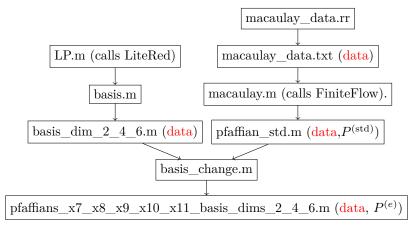
1 Program codes and data to obtain the Pfaffian matrices $P_i^{(e)}$ of one-loop massless pentagon with one massive leg



These program files are in mma_gkz/1L_Om_1os_pentagon of http://www.math. kobe-u.ac.jp/OpenXM/Math/amp-MM/pentagon-dist.zip. deq_check.m is a program to check the agreement of results by LiteRed and our method.

2 Notes

System configuration.

- 1. These codes are tested on Debian 11.1 with Mathematica 12.3.1, Risa/Asir, FiniteFlow, and LiteRed.
- 2. Risa/Asir download: http://www.openxm.org
- 3. FiniteFlow download: https://github.com/peraro/finiteflow
- 4. LiteRed download: https://www.inp.nsk.su/~lee/programs/LiteRed/
- The folder ~/.Mathematica/Applications contains LiteRed, LiteRed.m, RNL.
- 6. The file ~/.Mathematica/Kernel/init.m should be set properly to run FiniteFlow. A sample init.m is in the zip file pentagon-dist.zip
- 7. The command asir must be in the command search path and some environmental variables for asir should be set properly. Please use /usr/local/OpenXM/rc/dot.bashrc to set them.
- 8. gkz_utils/gkz.m of pentagon-dist.zip contains an interface (for Debian) to Risa/Asir and the package mt_gkz.m, tmp-MM.rr and linsolv. The interface requires that ~/.asirrc (asir initial run commands) is put

under the home folder. It is in the pentagon-dist.zip with the name dot.asirrc.

1L_Om_1os_pentagon/

Typewriter font is used for variables in codes and math-italic font is used for variables in the paper in the sequel. The page number of the paper is of the version 1 of the arxiv preprint.

- 1. macaulay.m: it constructs Pfaffian matrices $P_i^{(\text{std})}$ by the Macaulay matrix with FFSparseSolve function of FiniteFlow. $P_7^{(\text{std})}$ is stored in p[1], ..., $P_{11}^{(\text{std})}$ is stored in p[5].
- 2. basis_change.m: p7, p8, p9, p10, p11 are Pfaffian matrices $P_i^{(e)}$, $i = 7, \ldots, 11$ for the basis $\{e_i\}^1$. g and e give a Gauge transformation from $P_i^{(\text{std})}$ to $P_i^{(e)}$. The diagonal element Λ_{ii}^2 is proportional to e[[i,i]] in the code. Precisely speaking, e[[i,j]] is equal to Λ_{ii}/Λ'_i times prefactors³
- 3. basis.m: The "physical basis" ints is obtained by the IBPs section of LP.m, which is commented out in the code. mis is the output MIs[pent]. Each element of ints is of the form j[d0,nu], which is a generalized Feynman integral. The function diffop[] is defined in

gkz_utils/mma_asir_interface.m. It constructs a differential operator corresponding to a given differential form expressed by j[d0,nu] by calling mt_gkz.rr package. dbaseByD, which is equal to $e'_i{}^{\mathcal{D}}{}^4$ is an expression of dbase by differential operators.

gkz_util/

1. math.m: The function LPfac[] gives $c(d_0^{(i)}, \nu^{(i)})^5$ and the function kinfac[] gives $(-s_{12})^{d_0^{(i)}/2-\epsilon-|\nu^{(i)}|-5\epsilon\delta}$ 6

¹page 26, (5.36)

 $^{^{2}}$ page 26, (5.37)

³We call $(-s_{12})^{\epsilon}$,... in (5.36) of page 26 prefactors. The variable s12 in the program codes stands for $-s_{12}$ in the paper.

 $^{^{4}}$ page 22, (5.4) and page 27, (5.39)

 $^{{}^{5}}$ page 26, (5.37)

 $^{^{6}}$ page 26, (5.37)