

Tropical Feynman integration in the physical region

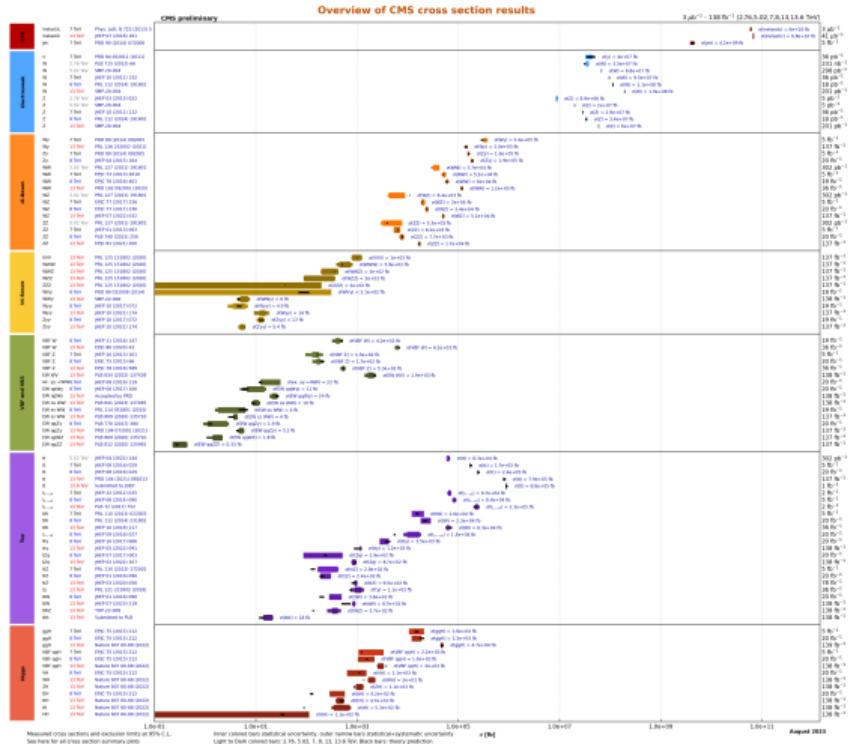
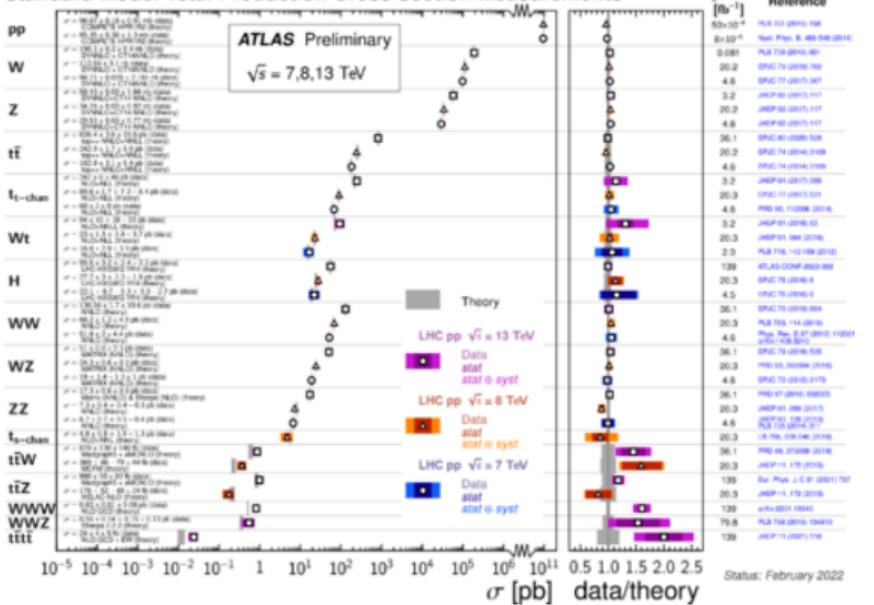
Computer Physics Communications 292 (2023) 108874, with M. Borinsky & H. Munch

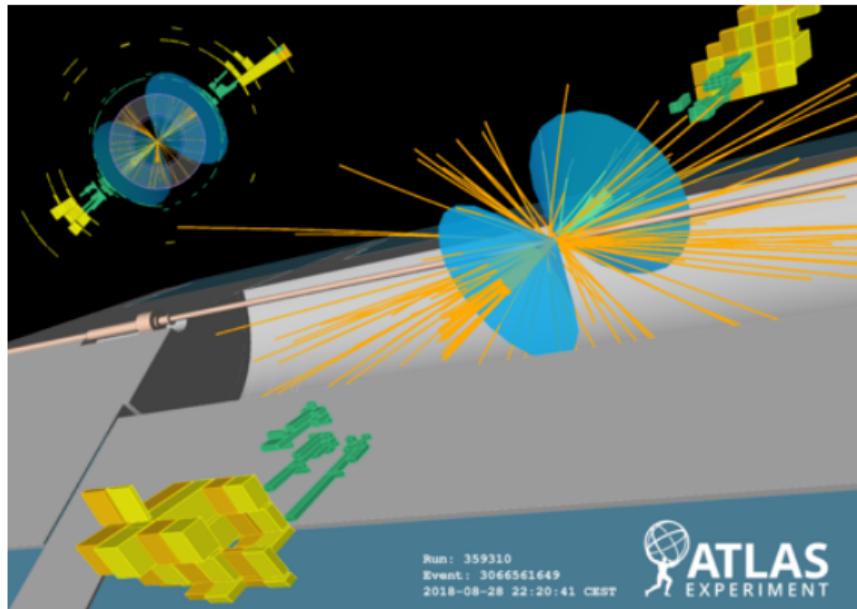
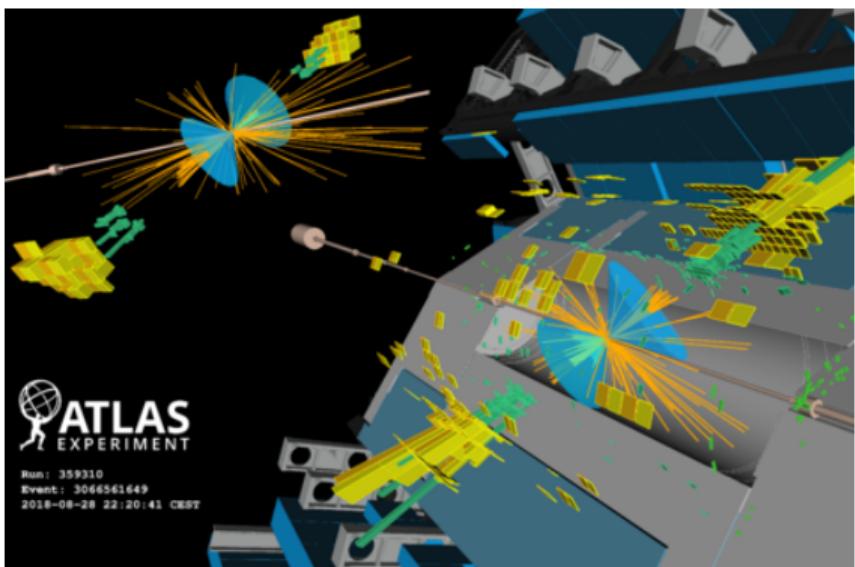
Felix Tellander

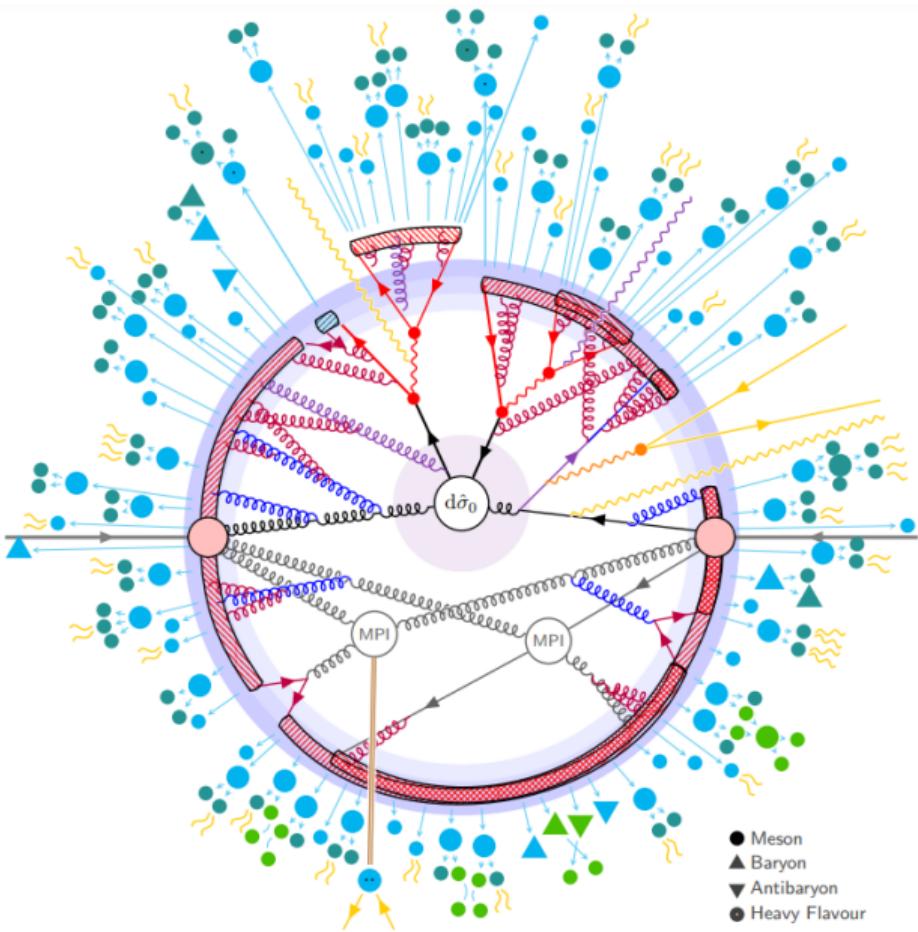
EPS--HEP, 25 Aug. 2023



Standard Model Total Production Cross Section Measurements







- Hard Interaction
- Resonance Decays
- MECs, Matching & Merging
- FSR
- ISR*
- QED
- Weak Showers
- Hard Onium
- Multiparton Interactions
- Beam Remnants*
- Strings
- Ministrings / Clusters
- Colour Reconnections
- String Interactions
- Bose-Einstein & Fermi-Dirac
- Primary Hadrons
- Secondary Hadrons
- Hadronic Reinteractions
- (*: incoming lines are crossed)



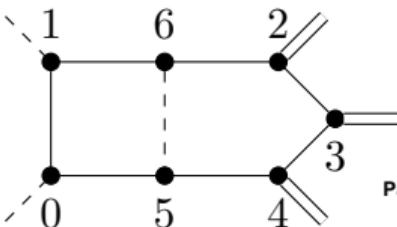
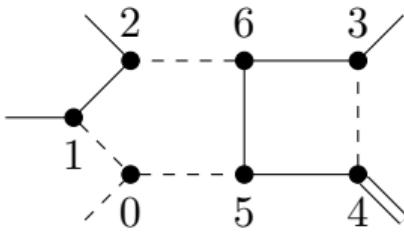
Two-loop box and pentagon integrals

indep. kinem. scales	massive/ off-shell legs	internal masses	process	full σ
$2 \rightarrow 2$				
2	0	0	$\gamma\gamma$	2011
2	0	0	$j j$ (lc)	2017
2	0	0	$\gamma + j$	2017
3	2	1	$t\bar{t}$	2013
3	2	0	VV	2014
4	2	0	VV'	2015
3	1	0	$V + j$	2015
3	1	0	$H + j$ (HTL)	2015
4	2	1	HH	2016
4	1	1	$H + j$	2018
3	0	1	$gg \rightarrow \gamma\gamma$	2019
4	2	1	$gg \rightarrow ZZ$	2020
4	2	1	$gg \rightarrow WW$	2020
5	2	1	$gg \rightarrow ZH$	2021
4	2	1	QCD-EW DY	2022
$2 \rightarrow 3$				
4	0	0	3γ	2019
4	0	0	$\gamma\gamma j$	2021
4	0	0	$3j$	2021
5	1	0	$Wb\bar{b}$	2022

Analytic results:
one off-shell leg

[Abreu et al. 2306.15431]

Today:



Why not NIntegrate?



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$$\int_0^1 \int_0^1 \frac{1}{x+y} dx dy = 2 \log 2 \approx 1.3863$$



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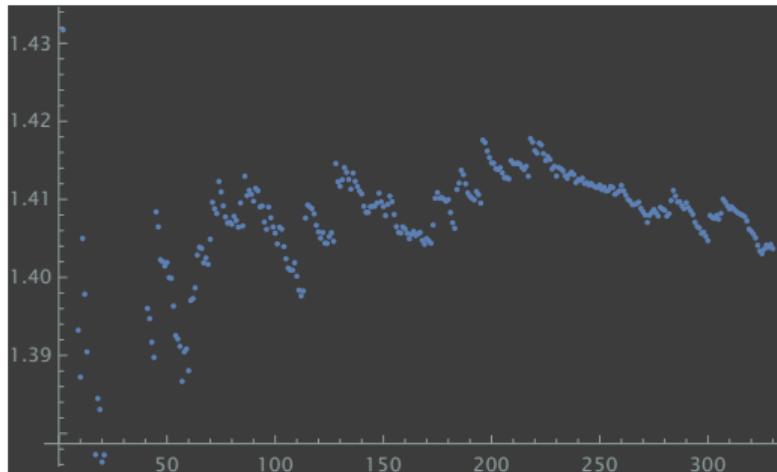
```
In[] NIntegrate[1/(x+y), {x,0,1}, {y,0,1},  
    Method -> {"MonteCarlo", "RandomSeed" -> 19950309}]  
Out[] 0.998259
```



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$$\text{Var} \left(\frac{1}{x+y} dx_{[0,1]} dy_{[0,1]} \right) = \infty$$



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This type of integrable boundary singularities are ubiquitous in Feynman integrals.



Feynman integration software

- > pySecDec [Borowka et al.]
 - > FIESTA [Smirnov]
 - > DiffExp [Hidding]
 - > AMFlow [Liu, Ma]
 - > SeaSyde [Armadillo et al.]
 - > HyperInt [Panzer]
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- > feyntrup [Borinsky, Munch, FT]

Uses [tropical Monte Carlo integration](#) and can be applied to Euclidean as well as Minkowski kinematics.

True power: On your laptop you can evaluate high-loop multi-scale integrals in [minutes](#) to [reasonable error](#).



The Feynman Integral

$$\mathcal{I} = \lim_{\varepsilon \rightarrow 0^+} \Gamma(\omega) \int_{\mathbb{R}_+^E} \prod_{e \in E} \left(\frac{x^{\nu_e} dx_e}{\Gamma(\nu_e) x_e} \right) \mathcal{U}^{-D/2} \frac{\delta(1 - x_1 - \dots - x_E)}{(\mathcal{V} - i\varepsilon \sum_{e \in E} x_e)^\omega}$$

with the superficial degree of divergence

$$\omega := \sum_{e \in E} \nu_e - LD/2$$

where ν_e are propagator powers and $\mathcal{V} = \mathcal{F}/\mathcal{U}$ with homogeneous graph/Symanzik polynomials

$$\mathcal{U} = \sum_{\substack{T \text{ a spanning tree of } G}} \prod_{e \notin T} x_e, \quad \deg(\mathcal{U}) = L$$

$$\mathcal{F} = \mathcal{F}_m + \mathcal{F}_0 = \mathcal{U} \sum_{e \in E} m_e^2 x_e - \sum_{\substack{F \text{ a spanning 2-forest of } G}} p(F)^2 \prod_{e \notin F} x_e, \quad \deg(\mathcal{F}) = L + 1$$



Contour Deformation

Why $i\varepsilon$?

Chooses the causal branch and ensures the convergence.



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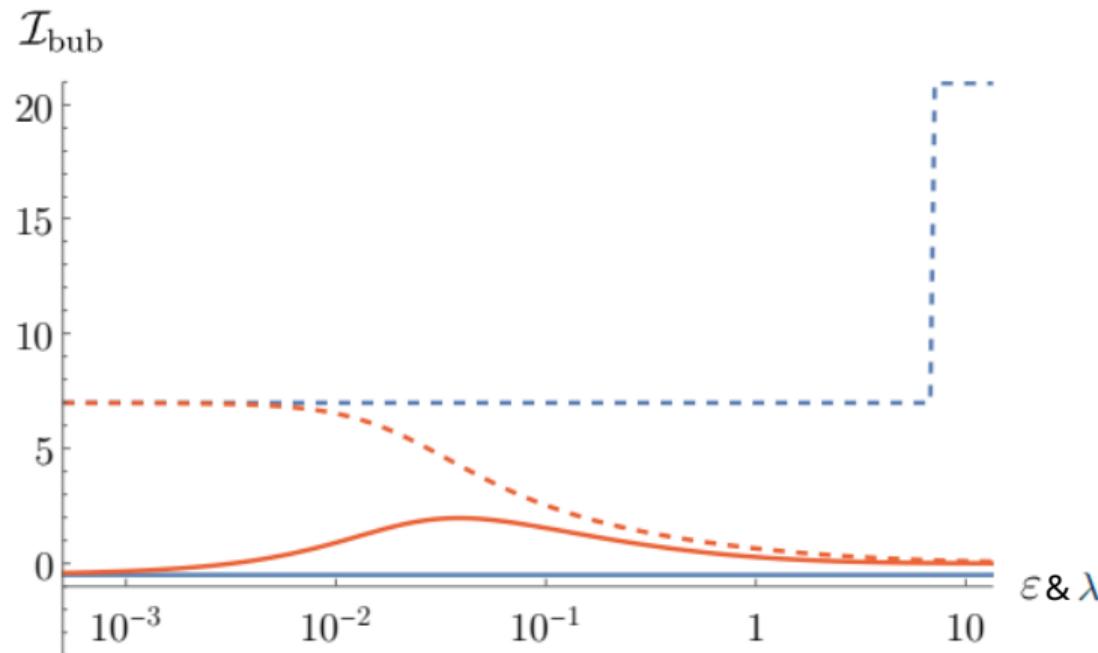
Instead: Change of variables

$$X_e = x_e \exp\left(-i\lambda \frac{\partial \mathcal{V}}{\partial x_e}\right)$$

Picks the same causal branch as $i\varepsilon$ as long as λ is sufficiently small and the Landau equations have no solutions.



Comparison with direct numerics on the Feynman parameterization with $i\varepsilon$ and with deformation:



Tropical Monte Carlo

The **tropical approximation** of a polynomial $p(\mathbf{x}) = \sum_{\alpha \in \text{supp}(p)} c_\alpha \mathbf{x}^\alpha$:

$$p^{\text{tr}}(\mathbf{x}) = \max_{\alpha \in \text{supp}(p)} \{\mathbf{x}^\alpha\}.$$



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Theorem

For a homogeneous polynomial $p \in \mathbb{C}[x_1, \dots, x_n]$ that is completely non-vanishing in \mathbb{P}_+^n there exists constants $C_1, C_2 > 0$ s.t.

$$C_1 \leq \frac{|p(\mathbf{x})|}{p^{\text{tr}}(\mathbf{x})} \leq C_2 \quad \text{for all } \mathbf{x} \in \mathbb{P}_+^n$$



Assumption

You can find bounds on a deformed polynomial with the un-deformed one.

That is, there are λ dependent constants $C_1(\lambda), C_2(\lambda) > 0$ s.t.

$$C_1(\lambda) \leq \left| \left(\frac{\mathcal{U}^{\text{tr}}(\mathbf{x})}{\mathcal{U}(\mathbf{X})} \right)^{D_0/2} \left(\frac{\mathcal{V}^{\text{tr}}(\mathbf{x})}{\mathcal{V}(\mathbf{X})} \right)^{\omega_0} \right| \leq C_2(\lambda) \quad \text{for all } \mathbf{x} \in \mathbb{P}_+^E$$

where the denominators are the deformed polynomials.



Expanding in ϵ :

Assuming that the only potential divergence comes from $\Gamma(\omega)$ we have:

$$\mathcal{I} = \Gamma(\omega_0 + \epsilon L) \sum_{k=0}^{\infty} \frac{\epsilon^k}{k!} \int_{\mathbb{P}_+^E} \left(\prod_{e \in E} \frac{X_e^{\nu_e}}{\Gamma(\nu_e)} \right) \frac{\det \mathcal{J}_\lambda(\mathbf{x})}{\mathcal{U}(\mathbf{X})^{D_0/2} \cdot \mathcal{V}(\mathbf{X})^{\omega_0}} \log^k \left(\frac{\mathcal{U}(\mathbf{X})}{\mathcal{V}(\mathbf{X})^L} \right) \Omega$$

where $\omega_0 = \sum_{e \in E} \nu_e - D_0 L / 2$.



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where $\omega_0 = \sum_{e \in E} \nu_e - D_0 L / 2$.

Writing the integral with these fractions in the integrand:

$$\mathcal{I} = \frac{\Gamma(\omega_0 + \epsilon L)}{\prod_{e \in E} \Gamma(\nu_e)} \sum_{k=0}^{\infty} \frac{\epsilon^k}{k!} \mathcal{I}_k$$

with

$$\mathcal{I}_k = I^{\text{tr}} \int_{\mathbb{P}_+^E} \frac{\left(\prod_{e \in E} (X_e/x_e)^{\nu_e} \right) \det \mathcal{J}_\lambda(\mathbf{x})}{(\mathcal{U}(\mathbf{X})/\mathcal{U}^{\text{tr}}(\mathbf{x}))^{D_0/2} \cdot (\mathcal{V}(\mathbf{X})/\mathcal{V}^{\text{tr}}(\mathbf{x}))^{\omega_0}} \log^k \left(\frac{\mathcal{U}(\mathbf{X})}{\mathcal{V}(\mathbf{X})^L} \right) \mu^{\text{tr}}$$

and

$$\mu^{\text{tr}} = \frac{1}{I^{\text{tr}}} \frac{\prod_{e \in E} x_e^{\nu_e}}{\mathcal{U}^{\text{tr}}(\mathbf{x})^{D_0/2} \mathcal{V}^{\text{tr}}(\mathbf{x})^{\omega_0}} \Omega, \quad \int_{\mathbb{P}_+^E} \mu^{\text{tr}} = 1.$$



Special structure of Feynman integrals

Q: How do we sample from μ^{tr} ?

A: Use that $\mathbf{N}[\mathcal{U}]$ and $\mathbf{N}[\mathcal{F}]$ are generalized permutohedra (some times).

Want details? Let's chat later.



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TLDR: Generalized permutohedra are special polytopal geometries revealing new structures in Feynman integrals. Fits very well in the modern program of polytopal geometries in amplitudes.

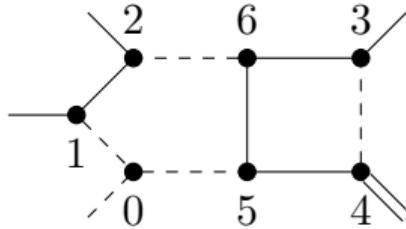


The program feyntrhop

Available at <https://github.com/michibo/feyntrhop>
A C++ program with Python and JSON interface.



Example:



Dashed lines: massless, the solid lines: mass m and double $p_4^2 \neq 0$.

```
edges = [((0,1), 1, '0'), ((1,2), 1, 'mm'), ((2,6), 1, '0'),
         ((6,3), 1, 'mm'), ((3,4), 1, '0'), ((4,5), 1, 'mm'),
         ((5,0), 1, '0'), ((5,6), 1, 'mm')]
```

Phase space point

$$p_0^2 = 0, \quad p_1^2 = p_2^2 = p_3^2 = m^2 = 1/2, \quad s_{01} = 2.2, \quad s_{02} = 2.3, \\ s_{03} = 2.4, \quad s_{12} = 2.5, \quad s_{13} = 2.6, \quad s_{23} = 2.7,$$

where $s_{ij} = (p_i + p_j)^2$. With $\lambda = 0.28$, $N = 10^7$, we obtain:

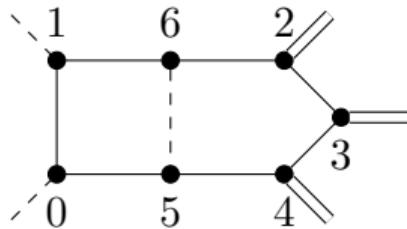
Analytic continuation: activated. Lambda = 0.28.
Started integrating using 4 threads and N = 1e+07 points.
Finished in 19.6719 seconds = 0.00546442 hours.

```
-- eps^0: [0.06474 +/- 0.00066] + i * [-0.08298 +/- 0.00077]
-- eps^1: [0.4066 +/- 0.0036 ] + i * [ 0.3202 +/- 0.0040 ]
-- eps^2: [-0.778 +/- 0.012 ] + i * [ 0.9419 +/- 0.0100 ]
-- eps^3: [-1.328 +/- 0.022 ] + i * [ -1.218 +/- 0.025 ]
-- eps^4: [ 1.395 +/- 0.033 ] + i * [ -1.237 +/- 0.041 ]
```

This is a **two-loop** integral with **seven** mass scales that you can integrate on your laptop in less than **20 seconds**.



Example:



```
edges = [((0,1), 1, 'mm_top'), ((1,6), 1, 'mm_top'),
         ((5,6), 1, '0'), ((6,2), 1, 'mm_top'),
         ((2,3), 1, 'mm_top'), ((3,4), 1, 'mm_top'),
         ((4,5), 1, 'mm_top'), ((5,0), 1, 'mm_top')]
```

Phase space point:

$$m_t^2 = 1.8995, \quad m_H^2 = 1,$$

$$s_{02} = -4.4, \quad s_{03} = -0.5, \quad s_{12} = -0.6, \quad s_{13} = -0.7, \quad s_{23} = 1.8,$$

Setting $\lambda = 0.64$ and $N = 10^8$ we get:

```
Prefactor: gamma(2*eps + 4).  
(Effective) kinematic regime: Minkowski (generic).  
Finished in 8.12 seconds.  
-- eps^0: [-0.0114757 +/- 0.0000082]  
          + i * [0.0035991 +/- 0.0000068]  
-- eps^1: [ 0.003250 +/- 0.000031 ]  
          + i * [-0.035808 +/- 0.000041 ]  
-- eps^2: [ 0.046575 +/- 0.000098 ]  
          + i * [0.016143 +/- 0.000088 ]  
-- eps^3: [ -0.01637 +/- 0.00017 ]  
          + i * [ 0.03969 +/- 0.00016 ]  
-- eps^4: [ -0.02831 +/- 0.00023 ]  
          + i * [-0.00823 +/- 0.00024 ]
```

- > feynntrop 8.12 seconds with relative error $\sim 10^{-3}$
- > pySecDec 3 hours with relative error $\sim 10^{-2}$



Thank you!

Contact

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