

Numerics for elliptic Feynman integrals

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- I: Review: Numerics for multiple polylogarithms**
- II: Example: The two-loop electron self-energy**
- III: Numerics: Single-scale elliptic Feynman integrals**

Motivation

- Precision physics for heavy particles (e.g. Higgs, top, W/Z -bosons).
- If we allow masses, already two-loop two-point functions go beyond multiple polylogarithms.
- What functions occur beyond multiple polylogarithm?
- Is there an efficient method to evaluate these functions numerically everywhere?

Outline

Efficient numerical evaluation methods rely on three ingredients:

1. An (iterated) **integral representation**, used to transform the arguments into the region of convergence.
2. A (nested) **sum representation** defined in the region of convergence, which can be truncated and gives a numerical approximation.
3. Methods, which **accelerate** the convergence of the truncated series.

Numerical evaluation of the dilogarithm

The **dilogarithm**:

$$\operatorname{Li}_2(x) = \int_0^x \frac{dt_1}{t_1} \int_0^{t_1} \frac{dt_2}{1-t_2} = \sum_{n=1}^{\infty} \frac{x^n}{n^2}$$

Map into region $|x| \leq 1$ and $-1 \leq \operatorname{Re}(x) \leq 1/2$, using

$$\operatorname{Li}_2(x) = -\operatorname{Li}_2\left(\frac{1}{x}\right) - \frac{\pi^2}{6} - \frac{1}{2}(\ln(-x))^2, \quad \operatorname{Li}_2(x) = -\operatorname{Li}_2(1-x) + \frac{\pi^2}{6} - \ln(x)\ln(1-x).$$

Acceleration using Bernoulli numbers B_j :

$$\operatorname{Li}_2(x) = \sum_{j=0}^{\infty} \frac{B_j}{(j+1)!} (-\ln(1-x))^{j+1},$$

Multiple polylogarithms

Definition based on nested sums:

$$\text{Li}_{m_1, m_2, \dots, m_k}(x_1, x_2, \dots, x_k) = \sum_{n_1 > n_2 > \dots > n_k > 0} \frac{x_1^{n_1}}{n_1^{m_1}} \cdot \frac{x_2^{n_2}}{n_2^{m_2}} \cdot \dots \cdot \frac{x_k^{n_k}}{n_k^{m_k}}$$

Definition based on iterated integrals:

$$G(z_1, \dots, z_k; y) = \int_0^y \frac{dt_1}{t_1 - z_1} \int_0^{t_1} \frac{dt_2}{t_2 - z_2} \dots \int_0^{t_{k-1}} \frac{dt_k}{t_k - z_k}$$

Conversion:

$$\text{Li}_{m_1, \dots, m_k}(x_1, \dots, x_k) = (-1)^k G_{m_1, \dots, m_k} \left(\frac{1}{x_1}, \frac{1}{x_1 x_2}, \dots, \frac{1}{x_1 \dots x_k}; 1 \right)$$

Short hand notation:

$$G_{m_1, \dots, m_k}(z_1, \dots, z_k; y) = G(\underbrace{0, \dots, 0}_{m_1-1}, z_1, \dots, z_{k-1}, \underbrace{0, \dots, 0}_{m_k-1}, z_k; y)$$

Numerical evaluation of multiple polylogarithms

Use the integral representation

$$G_{m_1, \dots, m_k}(z_1, z_2, \dots, z_k; y) = \int_0^y \left(\frac{dt}{t}\right)^{m_1-1} \frac{dt}{t-z_1} \left(\frac{dt}{t}\right)^{m_2-1} \frac{dt}{t-z_2} \dots \left(\frac{dt}{t}\right)^{m_k-1} \frac{dt}{t-z_k}$$

to transform all arguments into a region, where we have a **converging power series expansion**:

$$G_{m_1, \dots, m_k}(z_1, \dots, z_k; y) = \sum_{j_1=1}^{\infty} \dots \sum_{j_k=1}^{\infty} \frac{1}{(j_1 + \dots + j_k)^{m_1}} \left(\frac{y}{z_1}\right)^{j_1} \frac{1}{(j_2 + \dots + j_k)^{m_2}} \left(\frac{y}{z_2}\right)^{j_2} \dots \frac{1}{(j_k)^{m_k}} \left(\frac{y}{z_k}\right)^{j_k}.$$

Use the **Hölder convolution** to accelerate the convergent series:

$$G(z_1, \dots, z_w; 1) = \sum_{j=0}^w (-1)^j G\left(1-z_j, 1-z_{j-1}, \dots, 1-z_1; \frac{1}{2}\right) G\left(z_{j+1}, \dots, z_w; \frac{1}{2}\right).$$

The two-loop electron self-energy

As a proof-of-concept consider the two-loop electron self-energy (Sabry, '62).

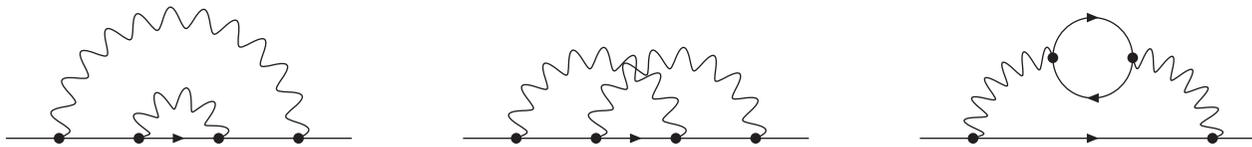
This is a quantity of physical relevance, although

- the self-energy is **not gauge-invariant**,
- the renormalised self-energy is **renormalisation-scheme dependent**,
- the bare self-energy is **not a pure function**.

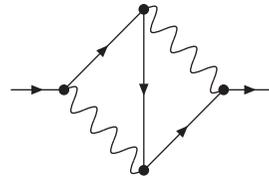
For concreteness, consider in **Feynman gauge** the **ϵ^0 -term** within dimensional regularisation and the **onshell-scheme**.

Diagrams

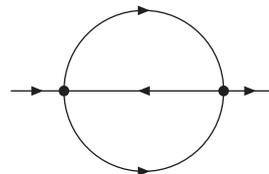
There are **three Feynman diagrams** contributing to the two-loop electron self-energy in QED with a single fermion:



All master integrals are (sub-) topologies of the **kite graph**:



One sub-topology is the **sunrise graph** with three equal non-zero masses:



The two-loop electron self-energy

Write for the two-loop electron self-energy

$$\Sigma_{\text{bare}}^{(2)} = \left(\frac{\alpha}{4\pi} \right)^2 \left(\Sigma_{\text{bare},V}^{(2)} \not{p} + \Sigma_{\text{bare},S}^{(2)} m \right),$$

Write the vector and the scalar part as a linear combination of master integrals I_j :

$$\Sigma_{\text{bare},V}^{(2)} = \sum_{j=1}^8 c_j^V I_j, \quad \Sigma_{\text{bare},S}^{(2)} = \sum_{j=1}^8 c_j^S I_j.$$

Laurent expansion in ϵ :

$$\Sigma_{\text{bare},V}^{(2)} = \sum_{k=-2}^{\infty} \epsilon^k \Sigma_{\text{bare},V}^{(2,k)}, \quad \Sigma_{\text{bare},S}^{(2)} = \sum_{k=-2}^{\infty} \epsilon^k \Sigma_{\text{bare},S}^{(2,k)}.$$

Pole terms and counterterms simple. Focus on $\Sigma_{\text{bare},V}^{(2,0)}$ and $\Sigma_{\text{bare},S}^{(2,0)}$.

The master integrals

There are 8 master integrals (Remiddi, Tancredi, '16).

Strategy for the master integrals:

- Derive the **differential equations** for the master integrals.
- Transform the system by a **non-algebraic redefinition** of the master integrals to an **ϵ -form**.
- Solve the system in terms of **iterated integrals**.

The differential equations

Set $x = p^2/m^2$.

Singularities of the differential equation are at $x \in \{0, 1, 9, \infty\}$.

Call these points the **cusps**.

The system of differential equations contains the **second-order irreducible** differential operator (Broadhurst, Fleischer, Tarasov '93)

$$L = x(x-1)(x-9) \frac{d^2}{dx^2} + (3x^2 - 20x + 9) \frac{d}{dx} + x - 3$$

Let ψ_1 and ψ_2 be **two independent solutions** of the homogeneous equation

$$L \psi = 0.$$

Periods

ψ_1 and ψ_2 can be taken as **periods of the elliptic curve**

$$E : w^2 - z(z+4) \left[z^2 + 2(1+x)z + (1-x)^2 \right] = 0.$$

Set

$$k^2 = \frac{16\sqrt{x}}{(1+\sqrt{x})^3(3-\sqrt{x})}, \quad k'^2 = 1 - k^2.$$

In a neighbourhood of $x = 0$ the periods may be taken as

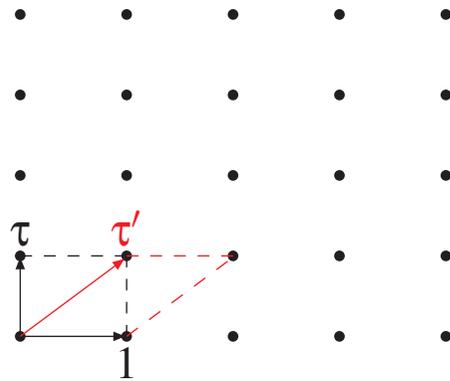
$$\psi_{1,0} = \frac{4K(k)}{(1+\sqrt{x})^{\frac{3}{2}}(3-\sqrt{x})^{\frac{1}{2}}}, \quad \psi_{2,0} = \frac{4iK(k')}{(1+\sqrt{x})^{\frac{3}{2}}(3-\sqrt{x})^{\frac{1}{2}}},$$

The **complete elliptic integral** $K(k)$ can be computed efficiently from the **arithmetic-geometric mean**

$$K(k) = \frac{\pi}{2 \operatorname{agm}(k', 1)}.$$

Bases of lattices

The periods ψ_1 and ψ_2 generate a lattice. Any other basis as good as (ψ_2, ψ_1) .
 Convention: Normalise $(\psi_2, \psi_1) \rightarrow (\tau, 1)$ where $\tau = \psi_2/\psi_1$.



Change of basis:
$$\begin{pmatrix} \psi'_2 \\ \psi'_1 \end{pmatrix} = \begin{pmatrix} a & b \\ c & d \end{pmatrix} \begin{pmatrix} \psi_2 \\ \psi_1 \end{pmatrix},$$

Transformation should be invertible:
$$\begin{pmatrix} a & b \\ c & d \end{pmatrix} \in \text{SL}_2(\mathbb{Z}),$$

In terms of τ and τ' :
$$\tau' = \frac{a\tau + b}{c\tau + d}$$

Modular forms

Denote by \mathbb{H} the **complex upper half plane**. A meromorphic function $f : \mathbb{H} \rightarrow \mathbb{C}$ is a **modular form** of modular weight k for $\mathrm{SL}_2(\mathbb{Z})$ if

(i) f transforms under Möbius transformations as

$$f\left(\frac{a\tau + b}{c\tau + d}\right) = (c\tau + d)^k \cdot f(\tau) \quad \text{for} \quad \begin{pmatrix} a & b \\ c & d \end{pmatrix} \in \mathrm{SL}_2(\mathbb{Z})$$

(ii) f is holomorphic on \mathbb{H} ,

(iii) f is holomorphic at ∞ .

Congruence subgroups

Apart from $\mathrm{SL}_2(\mathbb{Z})$ we may also look at congruence **subgroups**, for example

$$\Gamma_0(N) = \left\{ \begin{pmatrix} a & b \\ c & d \end{pmatrix} \in \mathrm{SL}_2(\mathbb{Z}) : c \equiv 0 \pmod{N} \right\}$$

$$\Gamma_1(N) = \left\{ \begin{pmatrix} a & b \\ c & d \end{pmatrix} \in \mathrm{SL}_2(\mathbb{Z}) : a, d \equiv 1 \pmod{N}, c \equiv 0 \pmod{N} \right\}$$

$$\Gamma(N) = \left\{ \begin{pmatrix} a & b \\ c & d \end{pmatrix} \in \mathrm{SL}_2(\mathbb{Z}) : a, d \equiv 1 \pmod{N}, b, c \equiv 0 \pmod{N} \right\}$$

Modular forms for congruence subgroups: Require “**nice**” transformation properties only for subgroup Γ (plus holomorphicity on \mathbb{H} and at the cusps).

Relevant to this talk: **Modular forms for $\Gamma_1(6)$.**

Variables

Recall

$$x = \frac{p^2}{m^2}.$$

Set

$$\tau_0 = \frac{\Psi_{2,0}}{\Psi_{1,0}}, \quad q_0 = e^{2i\pi\tau_0}.$$

Change variable from x to τ_0 (or q_0) (Bloch, Vanhove, '13).

The differential equation for the master integrals reads then

$$\frac{d}{d\tau_0} \vec{I} = \varepsilon A(\tau_0) \vec{I},$$

where $A(\tau_0)$ is an ε -independent 8×8 -matrix whose entries are modular forms.

Iterated integrals

For $\omega_1, \dots, \omega_k$ differential 1-forms on a manifold M and $\gamma: [0, 1] \rightarrow M$ a path, write for the **pull-back** of ω_j to the interval $[0, 1]$

$$f_j(\lambda) d\lambda = \gamma^* \omega_j.$$

The **iterated integral** is defined by (Chen '77)

$$I_\gamma(\omega_1, \dots, \omega_k; \lambda) = \int_0^\lambda d\lambda_1 f_1(\lambda_1) \int_0^{\lambda_1} d\lambda_2 f_2(\lambda_2) \dots \int_0^{\lambda_{k-1}} d\lambda_k f_k(\lambda_k).$$

Example 1: **Multiple polylogarithms** (Goncharov '98)

$$\omega_j = \frac{d\lambda}{\lambda - z_j}.$$

Example 2: **Iterated integrals of modular forms** (Brown '14): $f_j(\tau)$ a modular form,

$$\omega_j = 2\pi i f_j(\tau) d\tau.$$

Iterated integrals of modular forms

Modular forms have a *q-expansion*. Using

$$2\pi i d\tau = \frac{dq}{q}$$

we may *integrate term-by-term* and obtain the *q-expansion* of the master integrals.

For example, for the ε^2 -term of the sunrise integral one finds

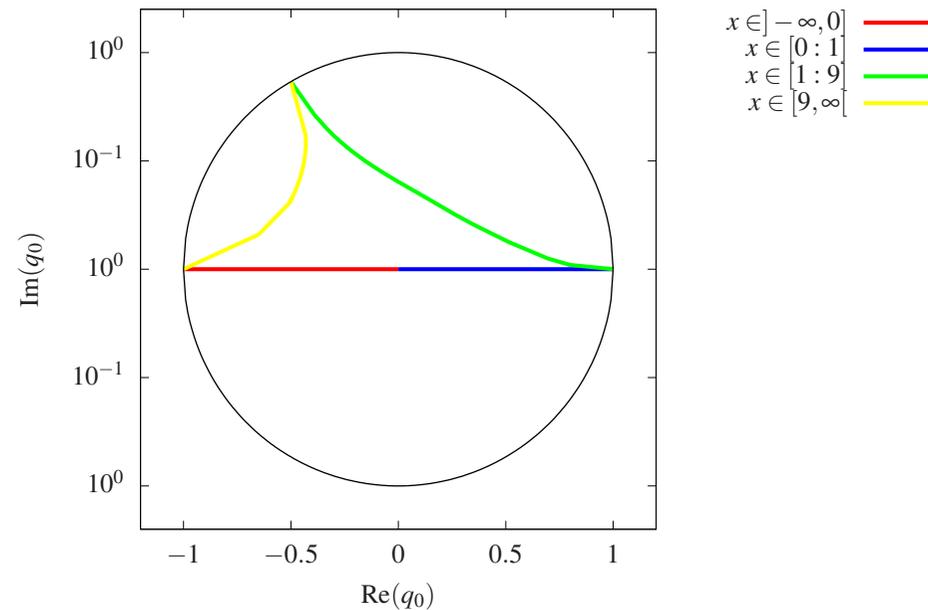
$$I_6^{(2)} = 3 \operatorname{Cl}_2\left(\frac{2\pi}{3}\right) - 3\sqrt{3} \left[q - \frac{5}{4}q^2 + q^3 - \frac{11}{16}q^4 + \frac{24}{25}q^5 - \frac{5}{4}q^6 + \frac{50}{49}q^7 - \frac{53}{64}q^8 + q^9 \right] + O(q^{10})$$

We may *truncate* the *q-series* and *evaluate* the resulting polynomial *numerically*.

Convergence

We defined q_0 such that $q_0 = 0$ for $x = 0$.

For which values $x \in \mathbb{R}$ do we have $|q_0| < 1$?



We have $|q_0| < 1$ for $x \in \mathbb{R} \setminus \{1, 9, \infty\}$.

Expansion parameters

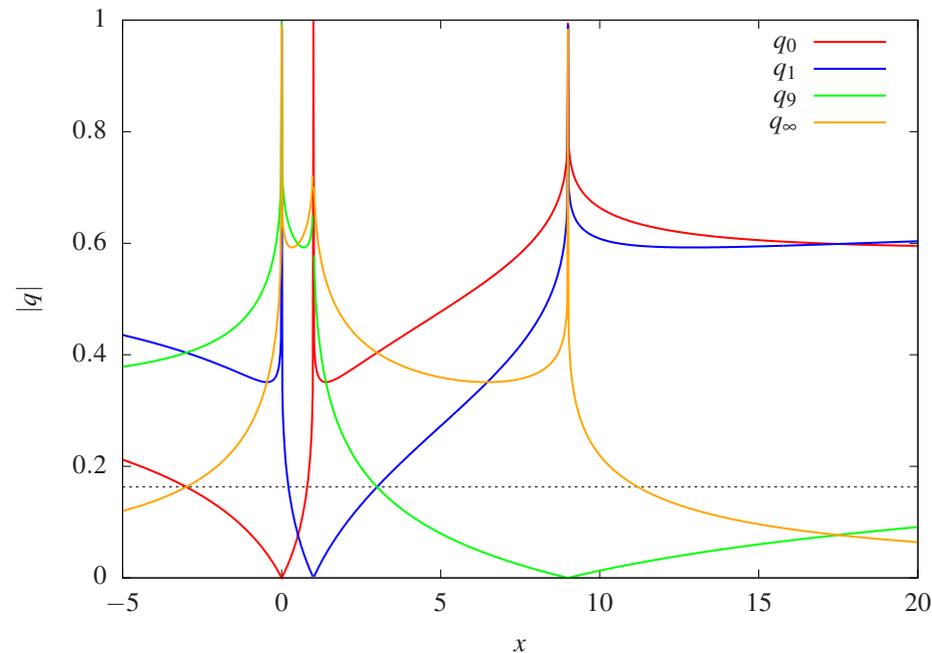
Recall that the periods ψ_1 and ψ_2 are **not uniquely determined**.

Can define q_0, q_1, q_9 and q_∞ such that

- the integration kernels are modular forms of $\Gamma_1(6)$,
- $q_j = 0$ for $x = j$.

Expansion parameters

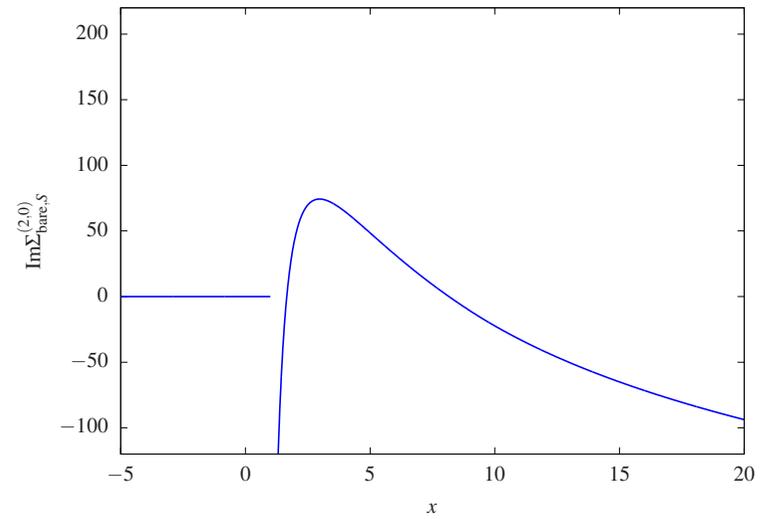
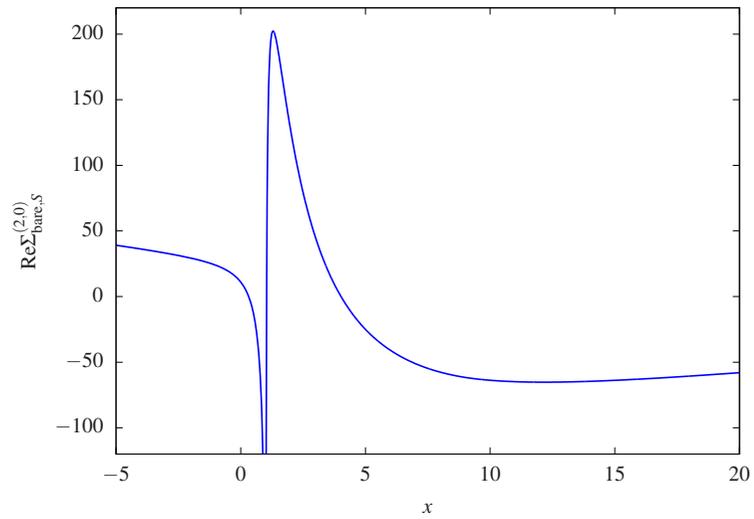
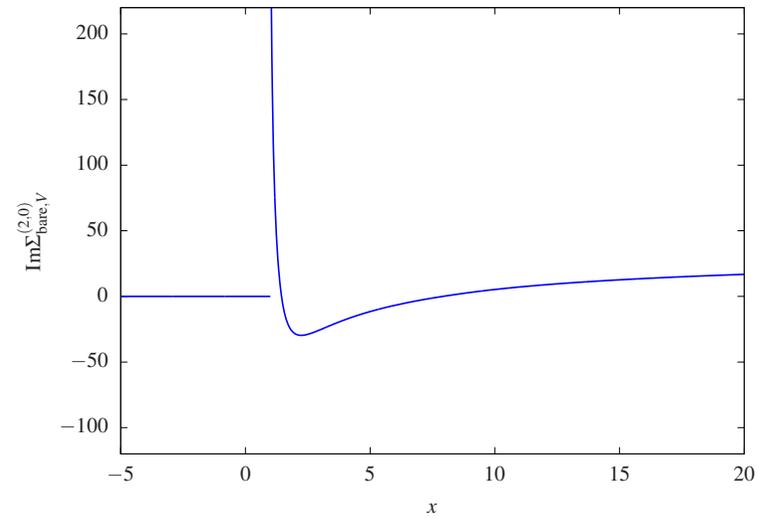
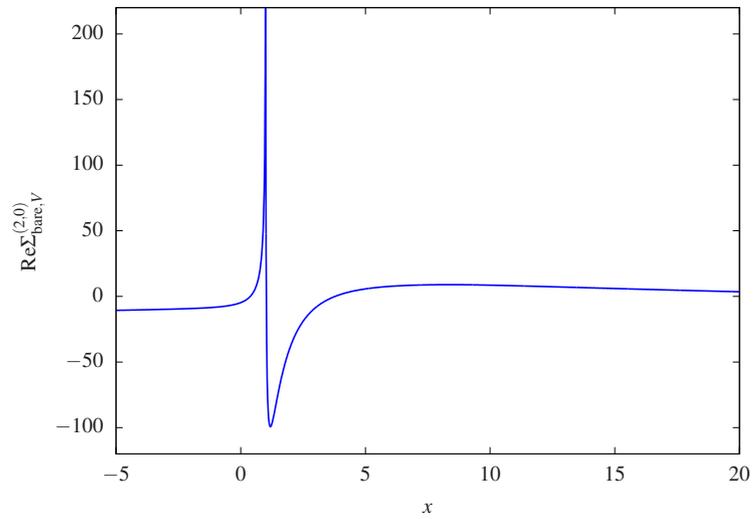
The absolute values of the variables q_j for $j \in \{0, 1, 9, \infty\}$ as a function of x .



There is always a choice such that $|q_j| \lesssim 0.163$ for all values of x .

Truncation of the q -series to order $O(q^{30})$ gives for the finite part of the self-energy a relative precision better than 10^{-20} for all real values p^2/m^2 .

Numerical results



Wrap-up

- Periods are calculated from the arithmetic-geometric mean.
- Master integrals are calculated from truncated q -expansions.

Conclusions

- Loop integrals with **masses important** for top, W/Z - and H -physics.
- May involve **elliptic sectors** from two loops onwards.
- There is a class of Feynman integrals evaluating to **iterated integrals of modular forms**.
- These can be **evaluated efficiently** from their q -expansions.
- We may expect more results in the near future.

Back-up

Relative precision for $\Sigma_{\text{bare},S}^{(2,0)}$ (similar for $\Sigma_{\text{bare},V}^{(2,0)}$):

