MBsums — A Mathematica PACKAGE FOR THE REPRESENTATION OF MELLIN–BARNES INTEGRALS BY MULTIPLE SUMS*

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Feynman integrals may be represented by the Mathematica packages AMBRE and MB as multiple Mellin–Barnes integrals. With the Mathematica package MBsums we transform these Mellin–Barnes integrals into multiple sums.

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1. Introduction

In recent years, there was a remarkable progress in the development of (semi-)automatized software for the numerical calculation of arbitrary, complicated Feynman integrals. Basically, two approaches are advocated. One relies on a sector decomposition. For an introduction given at this conference and for further references see [1]. We will report on the other approach, based on Mellin–Barnes representations [2–12]. When [9] appeared in 2005, several unsolved problems of different complexity existed. We mention non-planar diagrams, the massive case, multi-loop tensor integrals, Minkowskian kinematics. For all of the items, a progress is reported in [2], based on the source-open software AMBRE/MB [8, 13, 14]. An alternative is the direct analytical evaluation of MB-integrals. This is difficult. But in view of the recent progress in algebraically summing up infinite sums by the Linz group's computer algebra algorithms for nested sums and products, one might hope

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to achieve a breakthrough [15]; certainly only if the result leads to appropriate classes of functions. The idea is to apply the Linz group's algorithms (as well as those of others, *e.g.* [16]) to sums of residues after applying Cauchy's theorem [17]. A first attempt was reported in [18].

The automatized derivation of the multiple sums for a given MB-integral is certainly the easier part of the task, but it is the first step. We report here on a first version of the Mathematica program MBsums [19] for transforming MB-representations for Feynman integrals into multiple sums. The licence conditions of the source-open package are those formulated in the CPC non-profit use licence agreement of the Computer Physics Communications Program Library [20]. The authors expect that the potential users read and follow the licence agreement when using this code.

2. The Mathematica package MBsums

The package MBsums transforms Mellin–Barnes integrals into sums, by closing the integration contours and calculating the integrals by the residue theorem, *i.e.* by constructing sums over all residues inside the contours. The current version of MBsums is 1.0. The package MBsums works with Wolfram Mathematica 7.0 and later.

In order to obtain a sum from an MB-integral, the user should use the MBIntToSum function of MBsums:

or

where int is the MB-integral in the form as it is denoted in the Mathematica package MB $[9]^1$:

which corresponds to

$$\operatorname{int} = \frac{1}{(2\pi i)^d} \int_{-i\infty+c_1}^{i\infty+c_1} \cdots \int_{-i\infty+c_D}^{i\infty+c_D} f \prod_{k=1}^D \mathrm{d} z_k \,. \tag{4}$$

The integrand f can have the form

$$f = \sum_{j} f_{j} \tag{5}$$

¹ The package MBsums uses notations of the package MB, but can be run also independently.

and each f_i is assumed to be of the form of

$$f_{j} = \xi_{j} \frac{\prod_{m} \Gamma(N_{m}^{(j,1)}) \prod_{m} \Psi^{(n)} \left(N_{m}^{(j,2)}\right)}{\prod_{m} \Gamma\left(N_{m}^{(j,3)}\right)} \prod_{m} r_{m}^{N_{m}^{(j,4)}},$$
(6)

where $\Psi^{(0)}(z) = d \ln(\Gamma(z))/dz$, $\Psi^{(n>0)}(z) = d^n \Psi^{(0)}(z)/dz^n$, r_i are free parameters (usually kinematic parameters) in int and ξ_j is a factor independent of z-variables. The

$$N_{m}^{(j,k)} = \sum_{i} \alpha_{m,i}^{(j,k)} z_{i} + \beta_{m}^{(j,k)} + \gamma_{m}^{(j,k)} \epsilon \,, \tag{7}$$

where $\alpha_{m,i}^{(j,k)}$, $\beta_m^{(j,k)}$, $\gamma_m^{(j,k)}$ are rational numbers and ϵ (eps) is an infinitesimal dimensional shift, *e.g.* arising from $d = 4 - 2\epsilon$. All the singularities of the integrand of the MB-integral **f** are due (and only due) to Gamma and PolyGamma functions.

The values of c1, c2, ..., cD are converted to rational numbers by MBIntToSum before calculations.

Let us now focus on the case when the list kinematics is empty, *i.e.* we will consider (2). The list contours has the form

$$contours = \{z1 \rightarrow L/R, z2 \rightarrow L/R, ..., zD \rightarrow L/R\}.$$
 (8)

The order of the z-variables defines the order of integrations chosen by the user (from left to right). The L (R) means that the contour will be closed to the left (right). The L/R choice made by the user can be changed if kinematics is not empty and this will be covered later. The output of MBIntToSum in (2) is of the form

where

represents a sum with summand Sum_Coefficient_i that has non-negative indices given in the list List_i, and Conditions_i are conditions on those indices. The complete answer is the sum of all MBsum_i in the list.

The list kinematics has the form

where r_i are free parameters (usually kinematic parameters) in int and v_i are values of r_i. If kinematics is not empty, then MBIntToSum will

try to change the L/R choice made by the user in the list contours in order to obtain sums that have good asymptotic behaviour at given values of r_i . In any case the user is informed how the contour was closed. This will be explained later in detail. The values of v_i are converted to rational numbers by MBIntToSum before calculations. The user can turn off all messages printed by MBIntToSum by typing MBsumsInfo=False and turn them on by typing MBsumsInfo=True.

In addition, we provide function DoAllMBSums[sums,nmax,kinematics] that sums the sums in the form of (9). The nmax is the maximal value of each index, the minimal value is given by conditions on indices. The list kinematics is as above and may be empty. We used Wolfram Mathematica function ParallelMap inside DoAllMBSums to sum individual sums in the list sums in parallel.

3. Obtaining the sums

Let us now shortly explain how we obtain the sums. We point out the most important ingredients in our algorithm. Let us focus on the case when the list kinematics is empty, *i.e.* we will consider (2). The MB-integral is in the form as it was denoted in (3). Let us now assume that the user has chosen as first integration the z_2 ->L. As a first step we form a list, which we call NegArgsDoC, of arguments of the Gamma and PolyGamma functions in the numerators that give residues for $\operatorname{Re}(z_2) < c_2$ (see (3), the remaining contours are seen as a straight lines). We call that list NegArgsDoC. Next, we consider all possible cases: When all Gamma and PolyGamma functions that have arguments in NegArgsDoC contribute to a residue at the same time, and when only some subset of them contributes to a residue at the same time. We consider all possible subsets of NegArgsDoC. Additionally, we have to be careful when some Gamma functions in the denominator become singular at some points. If we have terms like $Gamma[2 \ z2]$, then the poles are at $z_2 = -n/2$ and we consider there 2 cases: n = 2n' and n = 2n' + 1, where n and n' are non-negative integers. Similarly, we proceed with arbitrary $M \times z2$ terms, where M is some integer value and, in general, with all n/Mterms which appear together with integration variable in the arguments of the Gamma and PolyGamma functions. So we produce a list of cases

$$\left\{\left\{f_1^{(1)}, c_1^{(1)}\right\}, \left\{f_2^{(1)}, c_2^{(1)}\right\}, \dots, \left\{f_{K_1}^{(1)}, c_{K_1}^{(1)}\right\}\right\},$$
(12)

where $f_i^{(1)}$ are expressions after taking residues of f and $c_i^{(1)}$ are conditions on the index that numerates terms (residues) in $f_i^{(1)}$. We obtain a list of K_1 elements after integrating over z_2 . Let us now assume that the user has chosen as second integration variable z_5 ->R. Then, we repeat the whole procedure on each $f_i^{(1)}$ taking into account conditions $c_i^{(1)}$. Thus, we produce analogous to (12) a list of cases

$$\left\{\left\{f_1^{(2)}, c_1^{(2)}\right\}, \left\{f_2^{(2)}, c_2^{(2)}\right\}, \dots, \left\{f_{K_2}^{(2)}, c_{K_2}^{(2)}\right\}\right\}.$$
(13)

We repeat the whole procedure for each integration variable.

4. Contours and convergent sums

Let us now shortly explain how we obtain the sums if the list kinematics is not empty. We follow the order of integration given in the list contours. Our aim is to determine the L/R such that we obtain sums that have good asymptotic behaviour at given values of \mathbf{r}_i in the list kinematics. We do it in the following way. At each integration step s, we analyse the expressions $f_i^{(s)}$ that are to be integrated over some z_C . Each $f = f_i^{(s)}$ we decompose as

$$f = \sum_{j} g_j \tag{14}$$

and each $g = g_j$ is of the form of

$$g = r_1^{a_1} r_2^{a_2} \dots r_K^{a_K} F, \qquad a_j = \sum_i a_{j,i} z_i, \qquad (15)$$

where r_i are the kinematic parameters in the list kinematics and F contains the rest of g. If we integrate over z_C , we consider

$$c^{z_C}, \qquad c = r_1^{a_{1,C}} r_2^{a_{2,C}} \dots r_K^{a_{K,C}}.$$
 (16)

The value of c is calculated. MBIntToSum prints the error message:

Found c = c (not a number): please complete kinematic's list for each g_i in each $f_i^{(s)}$ when c is symbolic (not a number) and at the end MBIntToSum prints

Unable to find correct contour for
$$z_C$$

and returns {}. The user should complete the list kinematics.

For each g_i in each $f_i^{(s)}$ it is returned L if |c| > 1 or R if |c| < 1 indicating how to close the contour or $\{\}$ if |c| = 1.

If for each g_i in each $f_i^{(s)}$ it is returneded L (R) or {}, then the contour for z_C will be closed to left (right) if it is returned at least one L (R).

If for each g_i in each $f = f_i^{(s)}$ it is returned $\{\}$, then the choice of user given in the list contours is taken.

If for some g_i it is returned R and for some g_j it is returned L, then MBIntToSum prints the error message:

Unable to find correct contour for z_C

and returns {}. Otherwise, we compute the sums as described above. We repeat the whole procedure for each integration variable. We stress that this procedure as described above does not always give convergent sums.

There are MB-integrals for which no convergent sums can be found. One such example is the following MB-integral:

$$B_1 = \frac{1}{2\pi i} \int_{-i\infty}^{i\infty} dz \Gamma^2 (1+z) \Gamma^2 (1-z) \,. \tag{17}$$

Here we can apply the first Barnes lemma [3] and obtain $B_1 = 1/6$, but the reader can check that indeed the infinite series of residues diverge both for Re z > 0 and Re z < 0.

Consider the following integral (see also [3]):

$$B_x = \frac{1}{2\pi i} \int_{-i\infty}^{i\infty} dz x^z \Gamma^2(1+z) \Gamma^2(1-z) , \qquad x > 0 \land x \neq 1 .$$
 (18)

Closing the contour to the right ($\operatorname{Re} z > 0$) gives the following series

$$s_{\rm R} = -\sum_{n=1}^{\infty} n x^n (2 + n \ln(x)),$$
 (19)

convergent for 0 < x < 1, while closing the contour to the left (Re z < 0) gives the following series

$$s_{\rm L} = -\sum_{n=1}^{\infty} n x^{-n} (2 - n \ln(x)) , \qquad (20)$$

convergent for x > 1. Both $s_{\rm L}$ and $s_{\rm R}$ give the same formula after summing up, that is

$$s_{\rm LR} = \frac{x(2-2x+(1+x)\ln(x))}{(x-1)^3}, \qquad x > 0 \land x \neq 1, \tag{21}$$

so $B_x = s_{\text{LR}}$ and $\lim_{x \to 1} B_x = B_1$.

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