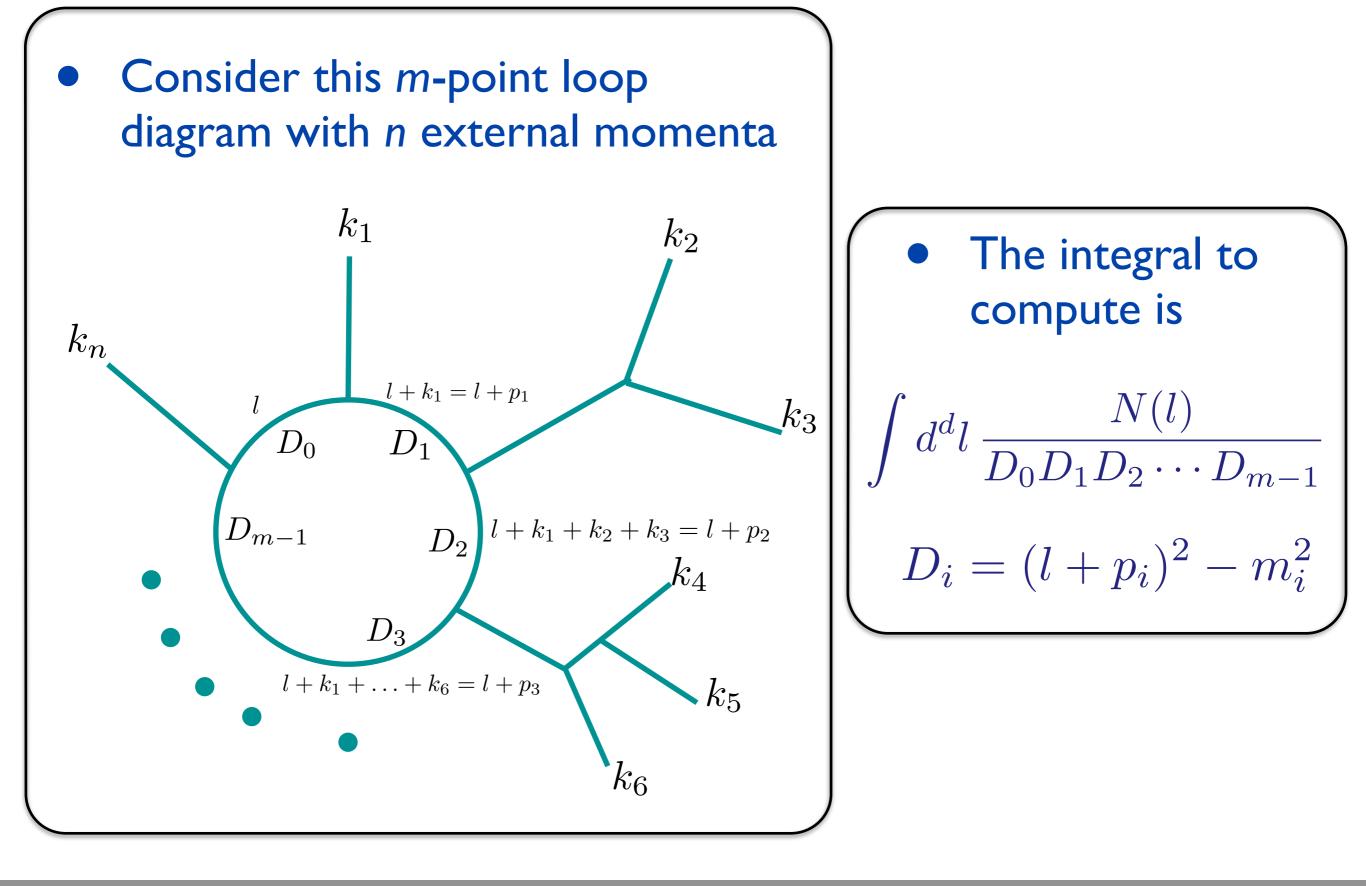


Loop Computation



One Loop











- Any one-loop integral can be decomposed in scalar integrals
- The task is to find these coefficients efficiently (analytically or numerically)

- Basis of scalar integrals

• The a, b, c, d and R coefficients depend only on external parameters and momenta

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

$$\mathcal{D}_i = (l + p_i)^2 - m_i^2 \\ \operatorname{Tadpole}_{i_0} = \int d^d l \frac{1}{D_{i_0}} \\ \operatorname{Bubble}_{i_0 i_1} = \int d^d l \frac{1}{D_{i_0} D_{i_1}} \\ \operatorname{Triangle}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i_1} D_{i_2} D_{i_1} D_{i_1} D_{i_1} D_{i_2}} \\ \operatorname{Box}_{i_0 i_1 i_2 i_3} = \int d^d l \frac{1}{D_{i_0} D_{i_1} D_{i$$

• All these scalar integrals are known and available in computer libraries (FF [v. Oldenborgh], QCDLoop [Ellis, Zanderighi], OneLOop [v. Hameren])





• The a, b, c, d and R coefficients depend only on external parameters and momenta

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

$$D_{i} = (l + p_{i})^{2} - m_{i}^{2}$$

$$Tadpole_{i_{0}} = \int d^{d}l \frac{1}{D_{i_{0}}}$$

$$Bubble_{i_{0}i_{1}} = \int d^{d}l \frac{1}{D_{i_{0}}D_{i_{1}}}$$

$$Triangle_{i_{0}i_{1}i_{2}} = \int d^{d}l \frac{1}{D_{i_{0}}D_{i_{1}}D_{i_{2}}}$$

$$Box_{i_{0}i_{1}i_{2}i_{3}} = \int d^{d}l \frac{1}{D_{i_{0}}D_{i_{1}}D_{i_{2}}}$$

- The coefficients d, c, b and a are finite and do not contain poles in $1/\epsilon$
- The I/ϵ dependence is in the scalar integrals (and the UV renormalization)
- When we have solved this system (and included the UV renormalization) we have the full dependence on the soft/collinear divergences in terms of coefficients in front of the poles. These divergences should cancel against divergences in the real emission corrections (according to KLN theorem)

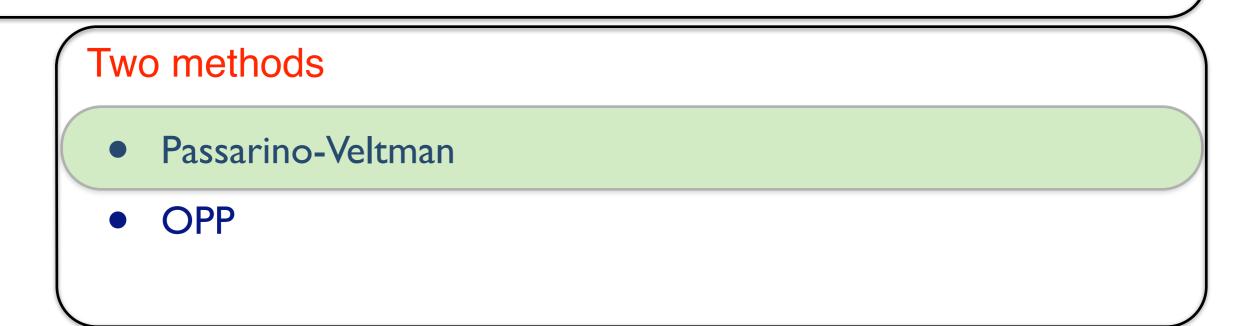
Virtual
$$\sim v_0 + \frac{v_1}{\epsilon} + \frac{v_2}{\epsilon^2}$$





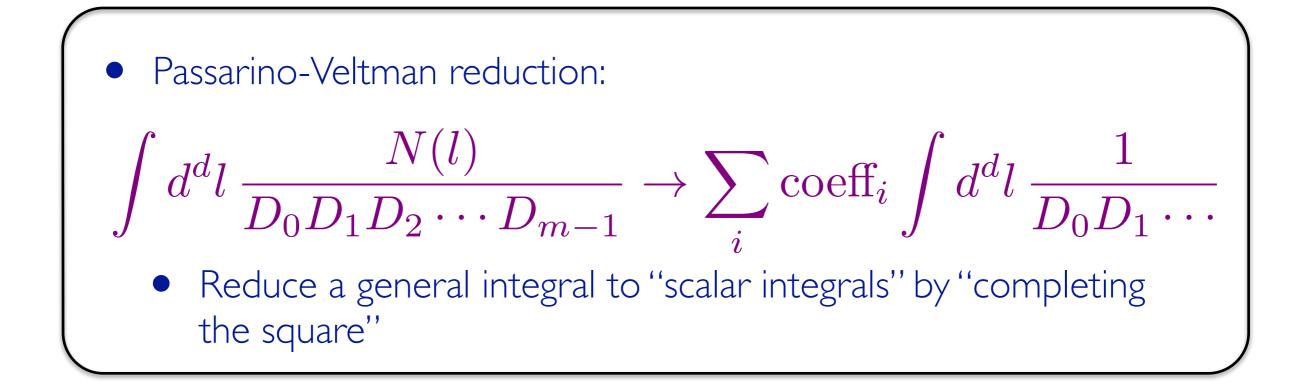


- Any one-loop integral can be decomposed in scalar integrals
- The task is to find these coefficients efficiently (analytically or numerically)







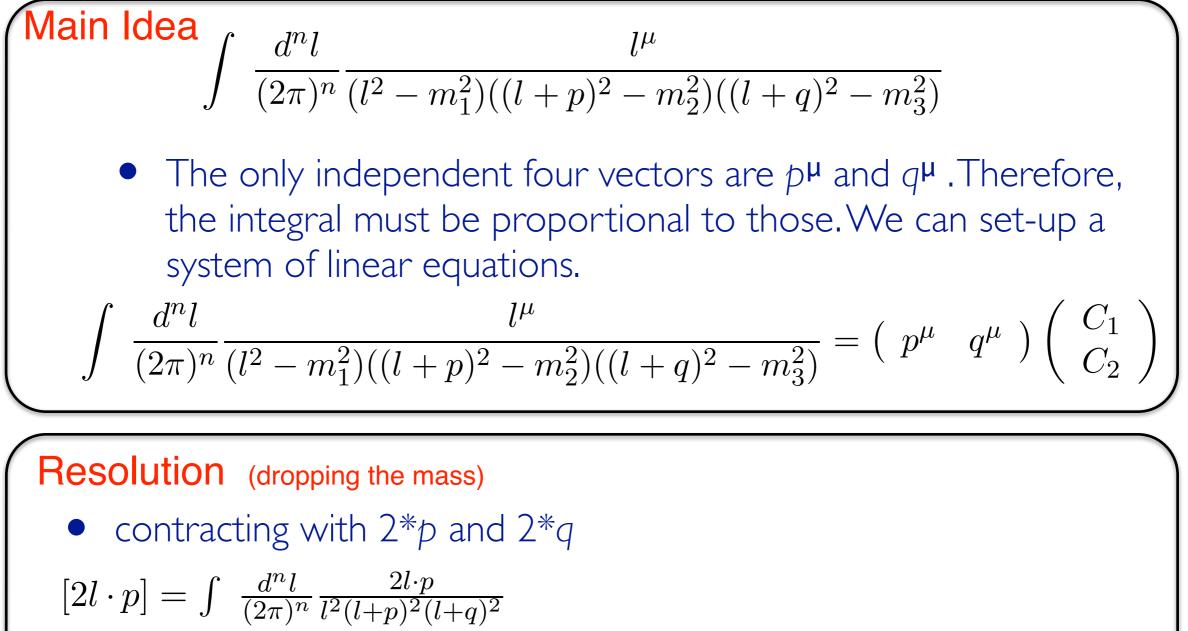


• Let's do an example:
Suppose we want to calculate this triangle integral
$$q = \int \frac{l}{(2\pi)^n} \frac{d^n l}{(l^2 - m_1^2)((l+p)^2 - m_2^2)((l+q)^2 - m_3^2)}$$



Passarino-Veltman



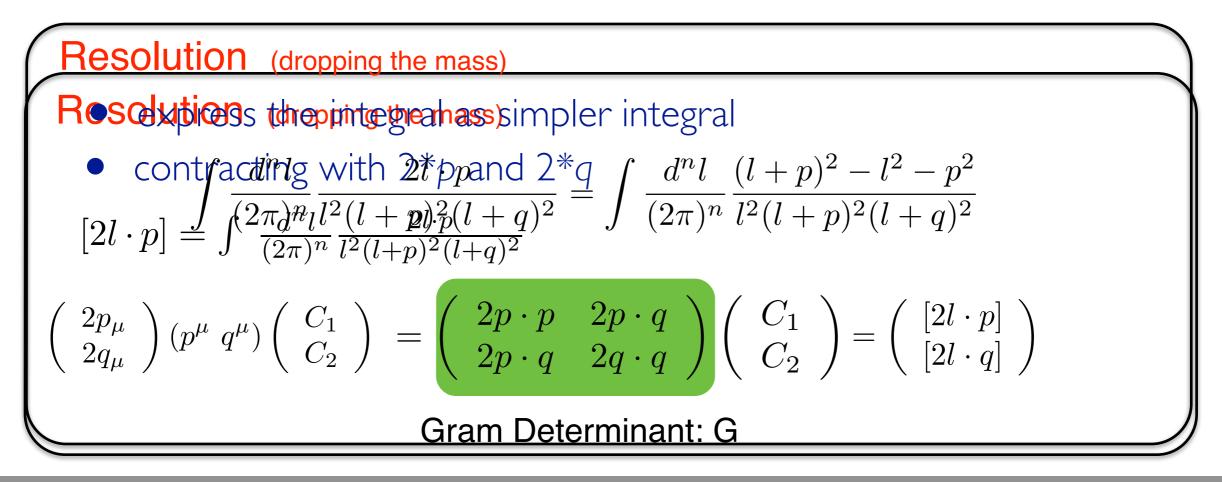


$$\begin{pmatrix} 2p_{\mu} \\ 2q_{\mu} \end{pmatrix} (p^{\mu} q^{\mu}) \begin{pmatrix} C_{1} \\ C_{2} \end{pmatrix} = \begin{pmatrix} 2p \cdot p & 2p \cdot q \\ 2p \cdot q & 2q \cdot q \end{pmatrix} \begin{pmatrix} C_{1} \\ C_{2} \end{pmatrix} = \begin{pmatrix} [2l \cdot p] \\ [2l \cdot q] \end{pmatrix}$$

Gram Determinant: G



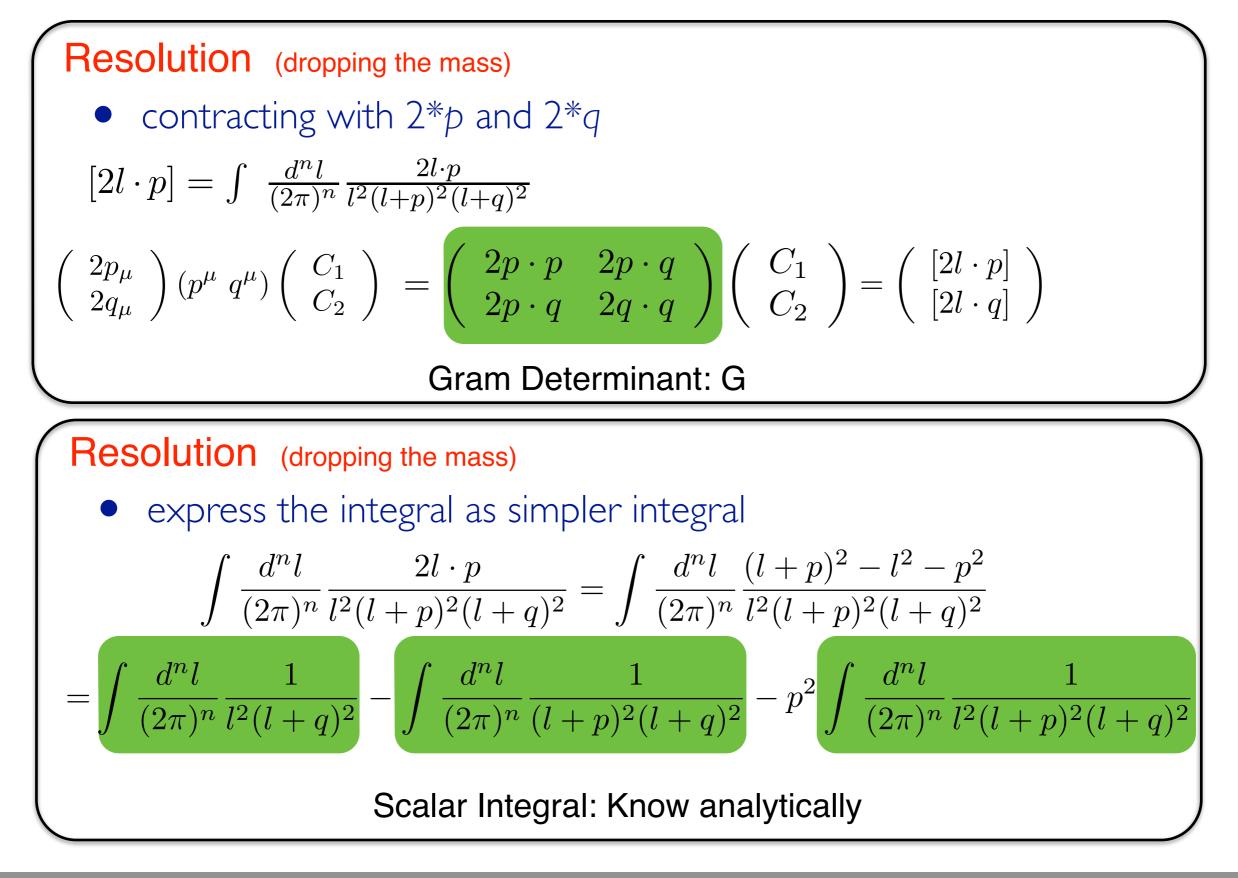






Passarino-Veltman

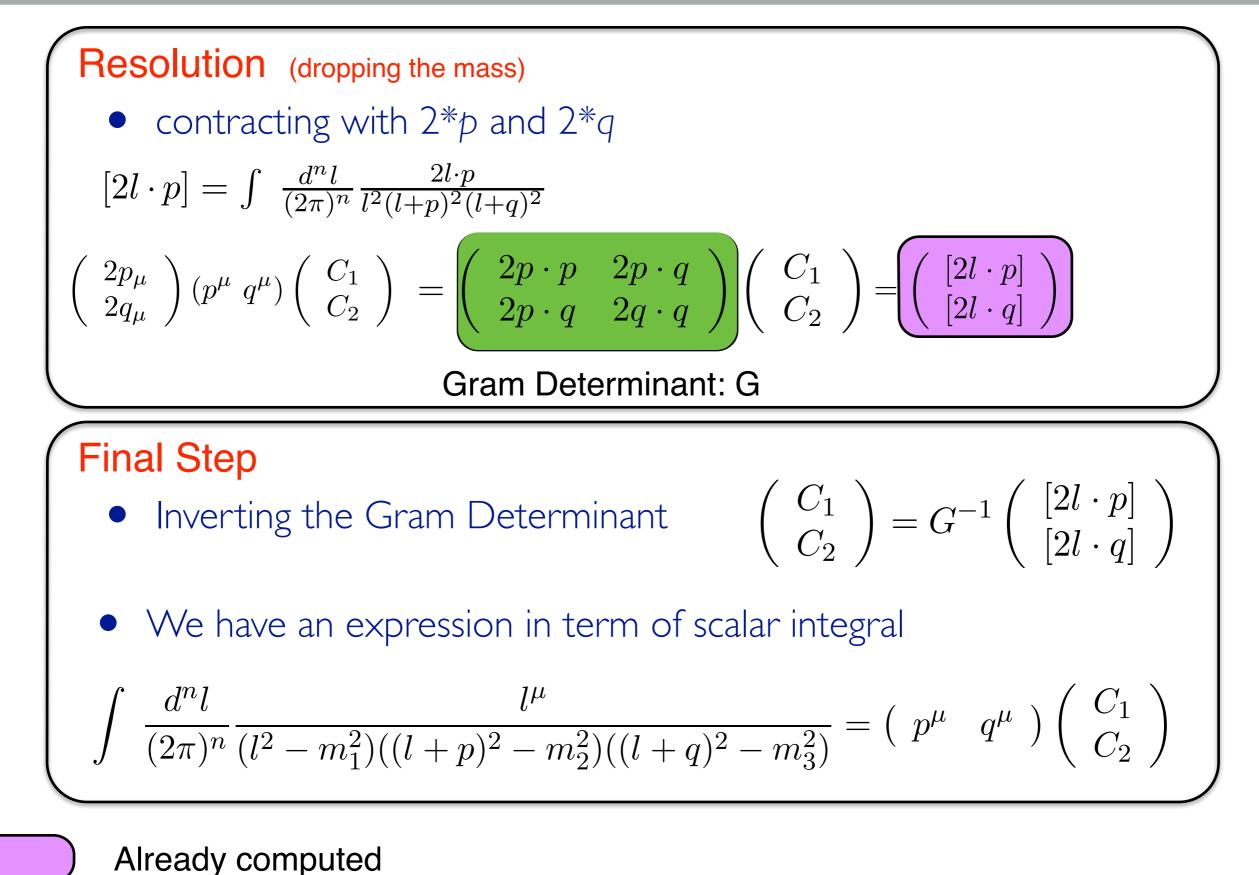






Passarino-Veltman



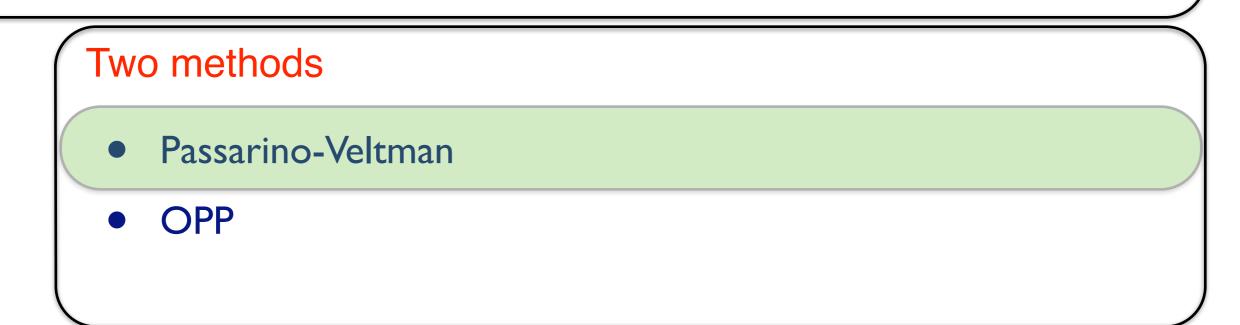








- Any one-loop integral can be decomposed in scalar integrals
- The task is to find these coefficients efficiently (analytically or numerically)



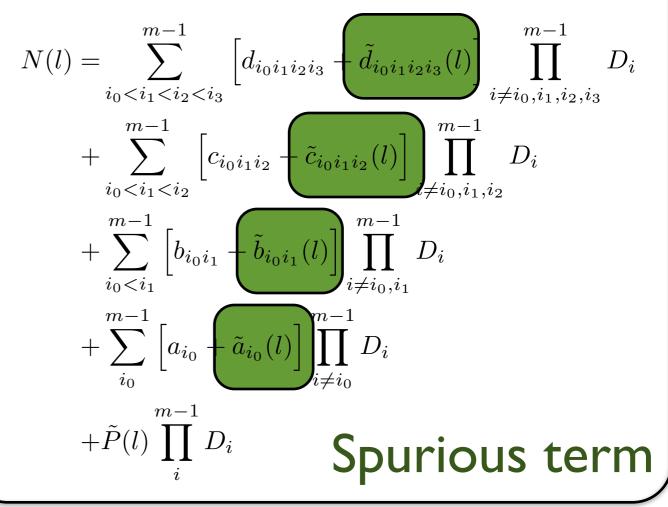




 The decomposition to scalar integrals presented before works at the level of the integrals

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

If we would know a similar relation at the **integrand** level, we would be able to manipulate the integrands and extract the coefficients without doing the integrals





- The functional form of the spurious terms is known (it depends on the rank of the integral and the number of propagators in the loop) [del Aguila, Pittau 2004]
 - for example, a box coefficient from a rank I numerator is

$$\tilde{d}_{i_0 i_1 i_2 i_3}(l) = \tilde{d}_{i_0 i_1 i_2 i_3} \epsilon^{\mu\nu\rho\sigma} l^{\mu} p_1^{\nu} p_2^{\rho} p_3^{\sigma}$$

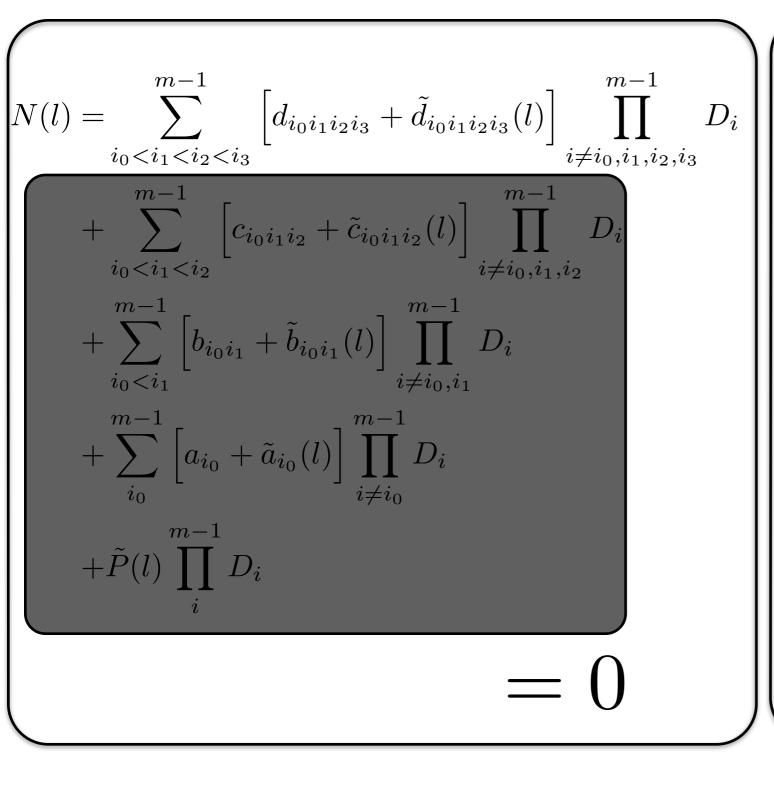
(remember that p_i is the sum of the momentum that has entered the loop so far, so we always have $p_0 = 0$)

• The integral is zero

$$\int d^d l \frac{\tilde{d}_{i_0 i_1 i_2 i_3}(l)}{D_0 D_1 D_2 D_3} = \tilde{d}_{i_0 i_1 i_2 i_3} \int d^d l \frac{\epsilon^{\mu\nu\rho\sigma} l^\mu p_1^\nu p_2^\rho p_3^\sigma}{D_0 D_1 D_2 D_3} = 0$$







To solve the OPP reduction, choosing special values for the loop momenta helps a lot

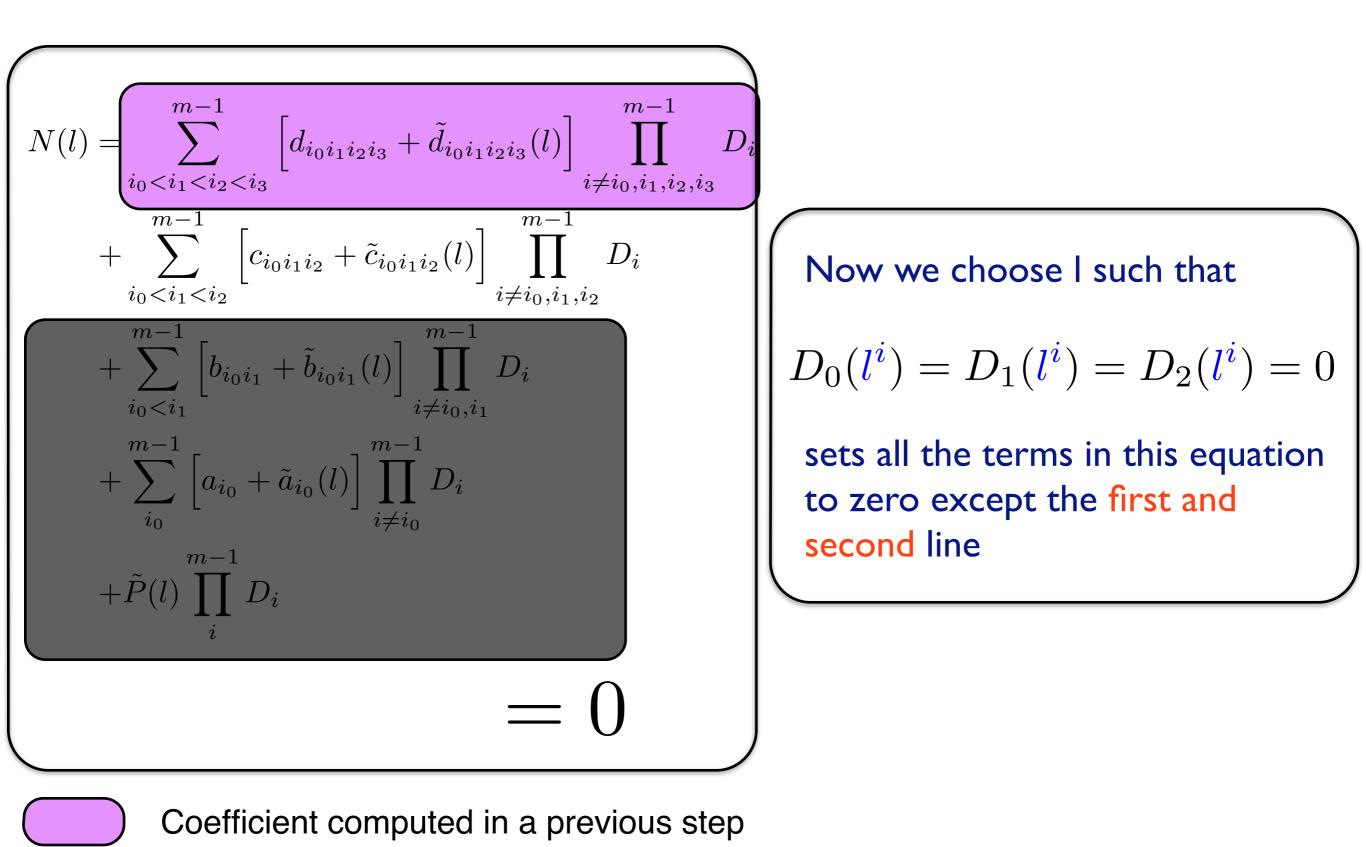
For example, choosing *l* such that $D_0(l^{\pm}) = D_1(l^{\pm}) =$ $= D_2(l^{\pm}) = D_3(l^{\pm}) = 0$

sets all the terms in this equation to zero except the first line

There are two (complex) solutions to this equation due to the quadratic nature of the propagators



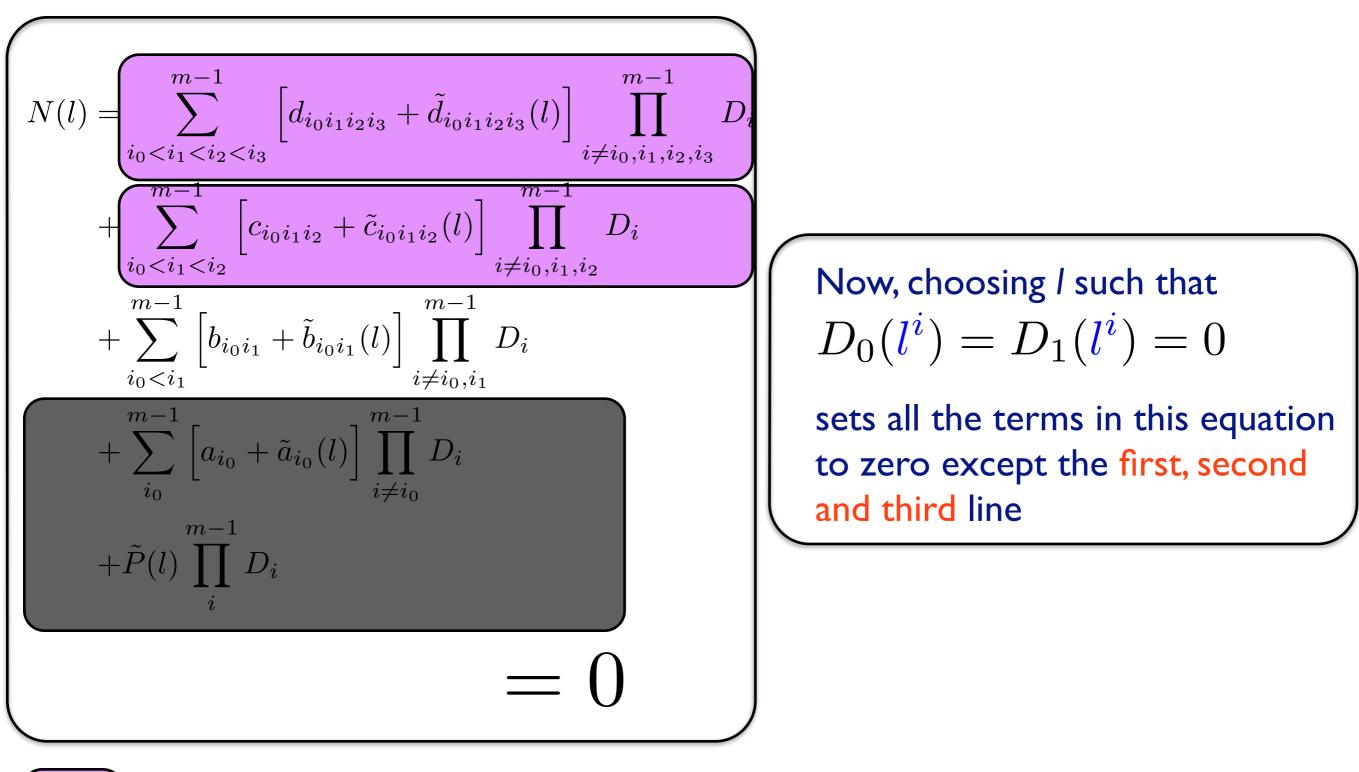




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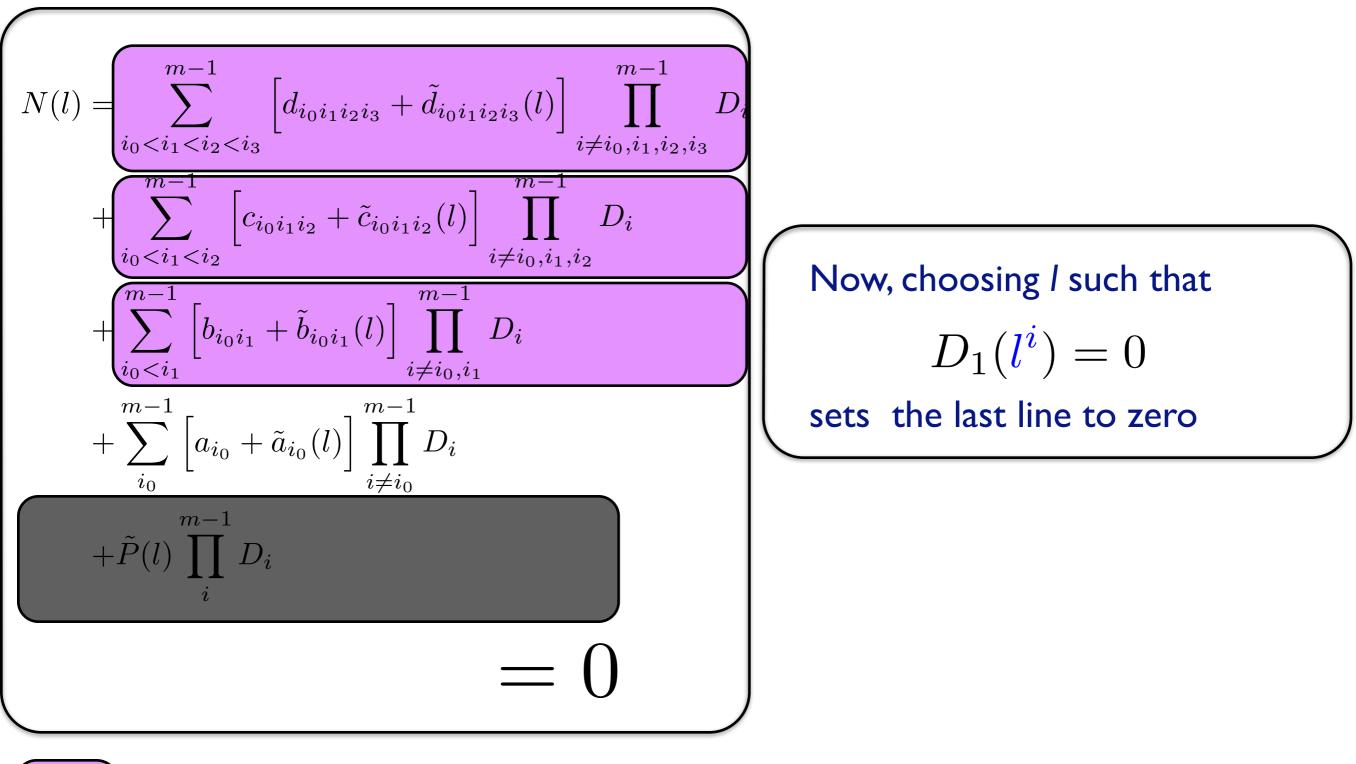




Coefficient computed in a previous step



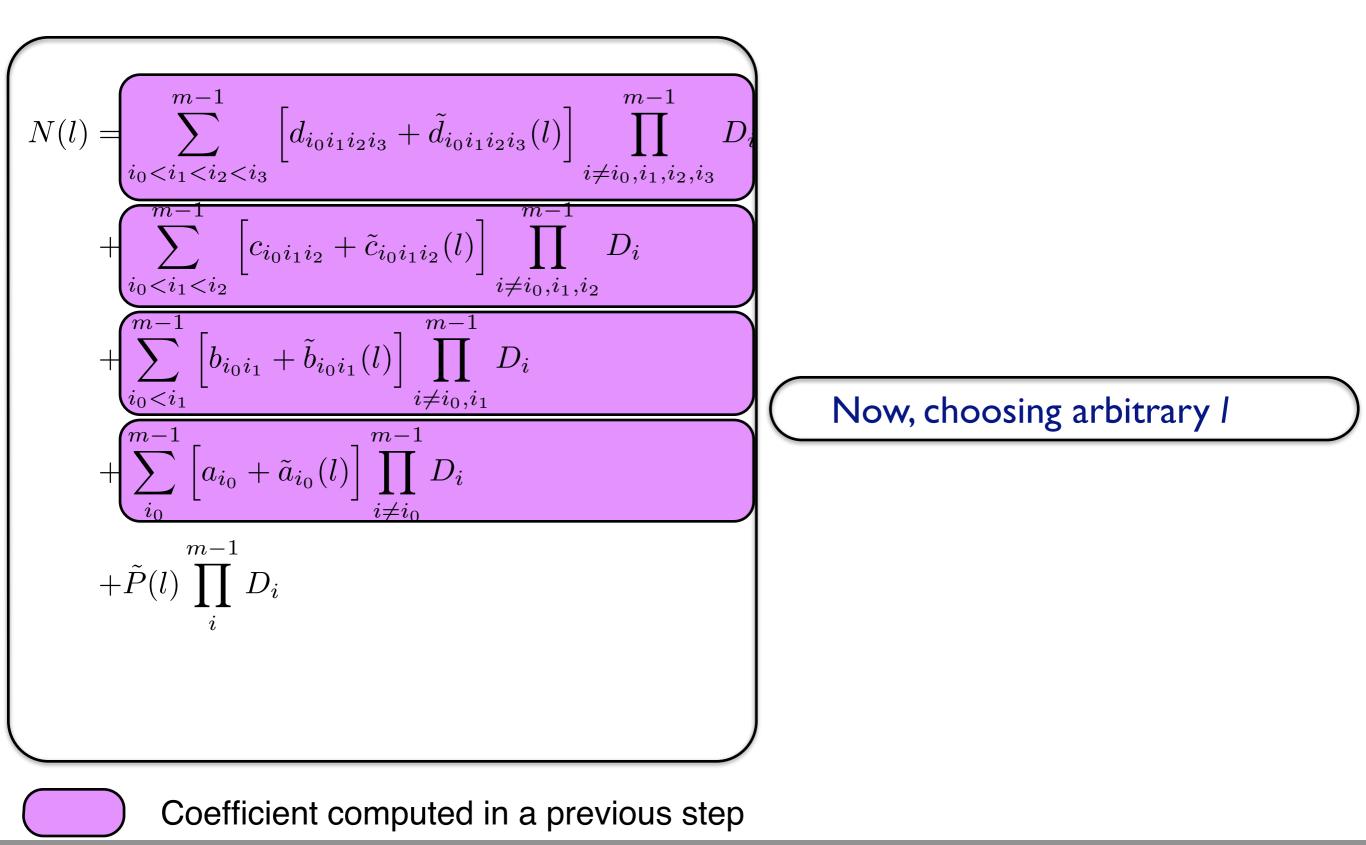




Coefficient computed in a previous step



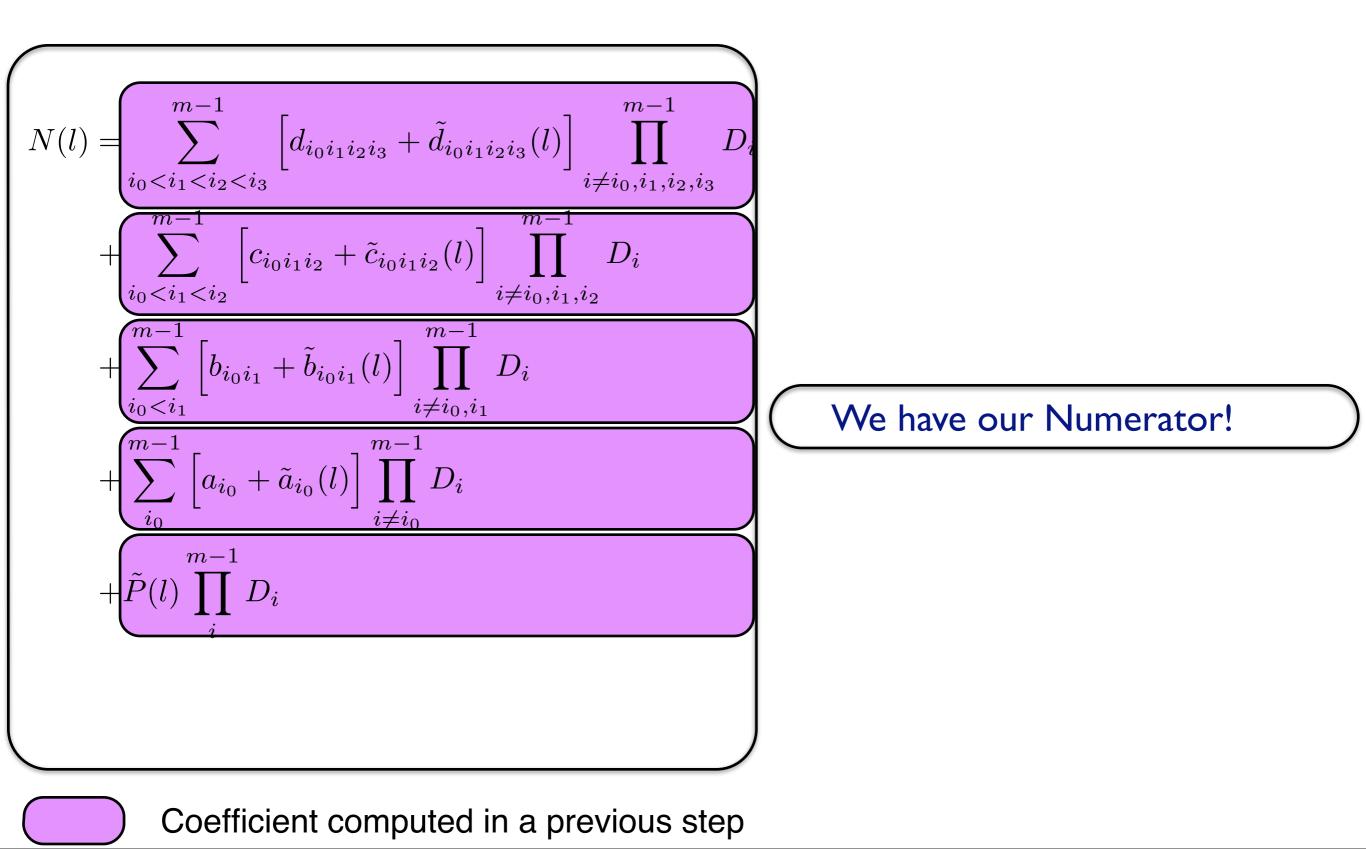




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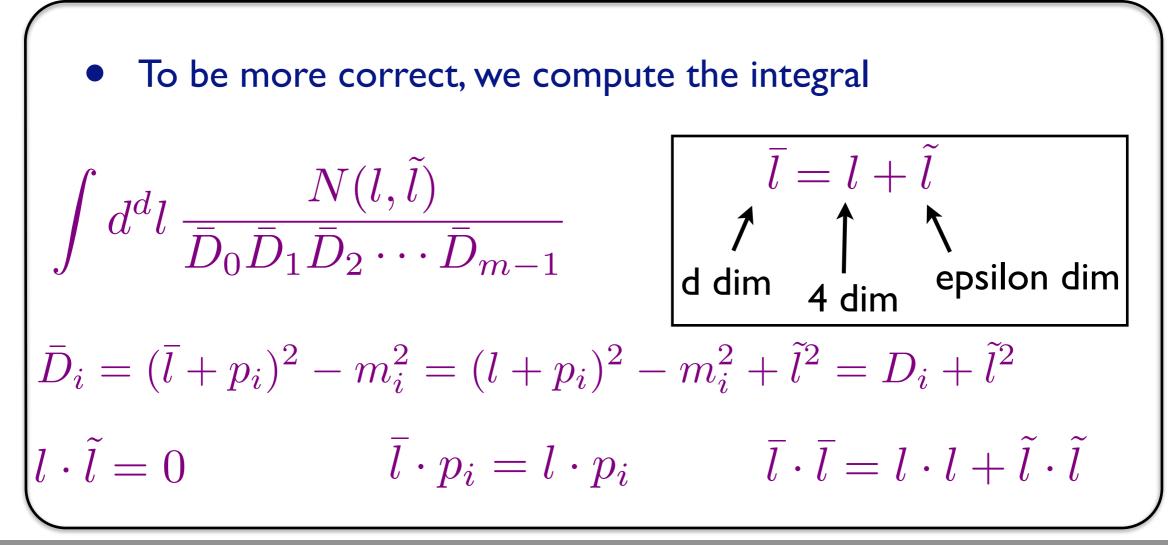
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d dimensions



- In the previous consideration I was very sloppy in considering if we are working in 4 or d dimensions
- In general, external momenta and polarization vectors are in 4 dimensions; only the loop momentum is in d dimensions







The decomposition in terms of scalar integrals has to be done in d dimensions

• This is why the rational part R is needed

$$\begin{split} & \sum_{\substack{0 \leq i_0 < i_1 < i_2 < i_3}}^{m-1} d(i_0 i_1 i_2 i_3) \int d^d \bar{\ell} \frac{1}{\bar{D}_{i_0} \bar{D}_{i_1} \bar{D}_{i_2} \bar{D}_{i_3}} \\ & + \sum_{\substack{0 \leq i_0 < i_1 < i_2}}^{m-1} c(i_0 i_1 i_2) \int d^d \bar{\ell} \frac{1}{\bar{D}_{i_0} \bar{D}_{i_1} \bar{D}_{i_2}} \\ & + \sum_{\substack{0 \leq i_0 < i_1}}^{m-1} b(i_0 i_1) \int d^d \bar{\ell} \frac{1}{\bar{D}_{i_0} \bar{D}_{i_1}} \\ & + \sum_{\substack{i_0 = 0}}^{m-1} a(i_0) \int d^d \bar{\ell} \frac{1}{\bar{D}_{i_0}} \\ & + R \,. \end{split}$$





 In the OPP method, they are split into two contributions, generally called

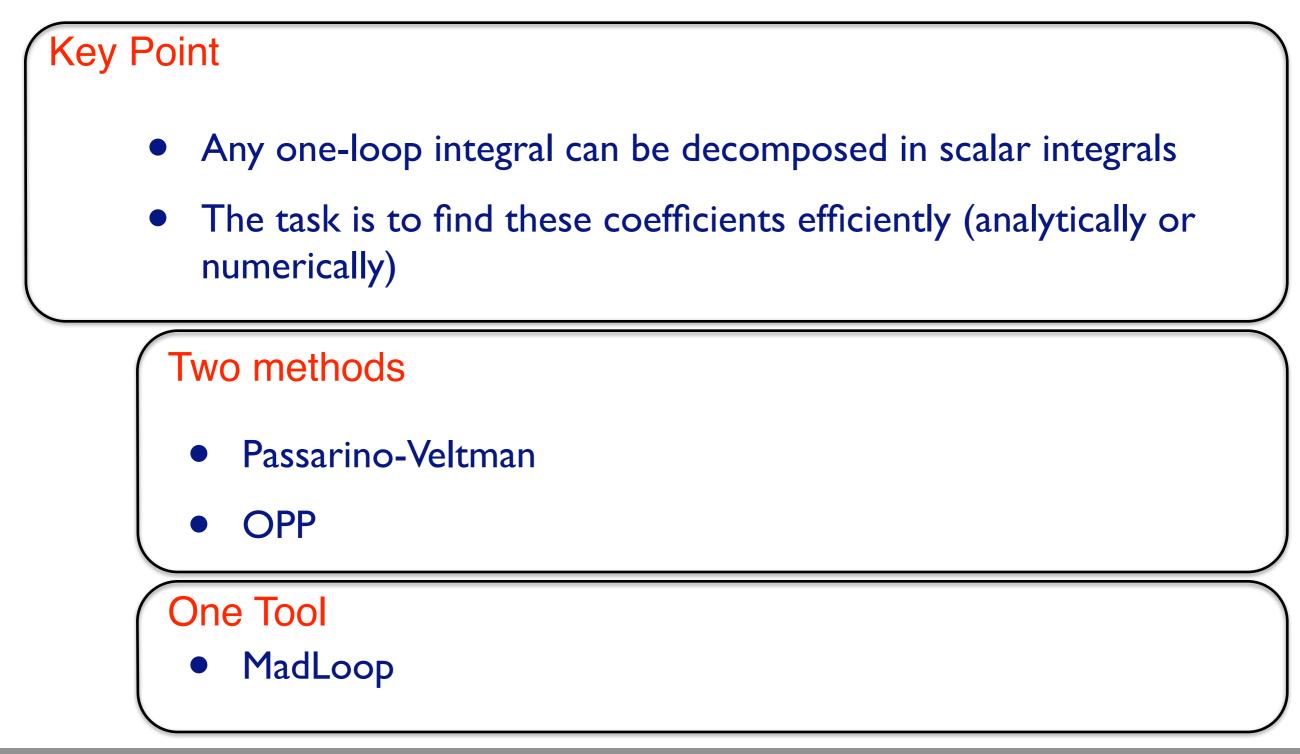
$$R = R_1 + R_2$$

- Both have their origin in the UV part of the model, but only R₁ can be directly computed in the OPP reduction and is given by the CutTools program
 - RI: originates from the propagator (calculate by CutTools)
 - R2: originates from the numerator (need in the model)





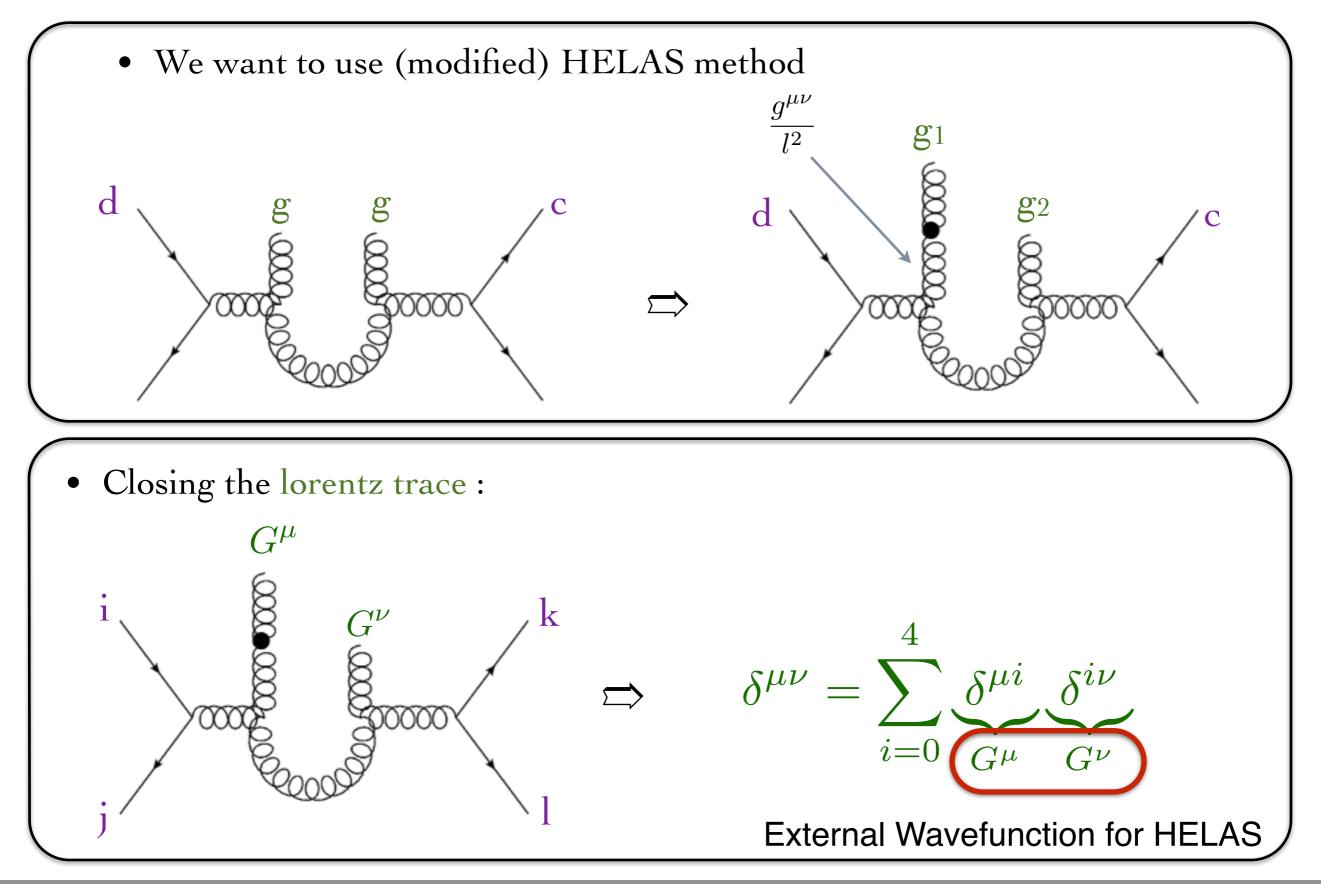






Numerator







MadLoop



Н

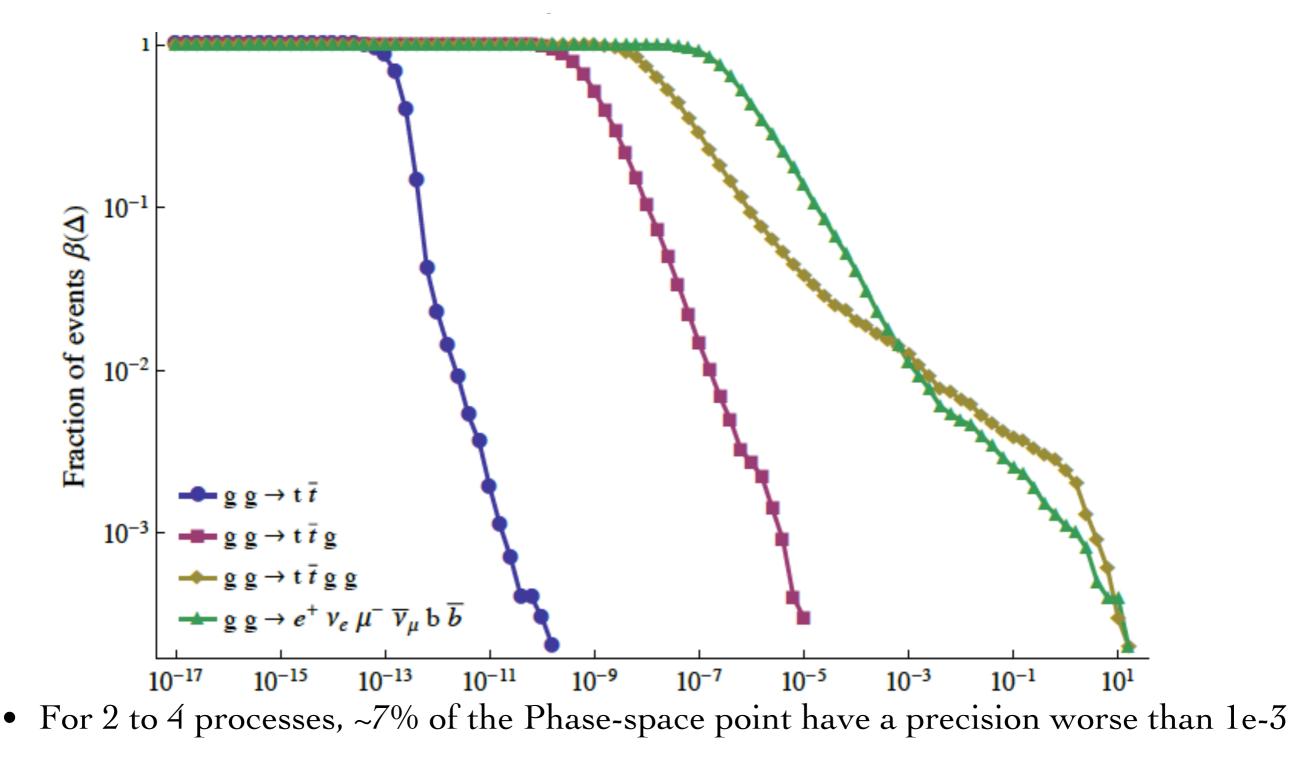
- Other modifications :
 - ► Allow for the loop momentum to be complex
 - ► Remove the denominator of the loop propagators
 - ➡ Close the color trace
 - Ok, now this gives you $\mathcal{N}(l^{\mu})$, the integrand numerator to be fed to CT!
 - But this is **SLOW**!!
 - We have to compute this numerator ~ 50 times for each phase-space point!
 - Idea instead of computing the numerator compute the polynomial form

$$\mathcal{N}(l^{\mu}) = \sum_{r=0}^{r_{max}} C^{(r)}_{\mu_0\mu_1\cdots\mu_r} l^{\mu_0} l^{\mu_1}\cdots l^{\mu_r}$$
[S. Pozzorini & al. hep-ph/1111.5206]



Numerical Stability





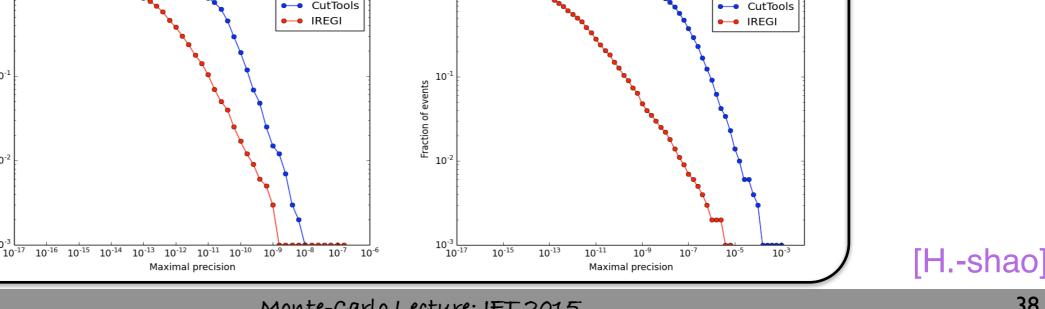
➡ Previous solution pass to quadruple precision (extremelly slow)



 10^{-2}

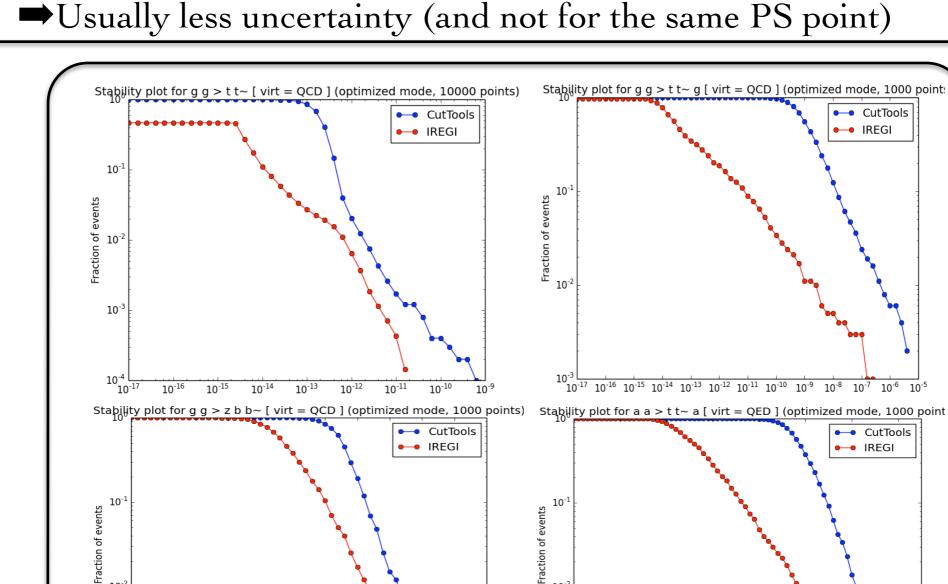
10

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CutTools

IREGI



Usually less uncertainty (and not for the same PS point)

IREGI

Slower than previous method but faster than quadruple precision

New Solution use IREGI: a TIR program

Maximal precision







To Remember



- The main trick is to decompose in scalar integral
- OPP: works at the integrand level
- TIR: works at the integral level
- Loop evaluation is very slow
- Loop evaluation can be "unstable"





Various package in MG5_aMC@NLO







exemple: HEFT

- Model Description
- Width Computation
- Decay Chain
- Interference

Will not be cover

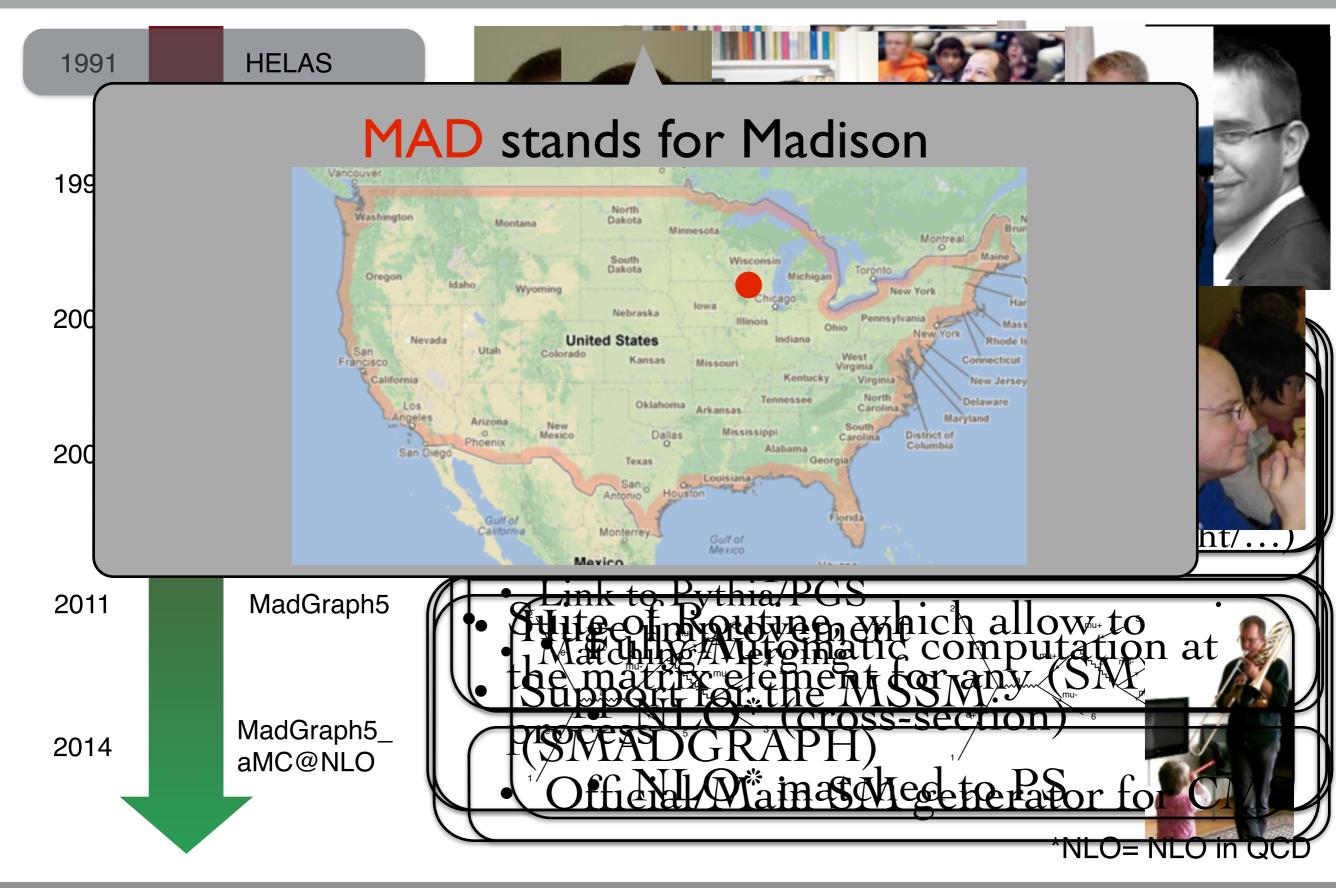
- Re-Weighting method
- Scale Variation
- TauDecay
- MadDM

- MadWeight
- Standalone
- external matrix element provider (Pythia8 and Matchbox)



MadGraph





- Effective Field Theory



$$\mathcal{L} = \mathcal{L}_{SM} + \sum rac{c_i}{\Lambda^2} \mathcal{O}_i$$

Only few Operators for one process and different effects



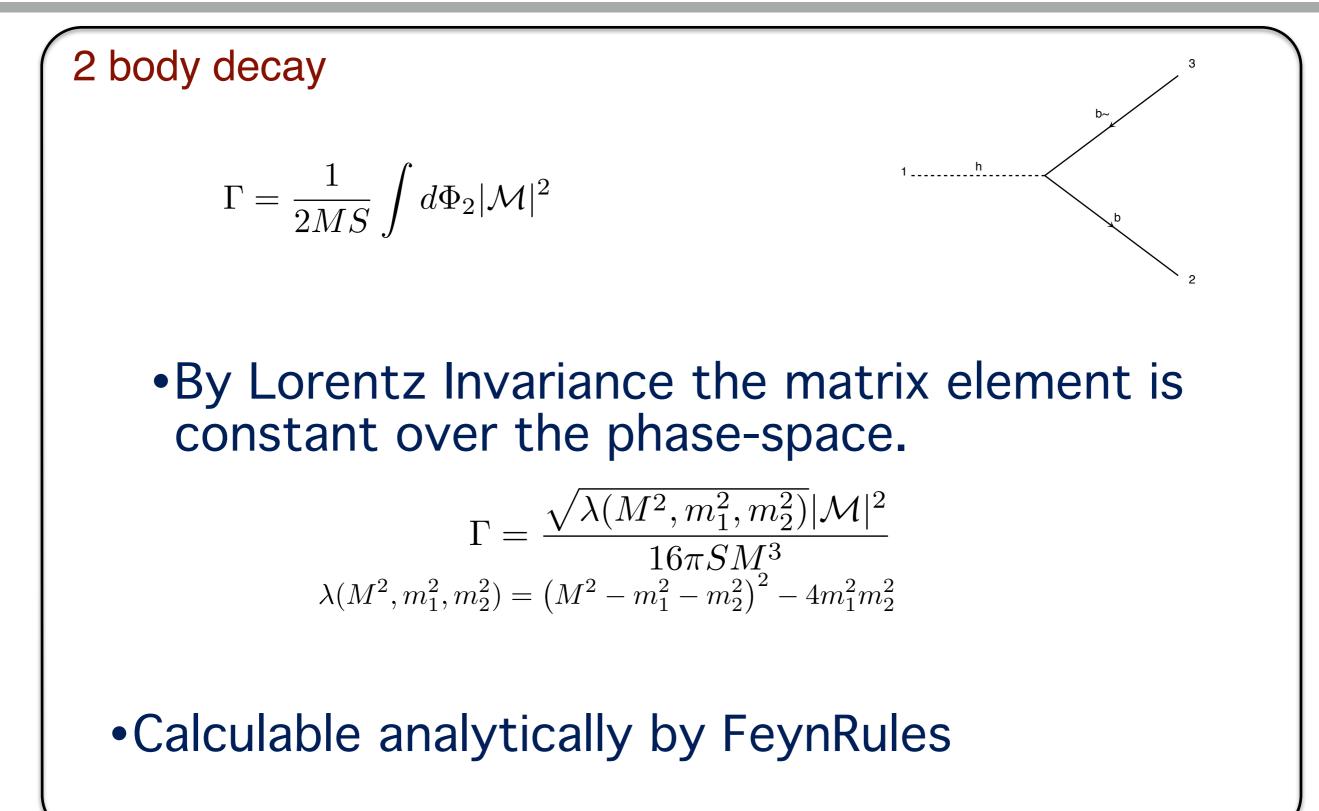
 $\mathcal{O}_{WWW} = \operatorname{Tr}[W_{\mu\nu}W^{\nu\rho}W^{\mu}_{\rho}]$ $\mathcal{O}_{W} = (D_{\mu}\Phi)^{\dagger}W^{\mu\nu}(D_{\nu}\Phi)$ $\mathcal{O}_{B} = (D_{\mu}\Phi)^{\dagger}B^{\mu\nu}(D_{\nu}\Phi)$

$$\mathcal{O}_{\tilde{W}WW} = \operatorname{Tr}[\tilde{W}_{\mu\nu}W^{\nu\rho}W^{\mu}_{\rho}]$$
$$\mathcal{O}_{\tilde{W}} = (D_{\mu}\Phi)^{\dagger}\tilde{W}^{\mu\nu}(D_{\nu}\Phi)$$



2-body decay

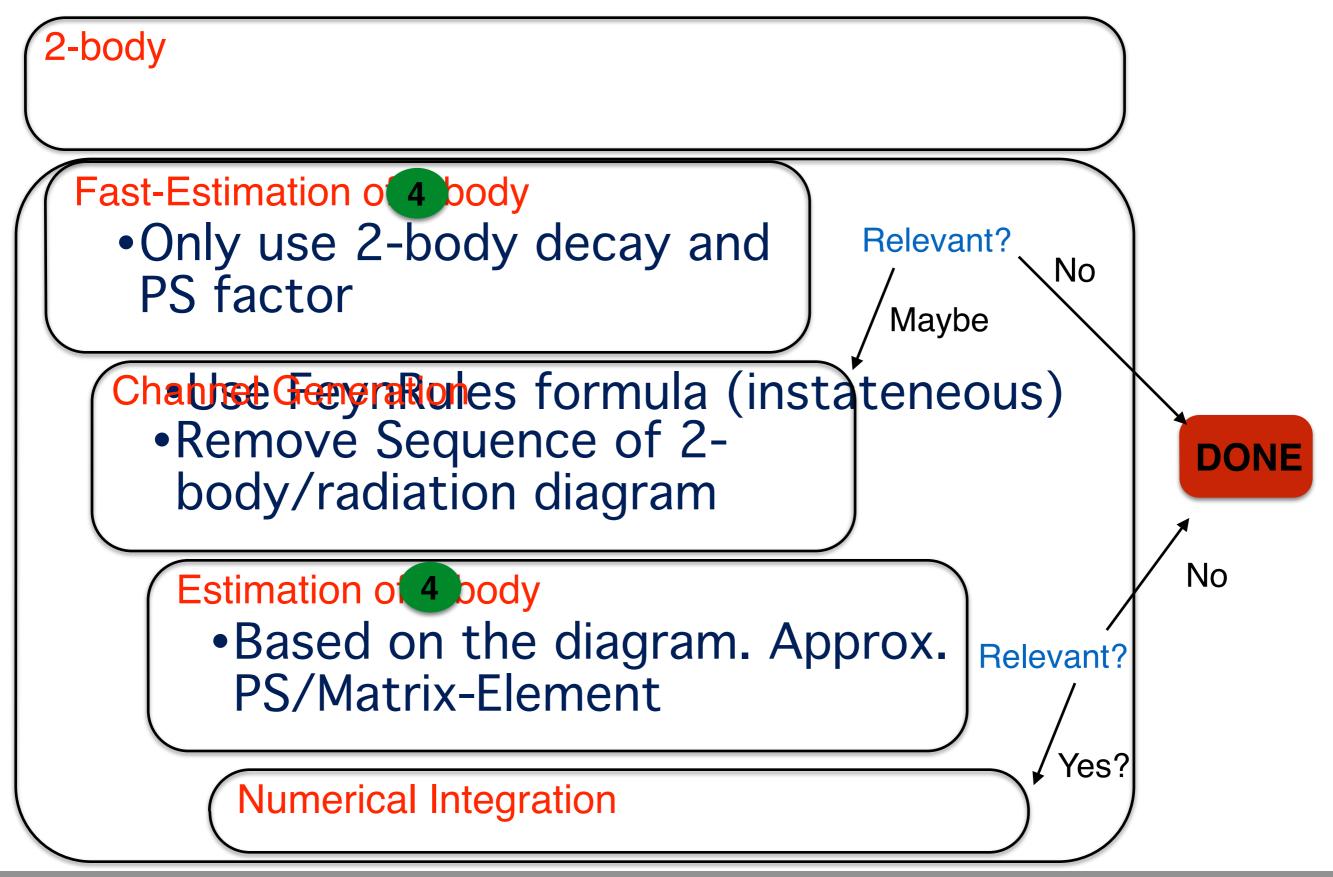






MadWidth hep-ph/1402.1178

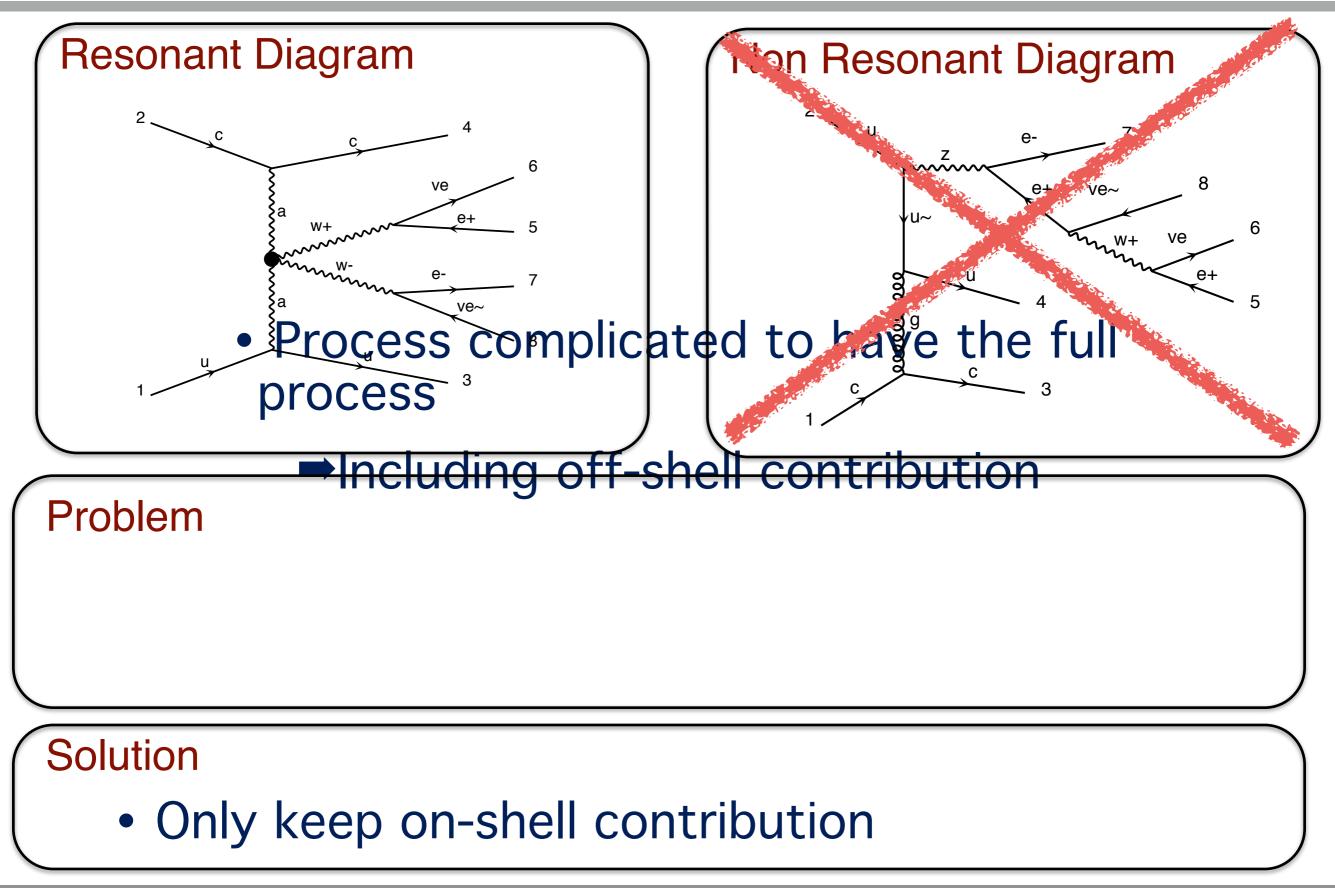










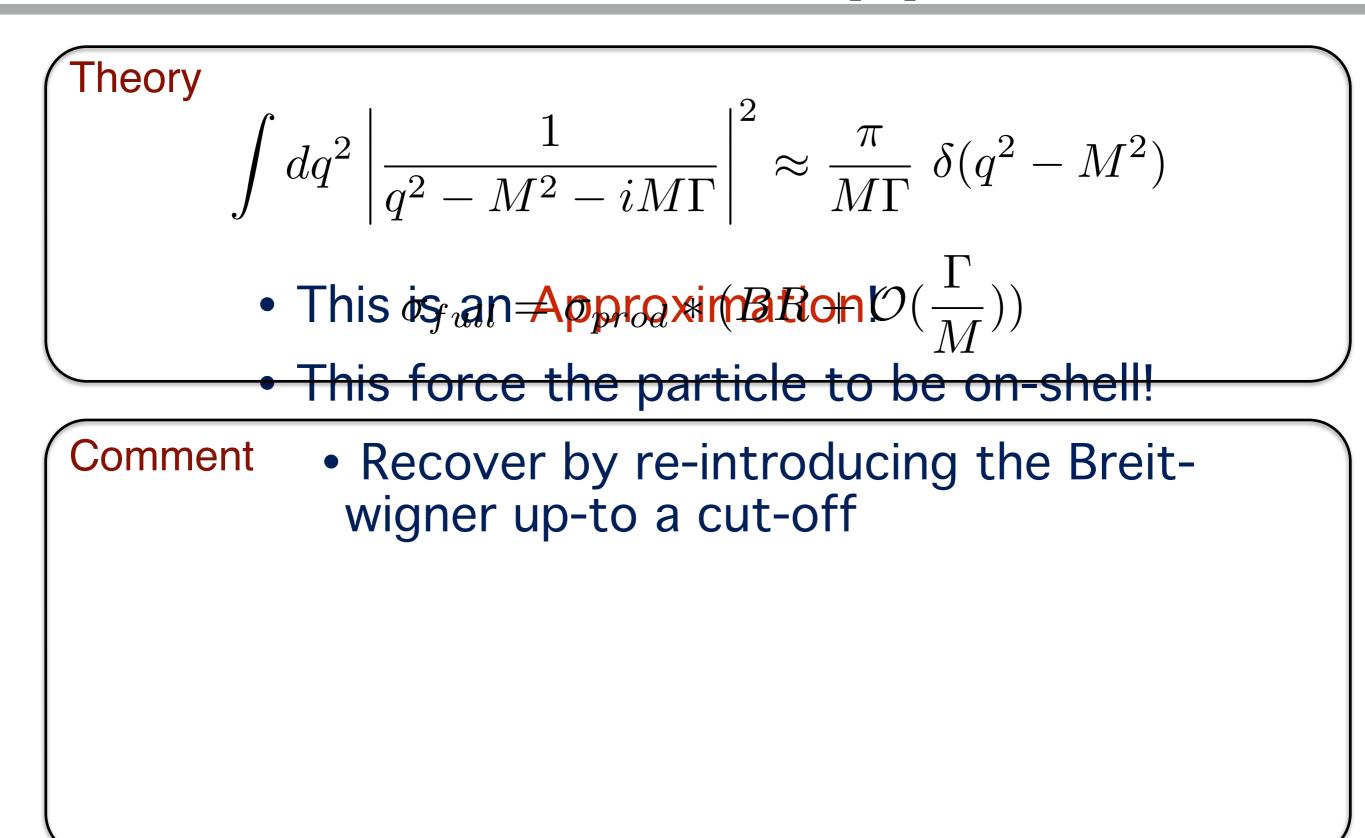


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Narrow-Width Approx.









Decay chains

•
$$p p > t t \sim w+, (t > w+ b, w+ > |+ v|), (t \sim > w- b \sim, w- > j j), (w+ > |+ v|)$$

- Separately generate core process and each decay
 Decays generated with the decaying particle as resulting wavefunction
- Iteratively combine decays and core processes
- Difficulty: Multiple diagrams in decays



Decay chains

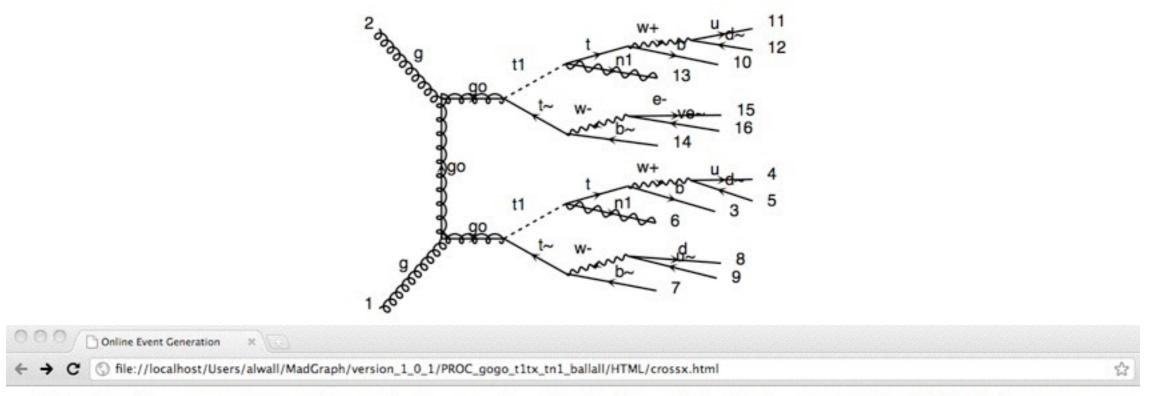


- •Decay chains retain full matrix element for the diagrams compatible with the decay
- Full spin correlations (within and between decays)
- •Full width effects
- •However, no interference with non-resonant diagrams
 - Description only valid close to pole mass
 - ➡ Cutoff at Im ± n□ where n is set in run_card.



Decay chains





Results for g g > go go , (go > t1 t~, t~> b~ all all / h+ , (t1 > t n1 , t > b all all / h+)) in the mssm

Available Results

Links	Events	Tag	Run	Collider	Cross section (pb)	Events
results banner	Parton-level LHE	fermi	test	p p 7000 x 7000 GeV	.33857E-03	10000

Main Page

Thanks to developments in MadEvent, also (very) long decay chains possible to simulate directly in MadGraph!

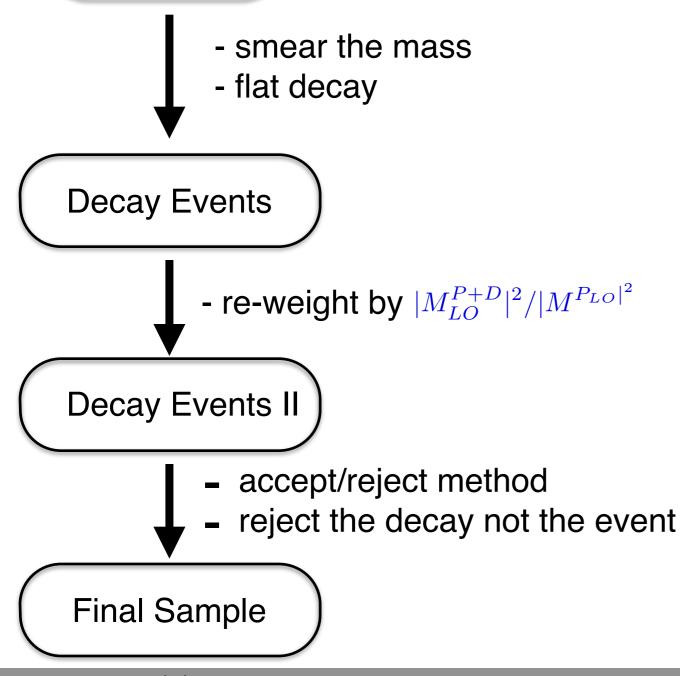
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MadSpin

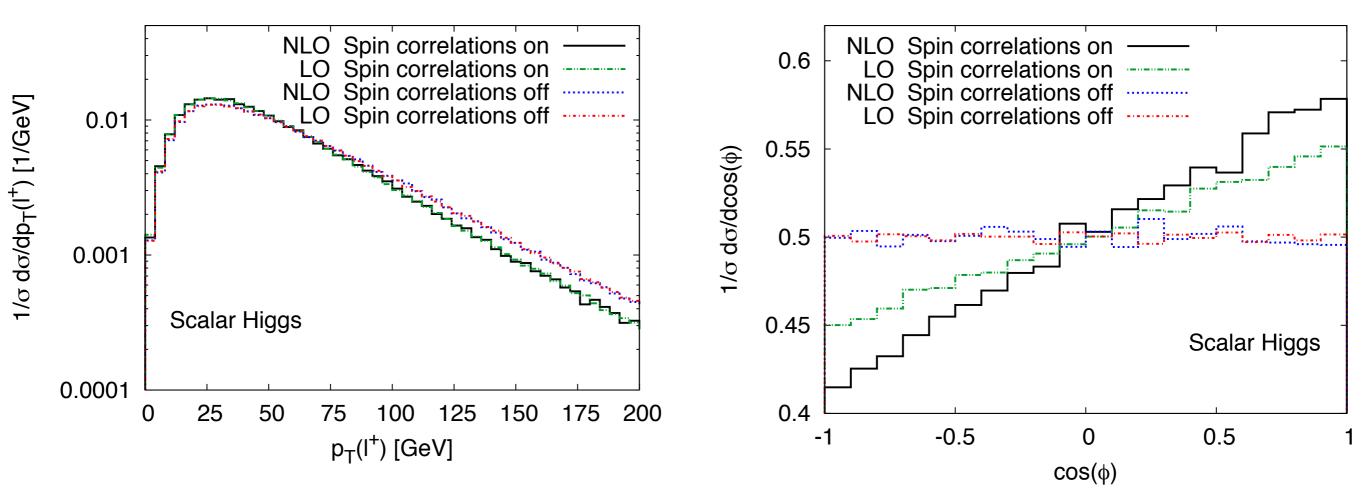


[Frixione, Leanen, Motylinski, Webber (2007)] (Frixione, Leanen, Motylinski, Webber (2007)] (One Event) (One Event) (Artoisenet, OM et al. 1212.3460] (



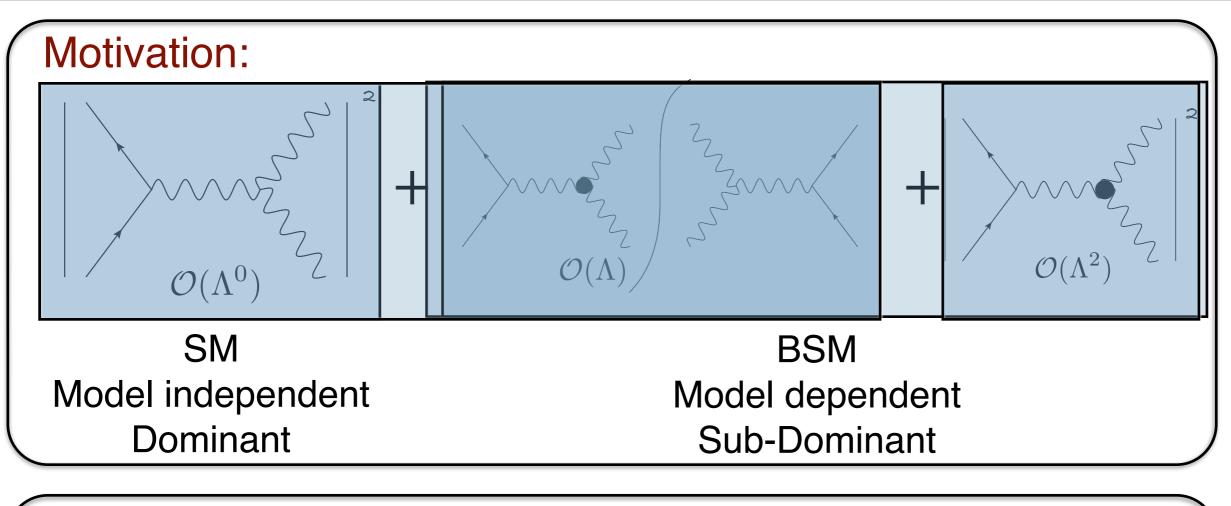
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Idea:

- Compute them separately
- Have a new syntax for such selection (NP^2=)

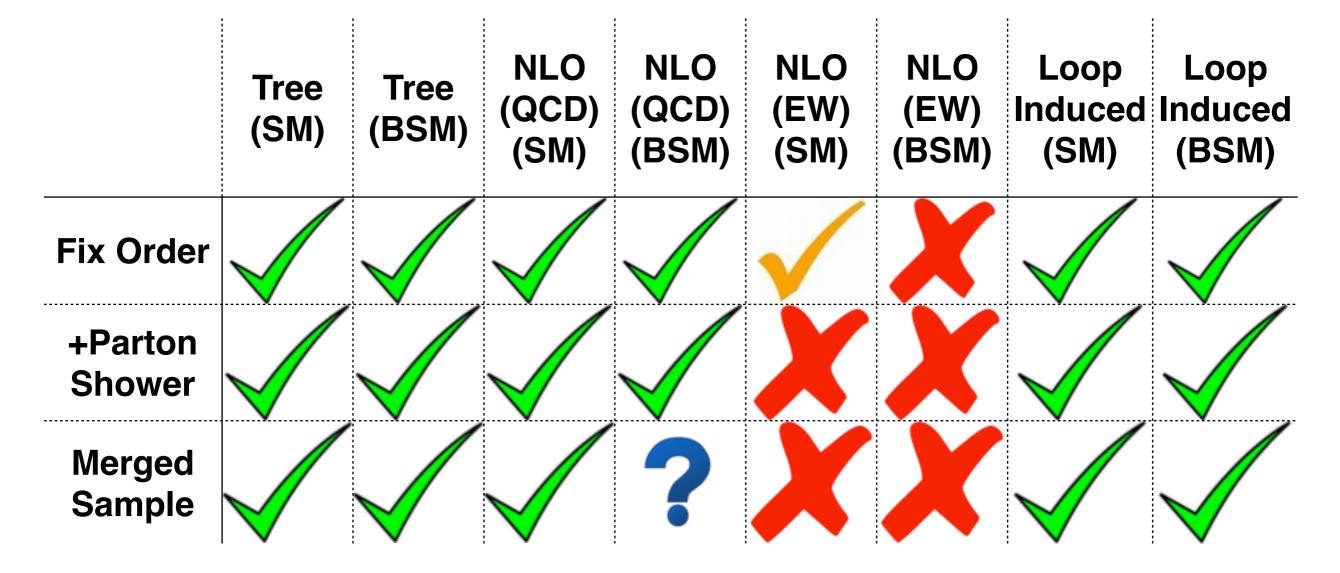
Status:

•Not compatible with decay chains



Type of generation





Loop Induced:

- 2 to 2 processes: OK on a laptop
- 2 to 3 processes: OK on a small size cluster
- 2 to 4 processes: Specific case

MadGraph Functionality





Leading Order Option

- Support of BSM
- Computation of the Width
- Narrow width Approximation
 - Decay Chain
 - MadSpin
- Systematics
- NLO
 - SM with merging