

Automated computation of scattering amplitudes

Tiziano Peraro

Max-Planck-Institut für Physik
Föhringer Ring 6, D-80805 München, Germany

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Outline

- 1 Introduction and motivation
- 2 Amplitudes, trees and loops
- 3 One-loop amplitudes
- 4 Higher loops
- 5 Summary & Conclusions

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Motivation

- The **Large Hadron Collider (LHC)** proved to be capable of
 - **validating SM** in unexplored regions of the phase space
 - making **new discoveries**, e.g. a Higgs(-like?) boson
- LHC high-energy events are characterized by
 - large **SM background** (could hide **new physics**)
 - **multi-particle** final states

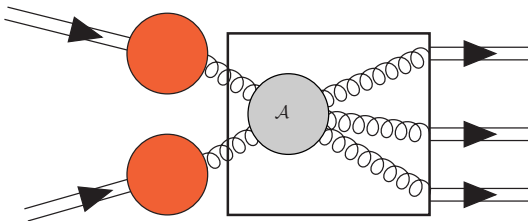
How to make the best use of LHC data?

We need **theoretical predictions** with

- high accuracy
- multi-particle interactions

Scattering Amplitudes

- Hadron collider interactions



- **Scattering amplitudes** represent
 - the **process-dependent part** of a physical event
 - the main point of contact btw. **theoretical models** and **phenomenology**
- They can be computed in **perturbation theory**

$$\mathcal{A} \sim \mathcal{A}_{\text{LO}} + \alpha \mathcal{A}_{\text{NLO}} + \alpha^2 \mathcal{A}_{\text{NNLO}} + \dots$$

Why automation?

- avoid human mistakes
- the complexity of the computation is often too high (for a human)
 - ... but it can be managed by a computer
- every new process requires the computation of new amplitudes
 - ⇒ implement a universal algorithm once
and let the computer do the job every time
- testing new models becomes easier
- experimenting with new/different computational techniques is easier within automated frameworks
- ...

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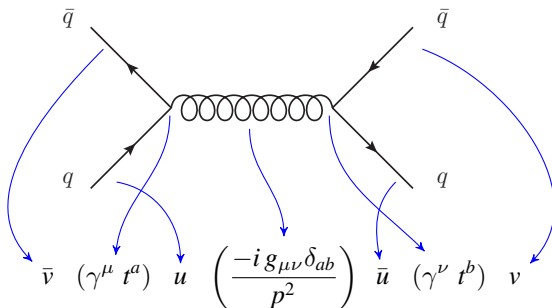
5 Summary & Conclusions

How to compute a scattering amplitude

- A amplitude is a sum of contributions called **Feynman diagrams**

$$\mathcal{A} = \sum_i \mathcal{D}_i$$

- Each **diagram** has a representation as a graph
 - Feynman rules:** each vertex, external and internal line correspond to an algebraic factor



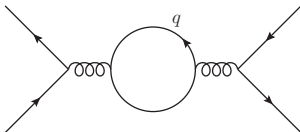
How to compute a scattering amplitude

A strategy:

- draw the **Feynman diagrams**
 - **easy** with the right software (FEYNARTS [T. Hahn], QGRAF [P. Nogueira], ...)
- substitute the **Feynman rules**
 - use a CAS (e.g. FORM [J. Vermaseren])
 - not always easy (because of renorm. schemes, counter-terms, etc. ...)
 - ... but they must be implemented only once per theoretical model
- work out the computation
 - typically **easy** at **leading order** (LO)
 - usually pure algebra
 - **difficult** at **next-to-leading order** (NLO) and beyond
 - involves the computation of **integrals**

Amplitudes and loop integrals

- **Integrals** come from diagrams with **loops**



- the momentum q running in the loop is not fixed by momentum conservation
- from QM: all the values of q give a contribution

$$\mathcal{A} \sim \langle \psi_{out} | \psi_{in} \rangle \sim \sum_q \langle \psi_{out} | q \rangle \langle q | \psi_{in} \rangle, \quad \text{continuum : } \sum_q \rightarrow \int d^4 q$$

- Integration should be done in d dimensions: $d^4 q \rightarrow d^d q$
 - **physical observables** are finite for $d \rightarrow 4$ (amplitudes are **not**!)

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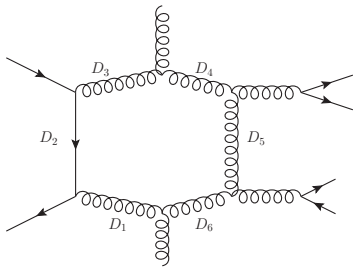
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One-loop amplitudes

- The **integrand** of a generic n -point one-loop integral:
 - is a **rational function** in the components of the **loop momentum** q
 - polynomial numerator** \mathcal{N}

$$\mathcal{M}_n = \int d^d q \, \mathcal{I}_n, \quad \mathcal{I}_n \equiv \frac{\mathcal{N}(q)}{D_1 \dots D_n}$$

- quadratic polynomial denominators** D_i
 - they correspond to Feynman loop propagators

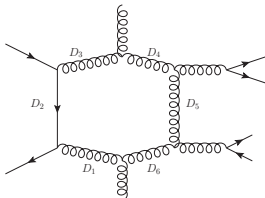


Loop denominators

From the **Feynman rules**:

$$D_i = (q + p_i)^2 - m_i^2$$

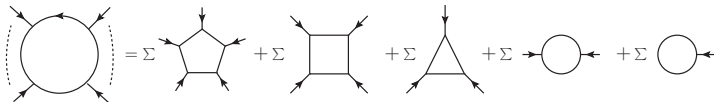
Integrand reduction of one-loop amplitudes



$$\mathcal{M}_n = \int d^d q \frac{\mathcal{N}(q)}{D_1 \dots D_n}$$

- In **multi-particle** scattering amplitudes
 - the number of diagrams (i.e. integrals to be computed) is higher
 - up to thousands or even more
 - the numerators \mathcal{N} are more complex
 - the number of loop denominators is high
 - can be equal to the number of external legs
- **Integrand reduction** of one-loop amplitudes
 - rewrite the **integrand** as a sum of “**simpler**” integrands
 - use an algorithm suited for **automation**

The Integrand reduction of one-loop amplitudes



- **Every** one-loop integrand, can be decomposed as
[Ossola, Papadopoulos, Pittau (2007); Ellis, Giele, Kunszt, Melnikov (2008)]

$$\frac{\mathcal{N}}{D_1 \cdots D_n} = \sum_{j_1 \dots j_5} \frac{\Delta_{j_1 j_2 j_3 j_4 j_5}}{D_{j_1} D_{j_2} D_{j_3} D_{j_4} D_{j_5}} + \sum_{j_1 \dots j_4} \frac{\Delta_{j_1 j_2 j_3 j_4}}{D_{j_1} D_{j_2} D_{j_3} D_{j_4}} + \cdots + \sum_{j_1} \frac{\Delta_{j_1}}{D_{j_1}}$$

- a sum of integrands with **5 or less** loop denominators
 - The **residues** $\Delta_{i_1 \dots i_k}$
 - are **polynomials** in the components of q
 - have a **known, universal parametric form**
 - are parametrized by **unknown, process-dependent coefficients**
- ⇒ can be completely determined with a **polynomial fit**

- Choice of 4-dimensional basis for an m -point residue

$$e_1^2 = e_2^2 = 0, \quad e_1 \cdot e_2 = 1, \quad e_3^2 = e_4^2 = \delta_{m4}, \quad e_3 \cdot e_4 = -(1 - \delta_{m4})$$

- Coordinates: $\mathbf{z} = (z_1, z_2, z_3, z_4, z_5) \equiv (x_1, x_2, x_3, x_4, \mu^2)$

$$q_{4\text{-dim}}^\mu = -p_{i_1}^\mu + x_1 e_1^\mu + x_2 e_2^\mu + x_3 e_3^\mu + x_4 e_4^\mu, \quad \bar{q}^2 = q_{4\text{-dim}}^2 - \mu^2$$

- Generic numerator

$$\mathcal{N} = \sum_{j_1, \dots, j_5} \alpha_j z_1^{j_1} z_2^{j_2} z_3^{j_3} z_4^{j_4} z_5^{j_5}, \quad (j_1 \dots j_5) \quad \text{such that} \quad \text{rank}(\mathcal{N}) \leq \# \text{ loop-denom.}$$

- Residues

$$\Delta_{i_1 i_2 i_3 i_4 i_5} = c_0 \mu^2$$

$$\Delta_{i_1 i_2 i_3 i_4} = c_0 + c_1 x_4 + \mu^2 (c_2 + c_3 x_4 + \mu^2 c_4)$$

$$\Delta_{i_1 i_2 i_3} = c_0 + c_1 x_3 + c_2 x_3^2 + c_3 x_3^3 + c_4 x_4 + c_5 x_4^2 + c_6 x_4^3 + \mu^2 (c_7 + c_8 x_3 + c_9 x_4)$$

$$\Delta_{i_1 i_2} = c_0 + c_1 x_2 + c_2 x_3 + c_3 x_4 + c_4 x_2^2 + c_5 x_3^2 + c_6 x_4^2 + c_7 x_2 x_3 + c_9 x_2 x_4 + c_9 \mu^2$$

$$\Delta_{i_1} = c_0 + c_1 x_1 + c_2 x_2 + c_3 x_3 + c_4 x_4$$

- It can be easily **extended** to **higher-rank** numerators

The Integrand reduction of one-loop amplitudes

- After integration
 - some terms vanish and do not contribute to the amplitude
⇒ **spurious** terms
 - non-vanishing terms give **Master Integrals (MIs)**
 - the amplitude is a **linear combination** of **known MIs**
- The **coefficients** of this linear combination
 - can be identified with some of the coefficients which parametrize the polynomial residues

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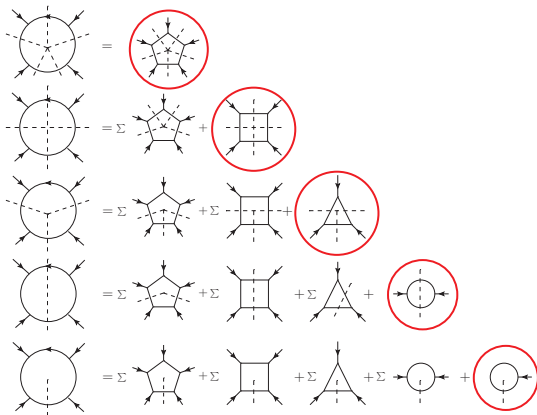
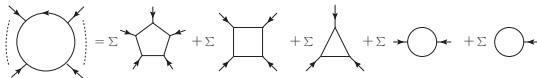
- After integration
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 - can be identified with some of the coefficients which parametrize the polynomial residues
 \Rightarrow **reduction to MIs** \equiv **polynomial fit** of the **residues**
- ★ **any one-loop amplitude** can be computed with a **polynomial fit**

$$\begin{aligned}
 & \text{Sun-like diagram} = c_{4,0} \text{ (square)} + c_{3,0} \text{ (triangle)} + c_{2,0} \text{ (circle)} + c_{1,0} \text{ (empty circle)} \\
 & + c_{4,4} \text{ (square with } d+4 \text{)} + c_{3,7} \text{ (triangle with } d+2 \text{)} + c_{2,9} \text{ (circle with } d+2 \text{)}
 \end{aligned}$$

Fit-on-the-cut at one-loop

[Ossola, Papadopoulos, Pittau (2007)]

Integrand decomposition:



Fit-on-the cut

- fit m -point residues on m -ple cuts
- **Cutting a loop propagator** means

$$\frac{1}{D_i} \rightarrow \delta(D_i)$$

i.e. putting it **on-shell**

Automation of one-loop computation

In several one-loop packages we can distinguish three phases:

- ➊ Generation
 - generate the integrand
 - cast it in a suitable form for reduction
 - write it in a piece of source code (e.g. FORTRAN or C/C++)
- ➋ Compilation
 - compile the code
- ➌ Run-time
 - use a **reduction** library in order to compute the integrals

Automation of one-loop computation in GoSAM

GoSAM is a PYTHON package which:

- generates analytic integrands
 - using QGRAF and FORM
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- can use different reduction algorithms at **run-time**

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 - SAMURAI (d -dim. integrand reduction)
 - faster than GOLEM95 but numerically less stable
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 - NINJA
 - implements a **new** integrand reduction algorithm
 - **fast** (2 to 5 times faster than SAMURAI)
 - **stable** (in worst cases $\mathcal{O}(1/1000)$ unstable points)
 - new default (in GoSAM-2.0)

Integrand reduction via Laurent expansion (NINJA)

P. Mastrolia, E. Mirabella, T.P. (2012)

The integrand reduction via **Laurent expansion**:

- **fits residues** by taking their **asymptotic expansions** on the **cuts**
 - elaborating ideas first proposed by Forde and Badger
- yields **diagonal systems of equations** for the coefficients
- requires the computation of **fewer coefficients**
- subtractions of higher point residues is simplified
 - implemented as **corrections at the coefficient level**

Integrand reduction via Laurent expansion (NINJA)

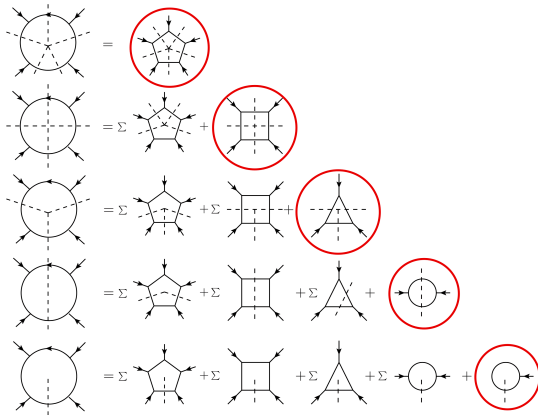
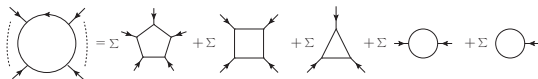
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- ★ Implemented in the semi-numerical C++ library **NINJA** [T.P. (2014)]
 - Laurent expansions via a **simplified polynomial-division algorithm**
 - interfaced with the package GOSAM
 - is a **faster and more stable** integrand-reduction algorithm

Integrand reduction via Laurent expansion (NINJA)

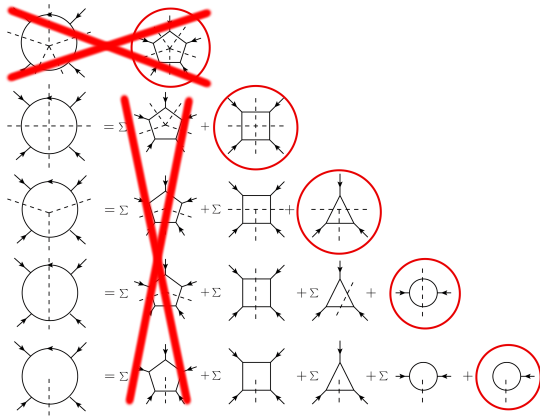
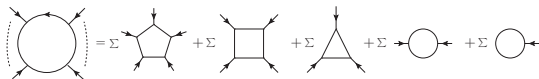
Integrand decomposition:



Laurent-expansion method

Integrand reduction via Laurent expansion (NINJA)

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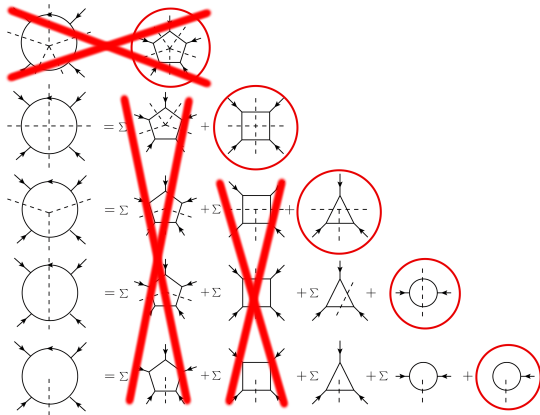
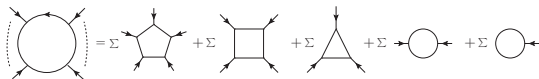


Laurent-expansion method

- pentagons not needed

Integrand reduction via Laurent expansion (NINJA)

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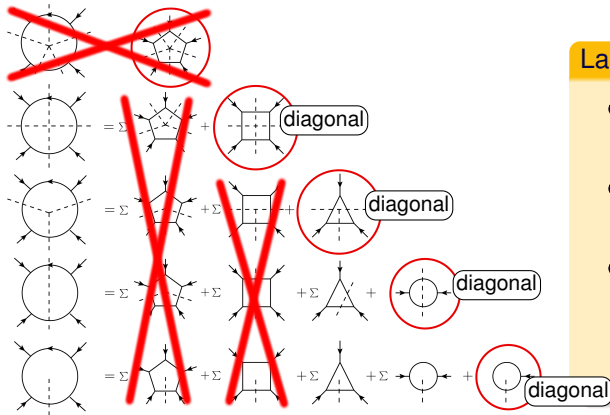
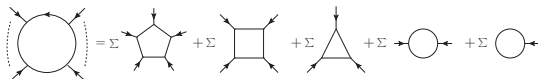


Laurent-expansion method

- pentagons not needed
- boxes never subtracted

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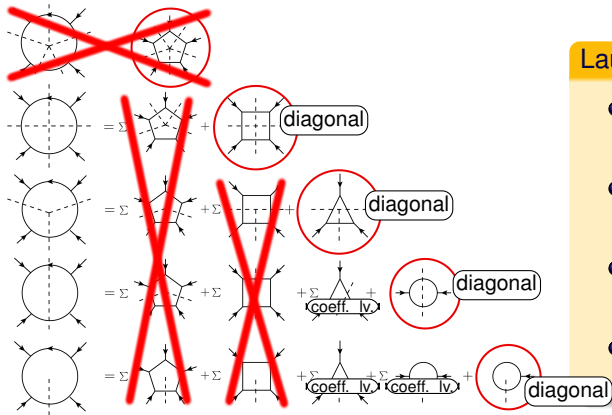
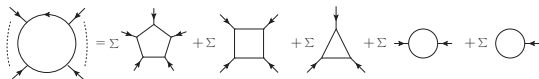


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Benchmarks of GoSAM + NINJA

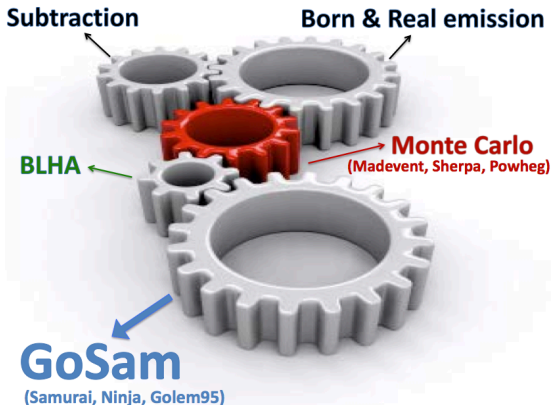
H. van Deurzen, G. Luisoni, P. Mastrolia, E. Mirabella, G. Ossola and T.P. (2013)

Benchmarks: GoSAM + NINJA			
Process		# NLO diagrams	ms/event ^a
$W + 3j$	$d\bar{u} \rightarrow \bar{\nu}_e e^- ggg$	1 411	226
$Z + 3j$	$d\bar{d} \rightarrow e^+ e^- ggg$	2 928	1 911
$t\bar{t}b\bar{b} (m_b \neq 0)$	$d\bar{d} \rightarrow t\bar{t}b\bar{b}$	275	178
	$gg \rightarrow t\bar{t}b\bar{b}$	1 530	5 685
$t\bar{t} + 2j$	$gg \rightarrow t\bar{t}gg$	4 700	13 827
$W b \bar{b} + 1j (m_b \neq 0)$	$u\bar{d} \rightarrow e^+ \nu_e b\bar{b}g$	312	67
$W b \bar{b} + 2j (m_b \neq 0)$	$u\bar{d} \rightarrow e^+ \nu_e b\bar{b}s\bar{s}$	648	181
	$u\bar{d} \rightarrow e^+ \nu_e b\bar{b}d\bar{d}$	1 220	895
	$u\bar{d} \rightarrow e^+ \nu_e b\bar{b}gg$	3 923	5 387
$H + 3j$ in GF	$gg \rightarrow Hggg$	9 325	8 961
$t\bar{t}H + 1j$	$gg \rightarrow t\bar{t}Hg$	1 517	1 505
$H + 3j$ in VBF	$u\bar{u} \rightarrow Hgu\bar{u}$	432	101
$H + 4j$ in VBF	$u\bar{u} \rightarrow Hggu\bar{u}$	1 176	669
$H + 5j$ in VBF	$u\bar{u} \rightarrow Hggu\bar{u}$	15 036	29 200

more processes in arXiv:1312.6678

^aTimings refer to full color- and helicity-summed amplitudes, using an Intel Core i7 CPU @ 3.40GHz, compiled with `ifort`.

From amplitudes to observables with GoSAM



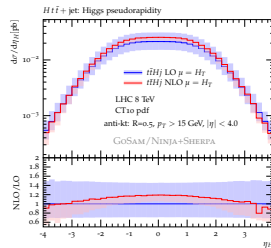
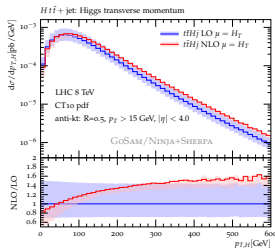
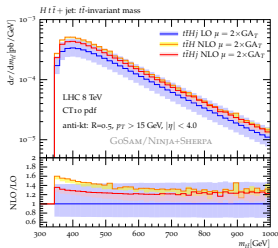
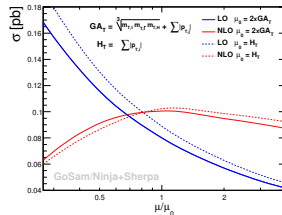
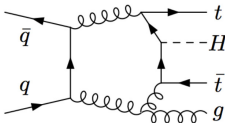
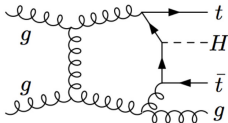
The GoSAM collaboration:

G. Cullen, H. van Deurzen, N. Greiner, G. Heinrich, G. Luisoni, P. Mastrolia, E. Mirabella,
G. Ossola, J. Reichel, J. Schlenk, J. F. von Soden-Fraunhofen, T. Reiter, F. Tramontano, T.P.

Application: $pp \rightarrow t\bar{t}H + jet$ with GoSAM + NINJA

H. van Deurzen, G. Luisoni, P. Mastrolia, E. Mirabella, G. Ossola, T.P. (2013)

● Interfaced with the Monte Carlo SHERPA



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Higher-loop techniques

- Integrand reduction can be extended to higher loop
 - P. Mastrolia, G. Ossola (2011)
 - using **multivariate polynomial division techniques**
 - Y. Zhang (2012); S. Badger, H. Frellesvig, Y. Zhang (2012-13);
 - P. Mastrolia, E. Mirabella, G. Ossola, T.P. (2012-13)
- The **Master Integrals** are **not** all known at 2 or higher loops
- Recent developments in the computation of higher-loop integrals:
 - **differential equations**
 - J. Henn (2013); Henn, Smirnov, Melnikov (2013-14);
 - M. Argeri, S. Di Vita, P. Mastrolia, E. Mirabella, J. Schlenk, U. Schubert, L. Tancredi (2014)
 - **numerical techniques** (e.g. SECDEC)
 - S. Borowka, G. Heinrich (2013-14)

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Summary & Conclusions

- Automated **one-loop** calculations via **integrand-reduction**
 - allow to compute **any** one-loop integral in **any** QFT
 - are implemented in several public codes (e.g. CUTTOOLS, SAMURAI, NINJA)
 - are producing phenomenological results for LHC (e.g. with GOSAM, FORMCALC [[T. Hahn et al.](#)])
- At higher loop
 - we have a framework which extends integrand reduction to any perturbative order
 - many interesting results from both **analytic** and **numerical** techniques
 - work is still in progress but things seem to be promising

THANK YOU
FOR YOUR ATTENTION