# On the Application of Intersection Theory to Feynman Integrals: The Univariate Case 

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A key step in modern high energy physics scattering amplitudes computation is to express the latter in terms of a minimal set of Feynman integrals using linear relations.
In this work we present an innovative approach (first introduced in [1] and further developed in $[2,3,4]$ ) based on intersection theory, in order to achieve this decomposition. This allows for the direct computation of the reduction, projecting integrals appearing in the scattering amplitudes onto an integral basis in the same fashion as vectors may be projected onto a vector basis. Specifically, we will derive and discuss few identities between maximally cut Feynman integrals, showing their direct decomposition.
This contribution will focus on the univariate part of the story, with the multivariate generalisation being discussed in a different contribution by Gasparotto and Mandal [5], and for other contributions see also $[6,7,8,9,10,11]$.

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## 1. Introduction

When analyzing the enormous amount of data gathered by particle scattering experiments such as those taking place at the Large Hadron Collider at CERN, high precision theoretical predictions based on the Standard Model of particle physics are needed, in order to be able to detect potential tiny discrepancies between theory and experiment, which may hold the key to new physics.

In the language of quantum field theory, precise predictions for the outcome of a particle scattering process require the computation of Feynman integrals, which are mathematical objects given as

$$
\begin{equation*}
I=\int \frac{\mathrm{d}^{d} k_{i}}{\pi^{d / 2}} \cdots \int \frac{\mathrm{~d}^{d} k_{L}}{\pi^{d / 2}} \frac{N(k)}{D_{1}^{a_{1}}(k) D_{2}^{a_{2}}(k) \cdots D_{P}^{a_{P}}(k)} \tag{1.1}
\end{equation*}
$$

where the $D$ s are propagators of the form $D_{i}=(k+p)^{2}-m^{2}, d=4-2 \varepsilon$ is the space-time dimensionality, $k$ and $p$ are $d$-dimensional momenta (internal and external), the $a_{i}$ are integer propagator powers, $L$ is the number of loops, and $P$ the number of propagators.

Higher precision theoretical predictions require the computation of Feynman integrals with more loops and legs than has been done previously. Adding a loop or a leg to a Feynman integral, implies a huge increase in complexity. For that reason it is important to reduce the set of integrals that has to be computed to a minimum number. Different Feynman integrals in the same family, i.e. integrals differing only by the values of the powers $a_{i}$, are related through linear relations, allowing for the expression of any integral in a integral family as a linear combination of a minimal set of integrals, known as master integrals

$$
\begin{equation*}
I=\sum_{i=1}^{v} c_{i} I_{i} \tag{1.2}
\end{equation*}
$$

with $v$ denoting the number of master integrals in the family.
The traditional way of deriving these linear relations, is through the use of integration-byparts relations [12], which may be combined in a systematic way using Laporta's algorithm [13]. This algorithm requires the solution of a huge linear system of equations relating the integrals in order to extract the relations needed. For more complicated Feynman integrals this becomes a significant computational challenge and therefore the search for a more direct way to extract relations on the form of eq. (1.2) becomes of higher interest and importance. The approach based on the intersection number, first applied to Feynman integrals in refs. [1] and further developed in $[2,3,4]$ as well as $[14,15,16]$, is a promising piece of progress in that direction.

## 2. The loop-by-loop Baikov representation

Before discussing the intersection theory, let us briefly discuss the loop-by-loop Baikov representation. The Baikov representation [17] (see also [18, 19, 20, 21, 22, 23]) for Feynman integrals is a parametric representation, meaning that the integrations are over a number of scalar variables, rather than over the $d$-dimensional loop-momenta of eq. (1.1). Other examples of parametric representations of Feynman integrals are Feynman and Schwinger parametrizations. Baikov representation is characterized by the fact that the integration variables equal the propagators of the original
integral. A Feynman integral in standard Baikov representation is given by

$$
\begin{equation*}
I=\frac{J(-i)^{L} \pi^{(L-n) / 2} \mathscr{G}^{(E-d+1) / 2}}{\prod_{l=1}^{L} \Gamma((d+1-E-l) / 2)} \int \frac{N(x) \mathscr{B}^{(d-E-L-1) / 2} \mathrm{~d}^{n} x}{x_{1}^{a_{1}} \cdots x_{\mathrm{P}}^{a_{\mathrm{P}}}} \tag{2.1}
\end{equation*}
$$

Here $L$ is the number of loops, and $E$ the number of independent external momenta.

$$
\begin{equation*}
n=E L+L(L+1) / 2 \tag{2.2}
\end{equation*}
$$

is the number of Baikov variables which again equal the number of scalar products that can be formed between the $L$ loop-momenta and the full set of internal and external momenta. The two polynomials are

$$
\begin{equation*}
\mathscr{G}=G\left(p_{1}, \ldots, p_{E}\right), \quad \mathscr{B}=G\left(k_{1}, \ldots, k_{L}, p_{1}, \ldots, p_{E}\right) \tag{2.3}
\end{equation*}
$$

where $G(y)$ is the Gram determinant of the set $y$ with itself, that is the determinant of the matrix formed by all the scalar products $y_{i} \cdot y_{j}$. Finally $J$ is the Jacobian of a variable change from the independent scalar products to the Baikov variables, that is $J^{-1}=\operatorname{det}\left(d P_{i} / d(k \cdot p)_{j}\right)$. The expression for $J$ depends on the exact definition of the propagators, but usually $J= \pm 2^{L-n}$. We see that in most cases $n>P$, which means that we have to introduce some artificial propagator-like objects to play the role of the remaining Baikov variables.

The loop-by-loop Baikov representation [23] applies the variable changes needed to go to Baikov representation to each loop individually rather than to the whole diagram at once. We define $E_{l}$ as the number of momenta external to the $l$ th loop, after performing the parametrization on the loops labelled with $k_{i>l}$. This number may include the loop-momenta of lower-numbered loops, and also external momenta not directly attached to the $k_{l}$-loop which may be induced by the previous parametrizations. See fig. 1 for an example.

In this loop-by-loop case we get the number of integration variables to be

$$
\begin{equation*}
\tilde{n}=L+\sum_{i=1}^{L} E_{i} \tag{2.4}
\end{equation*}
$$

which in most cases will be smaller (and never larger) than the number given by eq. (2.2).
In that case, the general loop-by-loop Baikov parametrization of the integral of eq. (1.1) is

$$
\begin{equation*}
I=\frac{J(-i)^{L} \pi^{(L-\tilde{n}) / 2}}{\prod_{l=1}^{L} \Gamma\left(\left(d-E_{l}\right) / 2\right)} \int \frac{N(x)\left(\prod_{l=1}^{L} \mathscr{G}_{l}^{\left(E_{l}-d+1\right) / 2} \mathscr{B}_{l}^{\left(d-E_{l}-2\right) / 2}\right) \mathrm{d}^{\tilde{n}} x}{x_{1}^{a_{1}} \cdots x_{\mathrm{P}}^{a_{\mathrm{P}}}} \tag{2.5}
\end{equation*}
$$

The corresponding polynomials are

$$
\begin{equation*}
\mathscr{G}_{l}=G\left(\left\{q_{l}\right\}\right), \quad \mathscr{B}_{l}=G\left(k_{l},\left\{q_{l}\right\}\right) \tag{2.6}
\end{equation*}
$$

where $\left\{q_{l}\right\}$ is the set of momenta external to the $l$ th loop. $J$ is defined as above, and is usually given as $J= \pm 2^{L-\tilde{n}}$.

If we take all loops to depend on all lower loop-momenta and all external momenta, we get $E_{l}=E+l-1$ implying that $\tilde{n}=n$, and it is not hard to see that eq. (2.5) will reduce to eq. (2.1) in that case.


Figure 1: This figure shows the massless pentabox integral, first studied in [24, 25]. It is a good example for showing the loop-by-loop Baikov parametrizartion. When doing loop-by-loop one has to choose which loop to start from. Starting from the $q_{1}$-loop-corresponding to $k_{1}=q_{2}, k_{2}=q_{1}$ gives the set of propagators containing $k_{2}$ to be $k_{2}^{2},\left(k_{2}+p_{1}\right)^{2},\left(k_{2}+p_{1}+p_{2}\right)^{2},\left(k_{2}-k_{1}\right)^{2}$. From this we see that the set of momenta external to that loop is $\left\{p_{1}, p_{2}, k_{1}\right\}$ corresponding to $E_{2}=3$. The Grammatrices $\mathscr{B}_{2}$ and $\mathscr{G}_{2}$ will contain the scalar products between $k_{1} \cdot p_{1}$ and $k_{1} \cdot p_{2}$ which together with the remaining propagators show that the $k_{1}$ integration depends on all four independent external momenta, so $E_{1}=4$. This gives by eq. (2.4) $\tilde{n}=2+4+3=9$ showing that only one auxiliary Baikov variable is needed, which we may choose as $x_{9}=\left(k_{1}+p_{1}\right)^{2}$.
Alternatively one could have chosen to start from the other side so $k_{1}=q_{1}, k_{2}=q_{2}$. In this case the propagators containing $k_{2}$ would be $k_{2}^{2},\left(k_{2}-p_{5}\right)^{2},\left(k_{2}-p_{5}-p_{4}\right)^{2},\left(k_{2}-p_{5}-p_{4}-p_{3}\right)^{2},\left(k_{1}-k_{2}\right)^{2}$ corresponding to $E_{2}=4$. This time $\mathscr{B}_{2}$ and $\mathscr{G}_{2}$ would introduce the scalar products $k_{1} \cdot p_{5}, k_{1} \cdot p_{4}$, and $k_{1} \cdot p_{3}$ which together with those coming from the remaining propagators gives $E_{1}=4$ corresponding to $\tilde{n}=10$ i.e. one more than in the other case. This shows the importance of a clever choice of parametrization of the momenta when applying the loop-by-loop Baikov representation.

## 3. Theory

Eq. (1.2) may remind the reader of writing a vector in a vector space in terms of a basis of vectors for that space. And indeed the space of Feynman integrals do live up to all requirements for being a vector space [1], so referring to the set of master integrals as a basis is not merely a metaphor.

For vector spaces equipped with an inner product, expressing a vector $\langle v|$ in terms of basis vectors $\left\langle v_{i}\right|$ is an easy task, since

$$
\begin{equation*}
\langle v|=\sum_{i j}\left\langle v \mid v_{j}\right\rangle\left(\mathbf{C}^{-1}\right)_{j i} i v_{i} \mid \quad \text { with } \quad \mathbf{C}_{i j}=\left\langle v_{i} \mid v_{j}\right\rangle \tag{3.1}
\end{equation*}
$$

and thus

$$
\begin{equation*}
\langle v|=\sum_{i} c_{i}\left\langle v_{i}\right| \quad \Rightarrow \quad c_{i}=\sum_{j}\left\langle v \mid v_{j}\right\rangle\left(\mathbf{C}^{-1}\right)_{j i} \tag{3.2}
\end{equation*}
$$

So we realize that the computationally intensive IBP-based algorithm traditionally used to extract the coefficients of eq. (1.2), may be avoided if we find a way to introduce (what corresponds to) an inner product between Feynman integrals.

To introduce this terminology, we must discuss parametric representations of Feynman integrals, such as those discussed in the previous section. In these representations Feynman integrals
are objects of the form

$$
\begin{equation*}
I=\int_{\mathscr{C}} u(\mathbf{z}) \varphi(\mathbf{z}) \quad \text { with } \quad u(\mathbf{z})=\prod_{i} \mathscr{B}_{i}(\mathbf{z})^{\gamma_{i}} \tag{3.3}
\end{equation*}
$$

where $\mathbf{z}$ is a set of variables, $\mathscr{B}_{i}$ are polynomials in the $\mathbf{z}$, the $\gamma_{i}$ are non-integers, and $\varphi$ is a differential form in the $\mathbf{z}$, with a prefactor that is a rational function $\hat{\varphi}(\mathbf{z})$ with the property that all potential poles of $u(\mathbf{z}) \varphi(\mathbf{z})$ are regulated by the $\gamma_{i}$.

Feynman integrals in the Baikov representation (eqs. (2.1) and (2.5)) are of that form, and so are Feynman integrals in the Lee-Pomeransky form [26] of Feynman parametrization, as long as potential poles in the individual variables are regulated [2] $z_{i}^{a_{i}} \rightarrow z_{i}^{a_{i}+\rho_{i}}$ where the regulators $\rho_{i}$ may be put to zero at the end of the computation. We will focus on the Baikov parametrization in the following.

In addition, also derivatives of Feynman integrals with respect to kinematical variables, and Feynman integrals in shifted dimensions ( $d \rightarrow d \pm 2 n$ ) are of the form of eq. (3.3), so a unified description of integrals of that form will be applicable to all these different objects.

Building on this terminology we may define

$$
\begin{equation*}
\omega=d \log (u) \tag{3.4}
\end{equation*}
$$

to be of use later. This object has the property that the number of master integrals $v$ may be found as [26]

$$
\begin{equation*}
v=\text { number of solutions to " } \omega=0 \text { " } \tag{3.5}
\end{equation*}
$$

something that will be used extensively in the following.
The integral of eq. (3.3) may be reexpressed as a pairing between a twisted co-cycle $\langle\varphi|$ and a twisted cycle $\mid \mathscr{C}]$

$$
\begin{equation*}
\left.I=\int_{\mathscr{C}} u(\mathbf{z}) \varphi(\mathbf{z})=\langle\varphi| \mathscr{C}\right] \tag{3.6}
\end{equation*}
$$

where the multi-valued function $u$ is implicit in the definition of the pairing. The $\langle\varphi|$ and $\mid \mathscr{C}]$ should be understood not as being necessarily equal to the $\varphi$ and $\mathscr{C}$ of eq. (3.3), but rather as representatives of the group of forms and integration contours respectively, which integrate to the same $I$.

It is likewise possible to define a dual twisted cycle and cocycle as $\int_{\mathscr{C}} u^{-1}(\mathbf{z}) \varphi(\mathbf{z})=[\mathscr{C}|\varphi\rangle$. This allows for the introduction of pairings between the cycles and the co-cycles themselves $\left\langle\varphi_{L} \mid \varphi_{R}\right\rangle$ and $\left[\mathscr{C}_{L} \mid \mathscr{C}_{R}\right]$ which are known as intersection numbers. For the complete definition of the intersection numbers, we must refer to the mathematical literature [27, 28]. In the following we will focus solely on the intersection number of the twisted co-cycles $\left\langle\varphi_{L} \mid \varphi_{R}\right\rangle$, which play the role of the inner product we were searching for.

With this definition in place, we now have a way of extracting the master integral coefficients of eq. (1.2). This can be done [1] using eq. (3.2) as

$$
\begin{equation*}
\left.\langle\varphi| \mathscr{C}]=\sum_{i} c_{i}\left\langle e_{i}\right| \mathscr{C}\right] \quad \Rightarrow \quad c_{i}=\sum_{j}\left\langle\varphi_{L} \mid h_{j}\right\rangle\left(\mathbf{C}^{-1}\right)_{j i} \tag{3.7}
\end{equation*}
$$

with $\mathbf{C}_{i j}=\left\langle e_{i} \mid h_{j}\right\rangle$ as in eq. (3.1) and with $\left\langle e_{i}\right|$ and $\left|h_{i}\right\rangle$ being the bases of cocycles and dual cocycles respecitively.

## 4. Univariate intersection numbers

Let us start by the univariate case in which the integral of eq. (3.3) is with respect to just one variable $z$. For that case the intersection number is given as

$$
\begin{equation*}
\left\langle\varphi_{L} \mid \varphi_{R}\right\rangle=\sum_{p \in \mathscr{B}} \operatorname{Res}_{z=p}\left(\psi_{p} \varphi_{R}\right) \tag{4.1}
\end{equation*}
$$

with

$$
\begin{equation*}
(d+\omega) \psi_{p}=\varphi_{L} \tag{4.2}
\end{equation*}
$$

The sum goes over $\mathscr{P}$, the set of poles of $\omega$, and $\psi$ is defined as the solution to eq. (4.2), but as indicated by the subscript all that is needed for practical applications is a solution that is valid around the point $p$, as the residue operation is local, so the defining equation can be solved using a series ansatz.

Let us illustrate this formula and eq. (3.7) on the Gauss' hypergeometric function

$$
\begin{equation*}
{ }_{2} F_{1}(a, b, c ; x)=\int_{0}^{1} u(z) \mathrm{d} z \quad \text { with } \quad u(z)=\frac{\Gamma(c)}{\Gamma(b) \Gamma(c-b)} z^{b-1}(1-z)^{c-b-1}(1-x z)^{-a} \tag{4.3}
\end{equation*}
$$

This corresponds to eq. (3.3) with $\mathscr{C}=[0,1]$ and $\varphi=1 \mathrm{~d} z$.
Hypergeometric functions are related through relations called contiguity relations, which relate different ${ }_{2} F_{1}$ for which the indices $a, b, c$ differ by integers. This is the equivalent of the linear relations relating Feynman integrals in the same family. One example of a contiguity relation is

$$
\begin{equation*}
{ }_{2} F_{1}(a, b+1, c+1 ; x)=c_{12} F_{1}(a, b+1, c ; x)+c_{2}{ }_{2} F_{1}(a, b, c ; x) \tag{4.4}
\end{equation*}
$$

The values of $c_{1}$ and $c_{2}$ are known, but we will try to derive them using the intersection-based formalism. With the $u$ given above, these three integrals correspond to

$$
\begin{equation*}
\hat{\varphi}=\frac{c z}{b}, \quad \hat{e}_{1}=\frac{(1+b-c) z}{b(z-1)}, \quad \hat{e}_{2}=1 \tag{4.5}
\end{equation*}
$$

with $\varphi=\hat{\varphi} \mathrm{d} z$ and $e_{i}=\hat{e}_{i} \mathrm{~d} z$.
The first step is to compute $\omega$ :

$$
\begin{equation*}
\omega=\mathrm{d} \log (u)=\left(\frac{b-1}{z}+\frac{b-c+1}{z-1}+\frac{a x}{1-z x}\right) \mathrm{d} z \tag{4.6}
\end{equation*}
$$

Solving for $\omega=0$ gives two solutions corresponding to the $v=2$ master integrals appearing on the RHS of eq. (4.4).

In order to be able to use the projecting formula eq. (3.7), we also need a set of dual forms $\left|h_{i}\right\rangle$. They can be chosen almost arbitrarily as long as no poles are present that are not poles of $\omega$. We chose

$$
\begin{equation*}
\hat{h}_{1}=\frac{1}{z}-\frac{1}{z-1}, \quad \hat{h}_{2}=\frac{1}{z}-\frac{x}{x z-1} . \tag{4.7}
\end{equation*}
$$

To apply eq. (3.7) we need to compute all intersection numbers between the forms of eqs. (4.5) and (4.7), i.e.:

$$
\begin{equation*}
\left\{\left\langle\varphi \mid h_{1}\right\rangle,\left\langle\varphi \mid h_{2}\right\rangle,\left\langle e_{1} \mid h_{1}\right\rangle,\left\langle e_{1} \mid h_{2}\right\rangle,\left\langle e_{2} \mid h_{1}\right\rangle,\left\langle e_{2} \mid h_{2}\right\rangle\right\} \tag{4.8}
\end{equation*}
$$

Let us focus on $\left\langle e_{2} \mid h_{1}\right\rangle$. To find the intersection number we need to perform the sum of eq. (4.1). It goes over the set of poles of $\omega$ i.e.

$$
\begin{equation*}
\mathscr{P}=\left\{0,1, \frac{1}{x}, \infty\right\} \tag{4.9}
\end{equation*}
$$

Let us start by the contribution from $z=0$. In that point we need to find a solution $\psi_{0}$ to eq. (4.2) on the form of a series ansatz, that is

$$
\begin{equation*}
\psi_{0}=\sum_{i} \kappa_{i} z^{i} \tag{4.10}
\end{equation*}
$$

solving eq. (4.2) order by order in $z$, we find the solution

$$
\begin{equation*}
\kappa_{\leq 0}=0, \quad \kappa_{1}=\frac{1}{b}, \quad \kappa_{2}=\frac{c-b-1-a x}{b(b+1)}, \quad \ldots \tag{4.11}
\end{equation*}
$$

and inserting that solution into the residue of eq. (4.1) gives

$$
\begin{equation*}
\operatorname{Res}_{z=0}\left(\psi_{0} h_{1}\right)=0 \tag{4.12}
\end{equation*}
$$

The residues in $z=1$ and $z=\frac{1}{x}$ are likewise 0 , only the pole in infinity contributes with the value $\frac{1}{c-a-1}$, and therefore

$$
\begin{equation*}
\left\langle e_{2} \mid h_{1}\right\rangle=\sum_{p \in \mathscr{P}} \operatorname{Res}_{z=p}\left(\psi_{p} h_{1}\right)=\frac{1}{c-a-1} . \tag{4.13}
\end{equation*}
$$

The same computation can be done for the other intersection numbers, with the results

$$
\begin{align*}
\left\langle\varphi \mid h_{1}\right\rangle & =\frac{c(a-a x+b x)}{b(c-a)(c-a-1) x}, & \left\langle\varphi \mid h_{2}\right\rangle & =\frac{c(c+x+b x-c x-1)}{b(c-a)(c-a-1) x^{2}} \\
\left\langle e_{1} \mid h_{1}\right\rangle & =\frac{b-a}{b(c-a-1)}, & \left\langle e_{1} \mid h_{2}\right\rangle & =\frac{1+b-c}{b(c-a-1) x},  \tag{4.14}\\
\left\langle e_{2} \mid h_{1}\right\rangle & =\frac{1}{c-a-1}, & \left\langle e_{2} \mid h_{2}\right\rangle & =\frac{1}{(c-a-1) x} .
\end{align*}
$$

Inserting these results into the projection formula eq. (3.7), we get the results

$$
\begin{equation*}
c_{1}=\frac{c(1-x)}{x(a-c)}, \quad c_{2}=\frac{-c}{x(a-c)} \tag{4.15}
\end{equation*}
$$

which makes the contiguity relation of eq. (4.4) true, something that can be checked with the literature or numerically.

The same approach can be applied to the most general univariate generalization of the ${ }_{2} F_{1}$, known as the Lauricella $F_{D}$ function [29] defined as

$$
\begin{align*}
& F_{D}\left(a, b_{1}, \ldots, b_{n}, c ; x_{1}, \ldots, x_{n}\right) \\
= & \frac{\Gamma(c)}{\Gamma(a) \Gamma(c-a)} \int_{0}^{1} z^{a-1}(1-z)^{c-a-1} \prod_{i}^{n}\left(1-x_{i} z\right)^{-b_{i}} \mathrm{~d} z \tag{4.16}
\end{align*}
$$

This integral family has $n+1$ master integrals, the coefficients of which can be extracted using eq. (3.7) as above.

## 5. The (maximal cut of the) double box



Figure 2: The massless doublebox

After the purely mathematical example of the ${ }_{2} F_{1}$ let us try (following ref. [2]) an example relevant for real life particle physics, the double box seen on fig. 2. This is a Feynman integral on the form of eq. (1.1), and it is given in terms of the propagators

$$
\begin{gather*}
D_{1}=k_{1}^{2}, \quad D_{2}=\left(k_{1}-p_{1}\right)^{2}, \quad D_{3}=\left(k_{1}-p_{1}-p_{2}\right)^{2}, \quad D_{4}=\left(k_{1}-k_{2}\right)^{2} \\
D_{5}=\left(k_{2}-p_{1}-p_{2}\right)^{2}, \quad D_{6}=\left(k_{2}+p_{4}\right)^{2}, \quad D_{7}=k_{2}^{2}, \quad z=\left(k_{2}-p_{1}\right)^{2} \tag{5.1}
\end{gather*}
$$

where the $z$ represents an irreducible scalar product (ISP).
We will analyze it on the maximal cut, in which case the loop-by-loop version [23] of Baikov representation given by eq. (2.5) allows for a univariate representation:

$$
\begin{equation*}
I_{\mathrm{dbox}}=\int_{\mathscr{C}} u \mathrm{~d} z \quad \text { with } \quad u \propto z^{\frac{d}{2}-3}(z+s)^{2-\frac{d}{2}}(z-t)^{d-5} \tag{5.2}
\end{equation*}
$$

We see that this integral is equivalent to the ${ }_{2} F_{1}$ defined in eq. (4.3) up to a variable change, and therefore the rest of the computation will proceed along similar lines.

Let us say we want to reduce the integral having two powers in the numerator of the ISP corresponding to the variable $z$, to a basis of master integrals with zero and one powers respectively. That is

$$
\begin{equation*}
I_{1111111 ;-2}=c_{0} I_{1111111 ; 0}+c_{1} I_{1111111 ;-1} \tag{5.3}
\end{equation*}
$$

corresponding to the co-cycles $\varphi=z^{2} \mathrm{~d} z, \varphi_{0}=\mathrm{d} z$, and $\varphi_{1}=z \mathrm{~d} z$. Going through the same procedure as above, tells us that for using eq. (3.7) we need to compute six intersection numbers, those of these three co-cycles with a set of dual co-cycles. Computing those intersection numbers and combining the according to eq. (3.7) gives the results

$$
\begin{equation*}
c_{0}=\frac{(d-4) s t}{2(d-3)}, \quad c_{1}=\frac{2 t-3(d-4) s}{2(d-3)} \tag{5.4}
\end{equation*}
$$

in agreement with the results obtained in the literature and by standard IBP-based codes such as FIRE [30].

With the same method we were able to get correct relations on the maximal cut for a large number of Feynman integrals, most of which are depicted in fig. 3.



























$$
\stackrel{p_{1}}{p_{1}} \text { Geaceec]}
$$



Figure 3: Univariate examples from ref. [2].

## 6. Discussion

In this work we elaborated on the relation between Feynman integrals and intersection theory, first presented in [1] and further developed in [2, 3, 4].
The focus of this contribution was on the univariate version of the theory, and for the multivariate generalization see [3, 4] or the contribution [5].

We showed how Feynman integrals have a natural vector space structure with an inner product, allowing for new ways to compute their reduction onto a basis of master integrals.
In particular, we described the evaluation of univariate intersection numbers for twisted cocycles, which are the key ingredients of the master decomposition formula eq. (3.7). We applied the master decomposition formula to derive contiguity relations between hypergeometric functions and, by means of the loop-by-loop Baikov representation (the properties of which are further outlined in this work), to a large number of Feynman integrals.
Interesting future studies would be to systematically apply intersection theory to integral parameterizations different from Baikov, as well as refining the algorithm for the computation of the intersection numbers, for instance in ways that remove the explicit solution of the differential equation of eq. (4.2). Additionally fruitful results might be obtained by combining these promising new techniques with finite field methods.

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