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Next to Leading Order Calculations for Higgs Boson + Jets

Simon Armstrong

A Thesis presented for the degree of
Doctor of Philosophy



Institute of Particle Physics Phenomenology
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August 22, 2017

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Abstract

Due to the recent Higgs boson discovery, an important target for particle physics is to investigate its properties to determine if it is the standard model Higgs boson or some other variety. The Large Hadron Collider is now in Run Two, collecting even more data at higher precisions, which requires predictions at next to leading order or higher orders. Therefore it is important to have an efficient and automatic calculation of the next to leading order amplitudes for the Higgs boson. This thesis discusses the methods needed to perform these calculations.

These calculations are specifically developed in order to add them to the BlackHat library, which already provides these types of calculations for amplitudes involving quarks, gluons and W and Z bosons. Both this thesis and BlackHat use recursive methods, as these are more efficient than using the Feynman rules directly. Specifically the BCFW recursion relation is used to calculate tree amplitudes and generalised unitarity is used to calculate one loop amplitudes. These methods are first used in 4 dimensions to calculate the cut constructable parts of the amplitudes, then the extension of these techniques to higher numbers of dimensions is discussed, allowing the rational terms to be extracted using D dimensional generalised unitarity. In general, two different even integer dimensions higher than 4 are required for a numeric implementation of D dimensional generalised unitarity with spinors, which therefore requires working in both 6 and 8 dimensions. To enable an efficient implementation, a technique is introduced that allows only 6 dimensional calculations to be used rather than both 6 and 8 dimensional calculations and a reduction of 6 dimensional calculations to be in terms of only 4 dimensional objects is developed. The methods presented in this thesis provide a solid groundwork for Higgs boson amplitudes to be implemented into BlackHat.

Declaration

The work in this thesis is based on research carried out at the Institute of Particle Physics Phenomenology, the Department of Physics, University of Durham, England. No part of this thesis has been submitted elsewhere for any other degree or qualification and it is all my own work unless referenced to the contrary in the text.

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Chapter 1

Introduction and Motivation

It is an exciting time for particle physics, with the Large Hadron Collider (LHC) now collecting data for Run Two at higher than ever energies. At these high energies there is much hope that some new physics, beyond the standard model, will be detected which will help with understanding the problems with the standard model, such as the lack of candidates for dark matter and the hierarchy problem. Supersymmetry, in which each standard model has a heavier partner particle, is one example of a beyond the standard model theory that could be detected. These types of theory are one of the main targets for Run Two.

Another of the main targets for Run Two is characterisation of the recently detected Higgs boson. This would show whether it is the standard model Higgs boson or some more exotic version. There is also hope that the Higgs boson could provide hints of new, beyond the standard model, theories through its interactions with them. Both high precision measurements and high precision predictions are needed to enable accurate conclusions to be drawn. The LHC in its second run is starting to collect high precision data for the relevant channels and analyses, so the situation is being approached where the limitation is the precision of the model predictions. Many of the channels at the LHC produce large numbers of jets. To produce predictions a hard process is combined with parton shower and hadronisation algorithms in a Monte Carlo simulation, which converts quarks and gluons produced in the hard process into large showers of particles, through soft and collinear radiation and then into the observable hadrons which can be detected as part of jets. A group of hadronised particles, produced by soft and collinear radiation, travelling in roughly the same direction and from the same location, will be detected as a jet. Each of these

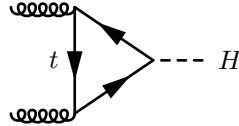


Figure 1.1: The leading contribution to the Higgs-gluon coupling which is mediated by a top quark loop.

jets is typically due to the radiation of a single particle produced in the hard process. Therefore, the leading contribution to processes with many jets will be from hard processes that produce a high multiplicity of particles. It is also important that the calculations are in a form that is efficient for use in a Monte Carlo simulation, so that they can be combined with the other steps to produce physically observable predictions.

The Higgs boson coupling is proportional to the mass of the particle involved and as such in quantum chromodynamics the strongest interaction is with the top and bottom quarks. Top and bottom quark masses are larger than the energy scale of the LHC and so can be treated as approaching infinite mass. All other quark masses are much lower than this scale and so can be approximated as massless. The exact leading order calculation for this type of amplitude is already a one loop amplitude, as the top quark runs in a loop. In this heavy top quark limit, as the mass is a large scale, the leading term in terms of the top quark mass can be approximated out and all other terms can be neglected. This only leaves the contribution from a top quark loop connecting two gluons and one Higgs boson, of the form in [Figure 1.1](#), from which the loop quark's degrees of freedom can be integrated out to give a new effective vertex between a Higgs boson and two gluons. The effective Lagrangian term this introduces is given by

$$\mathcal{L}_H^{\text{int}} = \frac{C}{2} H \text{tr} G_{\mu\nu} G^{\mu\nu} , \quad (1.1)$$

where H is the Higgs field, G is the gluon field strength tensor and C is the dimensionful coupling constant which can be calculated order by order in α_s . As this limit is being used, all particles other than the Higgs boson will be assumed to be massless from here on. Currently, amplitudes with a Higgs boson and jets at tree level in the high top mass limit are automated for any multiplicity.

As the Next to Leading Order (NLO) correction and using the true masses of the top and bottom quarks could be large contributions, both are needed to increase the

precision of our predictions. The accuracy of working in the infinite top mass limit at NLO has been directly investigated at low multiplicity by Harlander et al.[1] and Grazzini et al.[2] and has been found to give a relatively flat correction of the order of a few % for large areas of the phase space. This thesis examines the techniques and methods needed to calculate the NLO amplitude for a Higgs boson with many jets in the high top mass limit. The case of NLO without taking the high top mass limit is a two loop calculation and as such is beyond our current ability to compute in an efficient automated way for high multiplicities.

These NLO amplitudes are divergent in 4 dimensions and as such must be regulated. One of the most common methods, and the one used in this thesis, is to working in $d = 4 - 2\epsilon$ dimensions which causes the amplitude to have poles in ϵ which control and contain the divergences and will be cancelled with the soft and collinear contributions to calculate the finite cross section. There are many ways to perform this calculation. One scheme often used is the 't Hooft-Veltman scheme which allows all elements of the loop to extend into to the extra dimensions. Another scheme is the four dimensional helicity scheme[3] which keeps all spinor and polarisation vector states in 4 dimensions and allows only the internal loop momenta to enter the extra dimensions. This scheme has the advantage that the Ward identities are preserved which allows checks and tools to relate different amplitudes. This scheme is the scheme used in BlackHat and which will be used in this project. The different schemes are often related in ways that don't require extra loop calculations for example the 't Hooft-Veltman and four dimensional helicity scheme amplitudes for a pure gluon amplitude differ by a factor of

$$A^{HV} = A^{FDH} - \frac{\Gamma(1 + \epsilon)\Gamma(1 - \epsilon)}{3(4\pi)^{2-\epsilon}} \mu^{2\epsilon} A^{\text{tree}} \quad (1.2)$$

where A^{HV} is the amplitude in the 't Hooft-Veltman scheme, A^{FDH} is the amplitude in the four dimensional helicity scheme, A^{tree} is the tree amplitude and μ is a renormalisation scheme used to preserve the dimension of the amplitude.

It has already been shown that the amplitudes for any number of gluons in the Maximally Helicity Violating configuration are the very simple Parke-Taylor amplitudes[4]. In the Feynman diagram method, even for four gluons, the calculation contains three different diagrams. When the number of particles in an amplitude is increased, the number of terms increases factorially. This greatly limits the number

of particles that can be included before the calculation of the amplitude becomes too large to effectively implement. At 1-loop these problems become worse and even simple four particle amplitudes can become difficult to perform. Therefore, the Feynman diagram method and explicit loop integrations are not efficient for these types of amplitudes and more efficient methods are needed.

The BlackHat library[5, 6, 7] already includes calculations for amplitudes with quarks, gluons and optionally one of the vector bosons, W^\pm or Z , and other processes but does not currently include amplitudes for the Higgs boson. This thesis describes the methods needed to efficiently implement the calculation of the NLO 1-loop amplitude for Higgs boson with jets, in a generic way, for any number of quarks and gluons of any helicities. Once the calculations have been implemented into the BlackHat library it will be a very useful tool for adding NLO calculations into existing Monte Carlo Event Generators such as Herwig++ or Sherpa.

Following this introduction, [Chapter 2](#) gives a short review of the techniques of colour ordered amplitudes and the spinor helicity formalism. These form the basis of the techniques used in the remaining chapters. The notations used in this thesis are also introduced in this chapter. [Chapter 3](#) explores the BCFW recursion relation which is used to compute tree level amplitudes and discusses how it extends to amplitudes with a Higgs boson.

The next two chapters discuss the techniques used to calculate the loop amplitude. Firstly, [Chapter 4](#) demonstrates how to calculate the cut constructable parts using Generalised Unitarity. Explicit numerical formulas are derived to enable a systematic numeric calculation to be implemented and this is extended to support amplitudes with a Higgs boson. [Chapter 5](#) explores how to calculate the rational terms of amplitudes including a Higgs boson, using 6 dimensional spinors. After deriving a 6 dimensional extension of the spinor helicity formalism, tree amplitudes are calculated using it and simplified to allow an efficient calculation of the contributions needed for 6 dimensional Generalised Unitarity. [Chapter 6](#) discusses how the calculation has been implemented, how to use the implementations and where possible, how BlackHat has been tested against the implementations developed in this project. Finally, [Chapter 7](#) concludes the project and discusses the remaining steps needed to fully implement these calculations into BlackHat.

Chapter 2

An Introduction to Efficient Calculation Techniques

There are several techniques that can be used to efficiently calculate amplitudes which separate the amplitudes into simpler parts. The first method used is colour ordered amplitudes which separates the amplitudes' kinematics from the colour factors and splits the amplitudes into simpler “colour ordered” amplitudes. The second is to work in the spinor helicity formalism which separates the different helicity states and produces simpler formulas for different cases within the amplitude. Both of these techniques are discussed based on the versions used by Dixon[8]. A decomposition of the Higgs boson into a complex scalar is also introduced here which separates amplitudes into MHV like amplitudes. Finally, recursive calculations using unitarity techniques are introduced, which form the bedrock for the techniques discussed in the following two chapters.

2.1 Colour Ordered Amplitudes

Inclusion of colour factors in numerical calculations can greatly increase their complexity. If colour factors are included an amplitude is increased from a single complex number to a large tensor of complex numbers in colour space. Many of the elements in the matrix are zero and many of the rest are related, so it should be possible to extract and then combine elements to give a simpler form which has less redundancy and therefore leads to a more efficient calculation. One way this recombination can

be done is by expanding an amplitude as the coefficients of different colour terms. Each of these coefficients would be a single complex number. The modulus squared of any specific type of amplitude, as used in the cross section, can be calculated from the coefficients, the number of colours and the number of quarks.

To reduce an amplitude to the smallest possible set of colour terms requires converting all colour factors to chains of the fundamental generators with no internal colour or gluon indices. This reduction is possible as the gluon colour factor, f^{abc} , which is the structure constant for the fundamental generators, can be written as[8]

$$f^{abc} = -\frac{i}{\sqrt{2}} (\text{tr}[T^a T^b T^c] - \text{tr}[T^a T^c T^b]) , \quad (2.1)$$

where T^a is the fundamental generator which is normalised as[8]

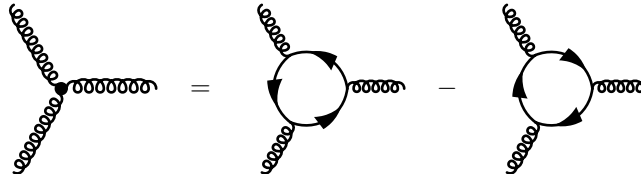
$$\text{tr}[T^a T^b] = \delta^{ab} . \quad (2.2)$$

To remove internal gluon indices the Fierz identity for the fundamental generators,[8]

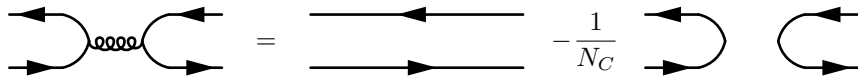
$$T^a_{i\bar{j}} T^a_{\bar{i}j} = \delta_{i\bar{j}} \delta_{\bar{i}j} - \frac{1}{N_c} \delta_i^j \delta_{\bar{i}\bar{j}} , \quad (2.3)$$

is used. These relations can be used to convert the colour factor for any pure gluon amplitude into a sum of terms that are each a single trace of fundamental generators, one generator for each gluon. For amplitudes with quarks, the colour factor can be converted into a sum of terms that are each the product of a single (possibly empty) trace of fundamental generators and a chain of fundamental generators for each quark pair, where again each gluon's generator only appears once in each term.

The relations given in Equations 2.1 and 2.3 can be represented graphically in terms of replacements of quarks and gluons as[8]



$$\text{Three-gluon vertex} = \text{Circular diagram 1} - \text{Circular diagram 2} \quad (2.4)$$



$$\text{Quark-gluon loop} = \text{Straight quark line} - \frac{1}{N_c} \text{Quark loop} \quad (2.5)$$

where the rules only apply for the colour factors not kinematics.

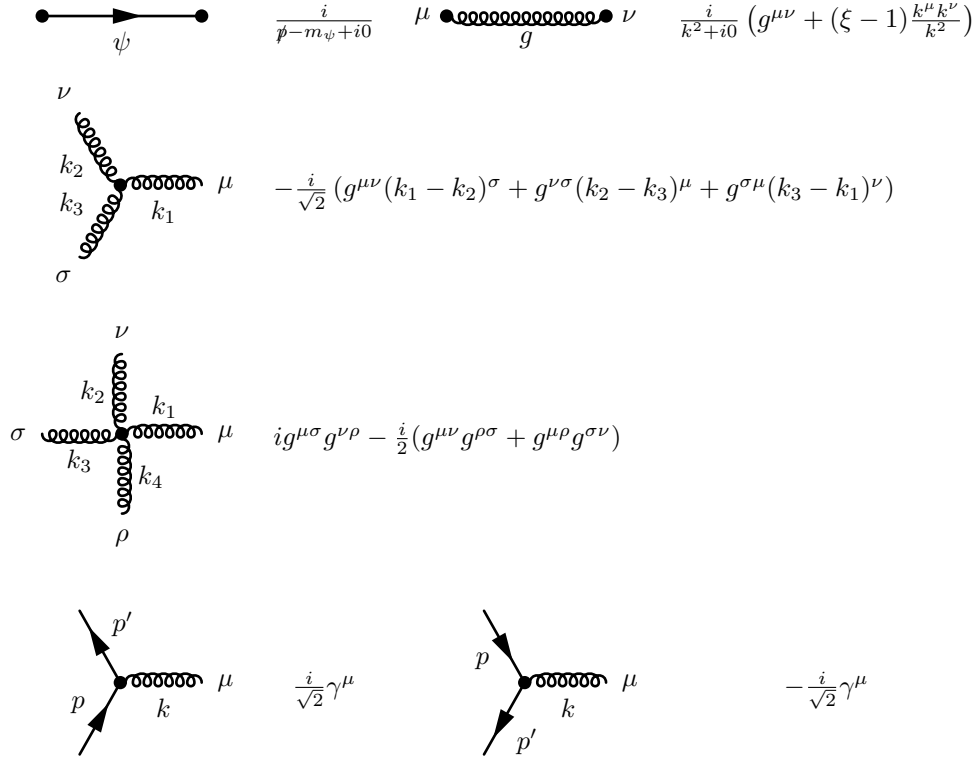


Table 2.1: Colour ordered Feynman rules in Faddeev-Popov gauge[8]. All momenta are inbound.

representation of the rearrangement is shown in Figure 2.1.

As all diagrams can be replaced with terms of the form shown in the last line of Equation 2.6, the entire tree amplitude can be changed to terms of that form. The coefficients of each trace structure can be extracted, which gives the colour ordered partial amplitudes. By combining them, the full amplitude can be constructed from the colour ordered partial amplitudes as

$$\mathcal{A}_{5, \text{gluon}, \text{tree}} = g^3 \sum_{\sigma \in S_5/Z_5} \text{tr}[T^{a_{\sigma(1)}} T^{a_{\sigma(2)}} T^{a_{\sigma(3)}} T^{a_{\sigma(4)}} T^{a_{\sigma(5)}}] A_{5, \text{gluon}, \text{tree}}(\sigma(1), \sigma(2), \sigma(3), \sigma(4), \sigma(5)). \quad (2.7)$$

Rather than calculating the full amplitude directly and then extracting the coefficient for each colour structure, it is possible to extract the coefficients for the colour structures from the Feynman rules and then evaluate the colour ordered partial amplitudes. The modified Feynman rules, given in Table 2.1, enable the colour ordered amplitudes to be evaluated by summing the expressions for all possible planar diagrams with a given fixed order of the external particles. There are two vertices for

$q\bar{q}g$ which differ in the direction of the quark line and are no longer equivalent, as the graphs are planar and the external particles have a fixed order.

For any number of gluons the formula extends to[8]

$$\mathcal{A}_{n \text{ gluon}, tree} = g^{n-2} \sum_{\sigma \in S_n/Z_n} \text{tr}[T^{a_{\sigma(1)}} \dots T^{a_{\sigma(n)}}] A_{n \text{ gluon}, tree}(\sigma(1), \dots, \sigma(n)) , \quad (2.8)$$

where S_n/Z_n is the set of permutations of n items that are not equivalent under cycles and $A_{n \text{ gluon}, tree}$ is a colour ordered amplitude. If there are external fermions then there will be chains of fundamental generators which connect with the quark colour indices. For example, with one pair of external fermions the amplitude takes the form[8]

$$\begin{aligned} \mathcal{A}_{n-2 \text{ gluon}, 2 \text{ quark}, tree} = \\ g^{n-2} \sum_{\sigma \in S_{n-2}} [T^{a_{\sigma(3)}} \dots T^{a_{\sigma(n)}}]_i^j A_{n-2 \text{ gluon}, 2 \text{ quarks}, tree}(1_p, 2_{\bar{p}}, \sigma(3), \dots, \sigma(n)) . \end{aligned} \quad (2.9)$$

The same method applies for 1 loop calculations where, for example, the n gluon amplitude is given by[8]

$$\begin{aligned} \mathcal{A}_{n \text{ gluon}, 1-loop} = g^n \left(\sum_{\sigma \in S_n/Z_n} N_c \text{tr}[T^{a_{\sigma(1)}} \dots T^{a_{\sigma(n)}}] A_{n;1}(\sigma(1), \dots, \sigma(n)) \right. \\ \left. \sum_{c=2}^{\lfloor \frac{n}{2} \rfloor + 1} \sum_{\sigma \in S_n/S_{n;c}} \text{tr}[T^{a_{\sigma(1)}} \dots T^{a_{\sigma(c-1)}}] \text{tr}[T^{a_{\sigma(c)}} \dots T^{a_{\sigma(n)}}] A_{n;c}(\sigma(1), \dots, \sigma(n)) \right) , \end{aligned} \quad (2.10)$$

where $S_n/S_{n;c}$ is the set of all permutations of n items that do not give the same trace terms, $\lfloor n \rfloor$ is the largest integer less than or equal to n and the $A_{n;c}$ are partial amplitudes. The $A_{n;1}$ are colour ordered partial amplitudes and are called primary amplitudes. The other partial amplitudes are not colour ordered. If there are only external gluons, these other partial amplitudes can be written as a sum of permutations of $A_{n;1}$. If there are external quarks, the $A_{n;c>1}$ can no longer be written in terms of only the $A_{n;1}$, but will need new colour ordered partial amplitudes which are defined by specifying the direction the quark travels around the loop. Again this decomposition can be performed diagrammatically and is shown for a two gluon and two fermion one loop diagram in Fig. 2.2.

The Higgs boson is a completely colourless particle, as are the W and Z bosons

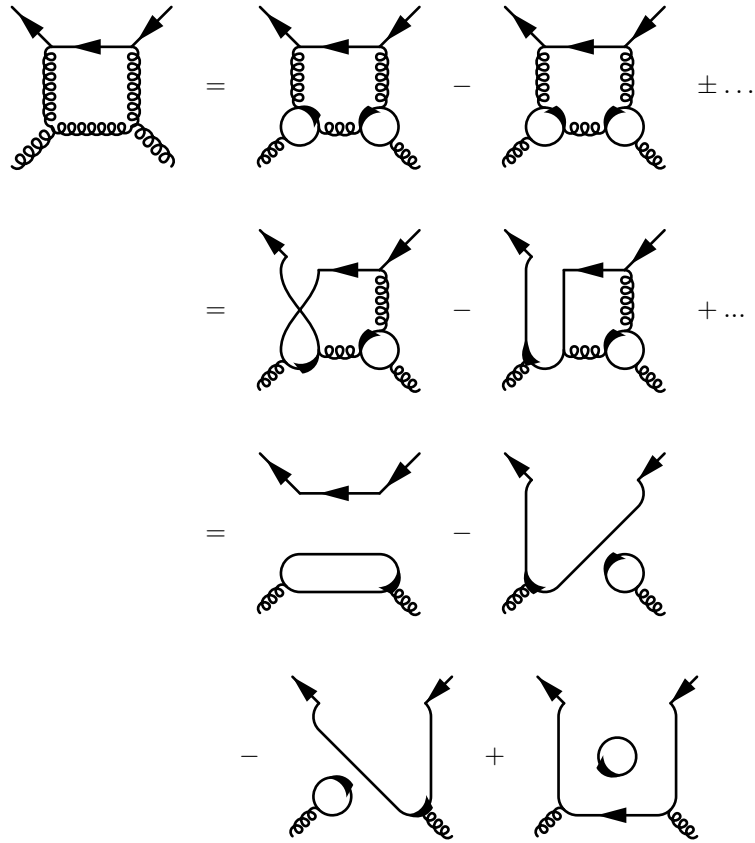


Figure 2.2: The colour decomposition of a one loop, two fermion and two gluon diagram. All $\frac{1}{N_c}$ contributions cancel at all stages.

which are already implemented into BlackHat. These particles do not have any colour factors attached and as such are not involved in any colour rearrangements. The Higgs boson can therefore be introduced into amplitudes without changing the colour ordering techniques and its ordering in partial amplitudes is not fixed, meaning that any Higgs boson must be included in all possible positions. The Higgs boson can therefore be included into colour ordered amplitudes without any extra complications.

2.2 Spinor Helicity Formalism

Spinors are defined as the solution of the equation

$$\not{p}u_p = 0 , \quad (2.11)$$

where $\not{p} = \gamma_\mu p^\mu$ and γ_μ are the gamma matrices. The solutions are normalised to

$$u \otimes \bar{u} = \not{p} , \quad (2.12)$$

where \bar{u} is the conjugate spinor for u and \otimes is an inner product summing over the different spinor states. The conjugate spinors are solutions to

$$\bar{u}_p \not{p} = 0 , \quad (2.13)$$

and for real momenta are related to the normal spinor by

$$\bar{u}_p = u_p^\dagger \gamma_0 . \quad (2.14)$$

In order to avoid problems with extending the definition of conjugate spinors to complex momenta, conjugate spinors defined without needing conjugation will be used for this project.

For 4 dimensions there is one degree of freedom left over after the spinors are normalised so there are two basis states in the spinor space. It was discovered that for massless QCD amplitudes are very simple when the basis states are helicity eigenstates. For massless spinors the helicity operator is defined by

$$\gamma^5 = i\gamma^0\gamma^1\gamma^2\gamma^3 . \quad (2.15)$$

The spinor eigen states are defined as

$$u_\pm(k) = P^\pm u(k) = \frac{1 \pm \gamma_5}{2} u(k) \quad v_\mp(k) = P^\pm v(k) = \frac{1 \pm \gamma_5}{2} v(k) , \quad (2.16)$$

where P^\pm is the projection operator for γ_5 and is defined as

$$P^\pm = \frac{1 \pm \gamma_5}{2} . \quad (2.17)$$

From the completeness relation for massless spinors

$$u(k) \otimes \bar{u}(k) = v(k) \otimes \bar{v}(k) = \not{k} , \quad (2.18)$$

it can be seen that the spinors for v and u are proportional to each other and as an overall phase is arbitrary the spinors can be defined such that $u_{\pm}(k) = v_{\mp}(k)$.

A more compact bra-ket notation can be defined which is given by[8]

$$\begin{aligned} u^+(k_i) = v^-(k_i) &\equiv |k_i^+\rangle \equiv |i\rangle & u^-(k_i) = v^+(k_i) &\equiv |k_i^-\rangle \equiv |i] \\ \bar{u}^+(k_i) = \bar{v}^-(k_i) &\equiv \langle k_i^+| \equiv [i| & \bar{u}^-(k_i) = \bar{v}^+(k_i) &\equiv \langle k_i^-| \equiv \langle i| , \end{aligned} \quad (2.19)$$

where the sign of the bra-ket type spinor is defined by the eigen value under the helicity operator. Some combinations of these spinors have products that are zero due to being opposite helicity eigen states, these products are $[1|2\rangle$ and $\langle 1|2]$. As opposite sign projection operators annihilate and the two spinors in the product are projected with opposite sign projection operators, these products vanish. As with all massless spinors, a product of any two of these spinors for the same momentum vanishes. The remaining spinor products can be combined to form the relation[8]

$$\langle ij \rangle [ji] = \text{Tr} \left[\frac{1 - \gamma^5}{2} \not{k}_i \not{k}_j \right] = 2k_i \cdot k_j = (k_i + k_j)^2 = s_{ij} , \quad (2.20)$$

which is derived using the completion relation for spinors, the anti-commutation relations for gamma matrices and, as the momenta are massless, $k_i^2 = k_j^2 = 0$. These spinor products are anti-symmetric so obey $\langle ij \rangle = -\langle ji \rangle$ and $[ij] = -[ji]$.

Gluon polarisation vectors can be written in terms of spinors as[8]

$$\epsilon_{\mu}^{\pm}(k; q) = \pm \frac{\langle k^{\pm} | \gamma_{\mu} | q^{\pm} \rangle}{\sqrt{2} \langle k^{\mp} q^{\pm} \rangle} , \quad (2.21)$$

or explicitly for each state

$$\epsilon_{\mu}^{+}(k; q) = \frac{[k | \gamma_{\mu} | q]}{\sqrt{2} \langle k q \rangle} , \quad \epsilon_{\mu}^{-}(k; q) = \frac{\langle k | \gamma_{\mu} | q]}{\sqrt{2} [q k]} , \quad (2.22)$$

where the vector, q , can be arbitrarily chosen independently for each gluon but must not be changed between uses of the same gluon and corresponds to a gauge choice. It can be shown to be a gauge by checking that a change in $q \rightarrow q'$ is proportional to

the vector k ,

$$\begin{aligned}
\epsilon_\mu^+(k; q') - \epsilon_\mu^+(k; q) &= \frac{[k|\gamma_\mu|q']}{\sqrt{2}\langle kq' \rangle} - \frac{[k|\gamma_\mu|q]}{\sqrt{2}\langle kq \rangle} \\
&= \frac{(-\langle qk \rangle)[k|\gamma_\mu|q'] - (-\langle q'k \rangle)[k|\gamma_\mu|q]}{2\langle kq' \rangle\langle kq \rangle} \\
&= \frac{-\langle q|k|\gamma_\mu|q' \rangle + \langle q'|k|\gamma_\mu|q \rangle}{2\langle kq' \rangle\langle kq \rangle} \\
&= \frac{\langle q'|\gamma_\mu|k|q \rangle + \langle q'|k|\gamma_\mu|q \rangle}{2\langle kq' \rangle\langle kq \rangle} \\
&= \frac{\langle q'(\gamma_\mu k + k\gamma_\mu)q \rangle}{2\langle kq' \rangle\langle kq \rangle} \\
&= \frac{\langle q'q \rangle}{2\langle kq' \rangle\langle kq \rangle} 2k^\mu \\
&= \frac{\langle q'q \rangle}{\langle kq' \rangle\langle kq \rangle} k^\mu, \tag{2.23}
\end{aligned}$$

where in the first few lines the properties of spinor products have been used, in the later lines the anti-commutation relations for slashed matrices have been used and the same method applies for the transformation of the other helicity polarisation vector. It is also possible to check that these expressions obey all the normal relations for polarisation vectors such as

$$\epsilon_\mu^\pm(k; q)k^\mu = \pm \frac{\langle k^\pm|k|q^\pm \rangle}{\sqrt{2}\langle k^\mp q^\pm \rangle} = 0, \tag{2.24}$$

and

$$\begin{aligned}
\epsilon_\mu^+(k; q)\epsilon^{-\mu}(k; q) &= \frac{[k|\gamma_\mu|q]\langle k|\gamma^\mu|q]}{\sqrt{2}\langle kq \rangle\sqrt{2}[qk]} \\
&= \frac{2[kq]\langle kq \rangle}{2\langle kq \rangle[qk]} \\
&= -1, \tag{2.25}
\end{aligned}$$

where identities from [8] have been used, which are not shown or proved here. Due to the ability to choose the reference vector freely for each gluon and the many spinor products that are 0, it is often possible with a clever choice of reference vectors to vastly simplify the calculation of amplitudes and to cause many of the diagrams contributing to an amplitude to vanish.

Due to the simple relations between spinor products, many of which are 0, very simple expressions for the amplitudes can be found and as shown by Parke and Taylor the pure gluon amplitudes with all helicities the same or all but one the same are

0. They found the first non-zero amplitude to be the so called Maximally Helicity Violating amplitude which has all but two gluons having the same helicity. The mostly positive MHV amplitude for any number of gluons is given by[8]

$$A_n^{\text{tree}}_{\text{gluon}}(1^-, 2^+, \dots, i-1^+, i^-, i+1^+, \dots, n^+) = i \frac{\langle 1i \rangle^4}{\langle 12 \rangle \langle 23 \rangle \dots \langle n1 \rangle}, \quad (2.26)$$

where one of the negative helicity gluons can be chosen to be labelled as 1 due to cyclic symmetry of colour ordered amplitudes. This simple form is the same for any number of gluons as long as they are in the MHV form, in contrast with amplitudes in other systems where the amplitude increases enormously in complexity as the number of particles increases. There is also an equivalent formula for the mostly negative case.

The amplitudes with one quark line are also found to be very simple in the MHV cases, which for these amplitudes is when all but one gluon have the same helicity. The mostly positive amplitude for this type of amplitude is given by[8]

$$A_{n-2}^{\text{tree}}_{\text{gluon}, q\bar{q}}(1^-, \dots, i_q^+, \dots, j_{\bar{q}}^-, \dots, n^+) = i \frac{\langle 1i \rangle^3 \langle 1j \rangle}{\langle 12 \rangle \langle 23 \rangle \dots \langle n1 \rangle}, \quad (2.27)$$

where again the choice has been made to label the negative helicity gluon as 1. The quark anti-quark pair must have opposite helicity as QCD conserves helicity along quarks lines and all momenta and helicities are taken as for outgoing particles.

For other cases the amplitudes are often not as simple, but they are still significantly simpler than in other forms and many cases have a value of zero.

2.3 Higgs Boson as part of a Complex Scalar Field

A further simplification of calculations for amplitudes involving a Higgs boson can be performed by defining a complex scalar field[9]

$$\phi = \frac{1}{2} (H + iA), \quad (2.28)$$

where A is a real pseudo-scalar field which interacts with an effective vertex of

$$\mathcal{L}_A^{\text{int}} = \frac{C}{2} iA \text{tr} G_{\mu\nu}^* G^{\mu\nu}, \quad (2.29)$$

where $*G^{\mu\nu}$ is the dual of the gluon field strength tensor and is given by

$$*G^{\mu\nu} = \frac{i}{2} \epsilon^{\mu\nu\sigma\rho} G_{\sigma\rho} . \quad (2.30)$$

The interaction terms for the pseudo-scalar, A , and for the Higgs boson, H , combine to produce an interaction term for the complex field ϕ of[9]

$$\mathcal{L}_\phi^{\text{int}} = C \left(\phi \text{tr} G_{\text{SD}\mu\nu} G_{\text{SD}}^{\mu\nu} + \phi^\dagger \text{tr} G_{\text{ASD}\mu\nu} G_{\text{ASD}}^{\mu\nu} \right) , \quad (2.31)$$

where G_{SD} and G_{ASD} are the self dual and anti self dual gluon field strength tensors and are given by[9]

$$G_{\text{SD}}^{\mu\nu} = \frac{1}{2} (G^{\mu\nu} + *G^{\mu\nu}) \quad G_{\text{ASD}}^{\mu\nu} = \frac{1}{2} (G^{\mu\nu} - *G^{\mu\nu}) . \quad (2.32)$$

Working with the complex field ϕ rather than directly with the Higgs field H is particularly useful when working with helicity amplitudes and the spinor helicity formalism as the helicity structures do not mix and the amplitudes are closely related to their pure quark and gluon equivalents. For example the mostly positive MHV amplitude with a ϕ is given by[9]

$$A_{\phi,n}^{\text{tree gluon}}(\phi, 1^-, 2^+, \dots, i-1^+, i^-, i+1^+, \dots, n^+) = i \frac{\langle 1i \rangle^4}{\langle 12 \rangle \langle 23 \rangle \dots \langle n1 \rangle} , \quad (2.33)$$

where the form of this amplitude is identical to the form of the amplitude without the ϕ , which is given in [Equation 2.26](#). The presence of the ϕ particle does impact the mostly negative amplitudes, as it causes them not to vanish. The all negative amplitude is given by[9]

$$A_{\phi,n}^{\text{tree gluon}}(\phi, 1^-, 2^-, \dots, n^-) = (-1)^n i \frac{m_H^4}{[12][23] \dots [n1]} , \quad (2.34)$$

where m_H is the mass of the Higgs boson.

Amplitudes involving the conjugate field ϕ^\dagger do not need to be calculated, as amplitudes involving it are related to amplitudes involving ϕ by a parity transformation, which gives the relation[9]

$$A_n(\phi^\dagger, 1^{-h}, \dots, n^{-l}) = (-1)^{n_{q\bar{q}}} A_n(\phi, 1^h, \dots, n^l) \Big|_{(ij) \leftrightarrow [ji]} , \quad (2.35)$$

where $n_{q\bar{q}}$ is the number of $q\bar{q}$ pairs in the amplitude and $\langle ij \rangle \leftrightarrow [ji]$ corresponds to complex conjugation of the amplitude if all momenta are real. Therefore when working in 4 dimensions, as is being done in Chapters 3 and 4, all Higgs boson amplitudes will be calculated as amplitudes in terms of ϕ and will then be combined afterwards.

2.4 Recursive Construction of Amplitudes

With colour ordered amplitudes and the spinor helicity formalism, amplitudes can often be written in simple forms but for an automated system a method is required to build amplitudes, systematically, from simpler building blocks. The traditional method is the Feynman diagram method where amplitudes are built by summing Feynman rules over different diagrams. This method includes all possible information about a process, including handling off-shell particles in a fully general way, but for producing amplitudes this information is unnecessary. The information about off-shell particles, along with the gauge redundancy that is in the Feynman rules but not in any final amplitude, increases the complexity of this method. An improvement would be to build amplitudes from lower multiplicity, already on-shell and numeric amplitudes and this is exactly what is done in the Blackhat library and in this project. The BCFW recursion relation is used to build tree amplitudes from lower multiplicity tree amplitudes and then Generalised Unitarity is used to build loop amplitudes from tree amplitudes. The next few chapters will explain and derive these methods and extend for the Higgs boson.

Both the BCFW recursion relation and Generalised Unitarity methods rely on unitarity. This is the statement that the residue of poles in amplitudes are the product of two lower multiplicity amplitudes up to a sum over internal particles. To derive this, conservation of probability needs to be applied to the S matrix. The S matrix is the matrix that contains information about the probability of an interaction and is defined in terms of the amplitude matrix as

$$S = 1 + iA , \tag{2.36}$$

where A is the matrix of amplitudes. If probability is conserved then $SS^\dagger = 1$.

Substituting [Equation 2.36](#) into this condition results in

$$\begin{aligned} 1 &= (1 + iA)(1 - iA^\dagger) \\ &= 1 + i(A - A^\dagger) - AA^\dagger \\ AA^\dagger &= i(A - A^\dagger) , \end{aligned} \tag{2.37}$$

which tells us that the imaginary part of an amplitude is related to the product of lower multiplicity amplitudes. From the definition of amplitudes it can be seen that the only imaginary part will come from the $i\epsilon$ terms in propagators and will only contribute when the propagator goes on-shell. From this it is clear that the residue of an amplitude at a pole is proportional to a product of tree amplitudes.

Chapter 3

BCFW Recursion Relation

In 2005, Britto, Cachazo, Feng and Witten[10] discovered a recursion relation that relates an on-shell amplitude to lower multiplicity, on-shell amplitudes, which is called the BCFW recursion relation. This recursion relation can be applied to any amplitude, from any theory, but is particularly useful for calculations using the Spinor Helicity Formalism, where the amplitudes being calculated have a helicity structure, for reasons that will become apparent through this chapter. The BCFW recursion relation works by selecting two momenta from the external momenta in the amplitude and shifting them by an arbitrary complex amount, z , of a fixed vector n^μ , such that the two shifted momenta stay on-shell and that momentum conservation is preserved. The shift is given by

$$p_i^\mu \rightarrow \hat{p}_i^\mu(z) = p_i^\mu + zn^\mu \qquad p_j^\mu \rightarrow \hat{p}_j^\mu(z) = p_j^\mu - zn^\mu, \qquad (3.1)$$

where $p_{i/j}$ are the two selected original momenta and $\hat{p}_{i/j}$ are the corresponding shifted momenta. To ensure the shifted momenta stay on-shell it is necessary to choose the shift vector n^μ such that it satisfies the conditions,

$$2n_\mu p_i^\mu + zn^2 = -2n_\mu p_j^\mu + zn^2 = 0 \qquad (3.2)$$

and as n and $p_{i/j}$ are independent of z , this relation must be true at all orders in z and the conditions simplify to

$$n_\mu p_i^\mu = n_\mu p_j^\mu = n^2 = 0, \qquad (3.3)$$

The solution to these conditions used for the BCFW relation is given by

$$n^\mu = \langle i\gamma^\mu j \rangle , \quad (3.4)$$

which satisfies the conditions explicitly. This shift is especially useful for calculations using the Spinor Helicity Formalism as it corresponds to a shift of the spinors given by[10]

$$|i\rangle \rightarrow |i(\hat{z})\rangle = |i\rangle + z|j\rangle \quad |j\rangle \rightarrow |j(\hat{z})\rangle = |j\rangle - z|i\rangle , \quad (3.5)$$

where all the other spinors are unchanged by the shift.

When the shift is applied the resulting amplitude, $\hat{A}(z)$, is a function of the complex parameter z . The recurrence relation is derived by performing a contour integration of $\hat{A}(z)/z$, around a circle of radius r , and taking the limit of $r \rightarrow \infty$. In this limit the Cauchy residue theorem shows that the integral is given by

$$\lim_{a \rightarrow \infty} \oint_{|z|=a} \frac{\hat{A}(z)}{z} dz = 2\pi i \left(\hat{A}(0) + \sum_{z_0} \frac{\text{Res}_{z=z_0} \hat{A}(z)}{z_0} \right) , \quad (3.6)$$

where the sum is over every pole, z_0 , in the shifted amplitude. By the analyticity properties of amplitudes, the only poles in the shifted amplitude will be due to internal propagators going on-shell and will be single poles whose residues, by the optical theorem, are shown to be proportional to the product of two lower multiplicity, on-shell amplitudes. The form of the pole in the amplitude expanded around the location of the pole is given by

$$\begin{aligned} \hat{A}(z) &= \frac{\hat{A}_l(z)\hat{A}_r(z)}{\hat{P}(z)^2} + \tilde{A}(z) \\ &= \frac{\hat{A}_l(z)\hat{A}_r(z)}{(P + zn)^2} + \tilde{A}(z) \\ &= \frac{\hat{A}_l(z)\hat{A}_r(z)}{P^2 + 2zP \cdot n} + \tilde{A}(z) \\ &= \frac{\hat{A}_l(z)\hat{A}_r(z)}{2P \cdot n} \frac{1}{z - \frac{-P^2}{2P \cdot n}} + \tilde{A}(z) , \end{aligned} \quad (3.7)$$

where $\hat{A}_{l/r}(z)$ are the two amplitudes either side of the selected propagator, $\hat{P}(z)$ is the shifted momentum in the propagator, n is the shift vector, $P = \sum p$ is the unshifted momentum of the selected propagator made up of a sum of external momenta including p_i but not p_j and $\tilde{A}(z)$ is the rest of the amplitude that does not have any

poles at $z \rightarrow z_0$. Without loss of generality, the left amplitude and the propagator momentum can be taken to include p_i and not p_j . This is because exchanging the left and right sides will give the same terms again and including both would be a double counting. The remaining cases are when both p_i and p_j are on the same side of the propagator, which causes the z dependence to cancel and hence such terms will not contribute. Taking the residue of this pole and substituting it into the equation for the integral produces the result

$$\begin{aligned}
A = \hat{A}(0) &= - \sum_m \frac{A_{l,m} A_{r,m}}{2P_i \cdot n z_{0,m}} - \frac{1}{2\pi i} \lim_{a \rightarrow \infty} \oint_{|z|=a} \frac{\hat{A}(z)}{z} dz \\
&= - \sum_m \frac{A_{l,m} A_{r,m}}{2P_m \cdot n \frac{-P_m^2}{2P_m \cdot n}} - \frac{1}{2\pi i} \lim_{a \rightarrow \infty} \oint_{|z|=a} \frac{\hat{A}(z)}{z} dz \\
&= \sum_m \frac{A_{l,m} A_{r,m}}{P_m^2} - \frac{1}{2\pi i} \lim_{a \rightarrow \infty} \oint_{|z|=a} \frac{\hat{A}(z)}{z} dz , \tag{3.8}
\end{aligned}$$

where the sum is over the poles labelled by m , and $A_{l/r,m} = \hat{A}_{l/r} \left(\frac{-P_m^2}{2P_m \cdot n} \right)$ are the on-shell amplitudes resulting from splitting the diagram at the propagator labelled by m and evaluating for the shift parameter $z = \frac{-P_m^2}{2P_m \cdot n}$. If the integral can be shown to be zero then the BCFW recursion relation results. To do this without explicit integration or using the residue theorem, power counting is used. In some cases power counting can show us that the integral is zero, for all other cases the integral's value cannot be determined in this way. For pure gluon amplitudes this gives the condition that[10]

$$h_i = + \text{ or } h_j = - \text{ or both } , \tag{3.9}$$

where $h_{i/j}$ are the helicities of the gluons i and j respectively. This is equivalent to saying that the helicities for the shifted gluons (h_i, h_j) must be either $(+, +)$, $(-, -)$ or $(-, +)$ but not $(+, -)$. If a shift needs to be performed on a pair of particles with helicities $(+, -)$ then the shift can be performed by swapping the roles of the gluons in the shift or equivalently by swapping $|\cdot\rangle \rightarrow |\cdot\rangle$ in the definition of the shift.

For MHV amplitudes it is easy to see that this condition is the correct condition. The amplitude only depends on $|\cdot\rangle$ spinors so the shift only has an effect through the spinor $|i\rangle$ being replaced with $|i\rangle + z|j\rangle$. If the gluon i has helicity $+$, then this spinor is only present in the amplitude in the denominator and as such, in the limit $|z| \rightarrow \infty$, the shifted amplitude will be proportional to $1/z^2$ or, if i and j are neighbouring particles, $1/z$ and the amplitude will vanish. For the case of a negative

helicity for gluon i , if the other gluon shifted is the other negative helicity gluon, then the numerator has the form $\langle ij \rangle^4 \rightarrow (\langle ij \rangle + z \langle jj \rangle)^4 = \langle ij \rangle^4$ which is unchanged, so again the only contribution from the shift is in the denominator and the amplitude will vanish in the limit $|z| \rightarrow \infty$. The remaining case, which is the amplitude that should not vanish, is that gluon i has negative helicity and j has positive helicity. In this case the numerator has the form $\langle ik \rangle^4 \rightarrow (\langle ik \rangle + z \langle jk \rangle)^4$, where gluon k is the other negative helicity gluon. As $|z| \rightarrow \infty$ this form tends to ∞ and the one or two powers of z in the numerator can not save the amplitude from tending to ∞ .

To work out the trend for a general amplitude is more complex and requires balancing the powers of z from the numerator and denominator of an amplitude. For a general pure gluon amplitude, from the Feynman rules, it can be seen that the contributions dependant on z in any diagram can only be from the single route through the diagram, from one of the shifted particles to the other one. The contributions along this route will be from propagators and from three particle vertices. The form of a gluon propagator from the internal momenta, P , has the form

$$\frac{1}{\hat{P}^2} = \frac{1}{P^2 + zP \cdot n} \xrightarrow{|z| \rightarrow \infty} \frac{1}{P \cdot n} \frac{1}{z} = 0, \quad (3.10)$$

where P was chosen to contain i and therefore not j , and the form of a quark propagator for the internal momenta, P , has the form

$$\frac{\hat{P}}{\hat{P}^2} = \frac{\not{P} + z(|j\rangle [i] + |i\rangle \langle j|)}{P^2 + zP \cdot n} \xrightarrow{|z| \rightarrow \infty} \frac{\not{P}}{P \cdot n} \frac{1}{z} + \frac{|j\rangle [i] + |i\rangle \langle j|}{P \cdot n} = 0. \quad (3.11)$$

The form of a three gluon vertex is

$$V_{3 \text{ gluon}}^{\mu \nu \sigma}_{k_1 k_2 k_3} = \frac{1}{\sqrt{2}} (z (2g^{\mu\nu} n^\sigma - g^{\nu\sigma} n^\mu - g^{\sigma\mu} n^\nu) + g^{\mu\nu} (k_1 - k_2)^\sigma + g^{\mu\nu} (k_2 - k_3)^\mu + g^{\mu\nu} (k_3 - k_1)^\nu), \quad (3.12)$$

where k_1 contains i and k_2 contains j and k_3 therefore contains no z dependence. From this it can be seen that for any diagram the propagator and vertex combination will depend on z at the worst as $\sim z$ and possibly as constant or lower powers of z if some of the numerator terms vanish. The polarisation vectors for i and j can each

either contribute $1/z$ or z , depending on their helicities, as they take the forms

$$\hat{\epsilon}_\mu^+(i; q) = \frac{[i|\gamma_\mu|q\rangle}{\sqrt{2}(\langle iq\rangle + z\langle jq\rangle)} \sim \frac{1}{z} \frac{[i|\gamma_\mu|q\rangle}{\sqrt{2}\langle jq\rangle} \quad (3.13)$$

$$\hat{\epsilon}_\mu^-(i; q) = -\frac{\langle i|\gamma_\mu|q\rangle + z\langle j|\gamma_\mu|q\rangle}{\sqrt{2}[iq]} \sim -z \frac{\langle j|\gamma_\mu|q\rangle}{\sqrt{2}[iq]} \quad (3.14)$$

$$\hat{\epsilon}_\mu^-(j; q') = -\frac{\langle j|\gamma_\mu|q'\rangle}{\sqrt{2}(\langle jq\rangle - z\langle iq'\rangle)} \sim \frac{1}{z} \frac{\langle j|\gamma_\mu|q'\rangle}{\sqrt{2}\langle iq'\rangle} \quad (3.15)$$

$$\hat{\epsilon}_\mu^+(j; q') = \frac{[j|\gamma_\mu|q'\rangle - z[i|\gamma_\mu|q'\rangle]}{\sqrt{2}\langle jq'\rangle} \sim -z \frac{[i|\gamma_\mu|q'\rangle}{\sqrt{2}\langle jq'\rangle}. \quad (3.16)$$

From this it can be seen that if gluon i has helicity $+$ and gluon j has helicity $-$, then the amplitude will at most go as $\frac{1}{z}$ for $|z| \rightarrow \infty$ and the integral at infinity will vanish. For the other valid helicity cases, the leading terms seem to go as z , but these leading terms can be shown to vanish, leaving a correct leading term of $\frac{1}{z}$.

This logic can be extended to amplitudes of quarks using the same methods. For this project it also needs to be extended to amplitudes with a complex scalar, ϕ , in them, which requires showing that these amplitudes vanish in the limit of $|z| \rightarrow \infty$. It is easy to see that if the amplitudes without the ϕ vanish, the ones with it will also vanish, as the amplitudes are related and the amplitudes that completely vanish without the ϕ will all vanish for any choice of shift particles.

Chapter 4

Generalised Unitarity Method

The Generalised Unitarity method has been used by many people to derive analytical forms for various one loop amplitudes[11, 12, 13]. It is a method that allows one loop amplitudes to be built up from tree level amplitudes without directly performing any integrations.

Any one loop amplitude can be expanded in terms of a basis of the scalar one loop amplitudes. These amplitudes are those containing only scalar particles. It is only necessary to use the subset of these amplitudes with all particles in the loop being massless and only the external particles entering the loop having mass, as in this project all particles are massless, other than the Higgs boson or complex scalar, ϕ , neither of which can enter the loop as that would introduce higher powers of the effective coupling and not be NLO. These scalar one loop integrals are of the form[14]

$$I_n(P_1, \dots, P_n) = \int \frac{d^D l}{i\pi^{\frac{D}{2}}} \frac{1}{(l)^2 \dots (l - P_1 - \dots - P_{n-1})^2}, \quad (4.1)$$

where n is the number of propagators in the loop which is the number of external scalars entering the loop, P_1, \dots, P_n are the momenta of the external scalars entering the loop which must sum to zero and l is the loop momenta. Any one loop integral with massless propagators can be written as a sum of these integrals multiplied by rational functions of spinor products and Minkowski products of the external momenta, as contributions that are divergent in the limit $d \rightarrow 4$ are all due to integrating propagators over the loop momentum. Therefore, when using this basis any contributions that depend on the nature of the particles in the amplitude will be included in coefficients of these integrals. This basis is chosen as it is one of the simplest possible sets

and any amplitude will have a unique expansion in terms of this basis. Any given integral can be simplified and numerator terms removed using the Passarino-Veltman reduction to write them as combinations of metric tensors and loop momenta squared, which reduces the integral to a combination of the basis integrals. Then any amplitudes with more than four propagators can then be reduced to amplitudes with at most four propagators. In four dimensions only loop integrals with at most four propagators are needed. For more than five propagators a simple partial fractioning of the amplitude will reduce it to a sum of terms with at most five propagators multiplied by rational functions of the Minkowski products of external momenta, as the set of external momenta is linearly dependent. The remaining terms with five propagators can again be reduced to give a combination of terms with at most four propagators along with terms that contain enough explicit powers of the dimension to cancel any poles from the integration and as such do not contribute to the cut producible parts of the amplitude. After this reduction an amplitude will be written in the form of sums of scalar one loop integrals with at most four propagators multiplied by rational functions of spinors and Minkowski products of external momenta plus a term with no poles in ϵ where $d = 4 - 2\epsilon$ is the dimension of space time. This expansion is shown in [Figure 4.1](#). The full set of these integrals have been evaluated explicitly as functions of the external momenta flowing into each vertex of the loop.

To evaluate any one loop amplitude, all that remains to be calculated are the coefficients of each of these scalar loop functions. These can be calculated by taking any amplitude and making use of various methods to reduce the amplitude to an explicit expansion in terms of these functions. However, this becomes intractable when the number of particles increases and is complex to implement numerically. For a more efficient method generalised unitarity can be used. This extends the unitarity method to multiple cuts for loop amplitudes. A cut is defined as the process of replacing a propagator by a delta function or more precisely of introducing the factor

$$2\pi (l - P)^2 \delta((l - P)^2) , \quad (4.2)$$

where l is the loop momenta and $\frac{1}{(l-P)^2}$ is the propagator that is being cut. This factor will integrate to zero unless multiplied by a function containing exactly the propagator to which it corresponds. If cuts are performed on a loop amplitude it will split it into a sum of products of tree amplitudes due to the factorisation properties

$$A = \sum_i d_i \text{---} \text{---} \text{---} \text{---} \text{---} + \sum_i c_i \text{---} \text{---} \text{---} + \sum_i b_i \text{---} \text{---} \text{---} + \sum_i a_i \text{---} \text{---} \text{---} + R$$

Figure 4.1: The expansion of an amplitude in terms of the scalar basis functions. The sums are over each possible scalar loop function of that form, defined by the momenta at each corner and R contains all contributions that do not contain poles in ϵ .

of amplitudes. The same cuts, when performed on the expansion, explicitly isolate coefficients of different loop functions. Four cuts will isolate the coefficient of a single box integral, as no integrals have more than four propagators and no other scalar loop functions will be extracted by this set of cuts because any other scalar loop integral will not contain all four of the propagators. Therefore, at least one of the cuts will not be matched by a propagator and will vanish when integrated. By performing each possible set of four cuts, all box coefficients can be extracted. For three cuts, in the expansion in terms of scalar loop functions, a single triangle function will be isolated, but there will also be contributions from box functions. However, triangle and lower cuts are not required for boxes to be extracted, so their contributions can be removed before attempting to calculate triangle coefficients. This same process applies for one or two cuts and allows us to isolate each coefficient in turn.

After applying three cuts to the raw amplitude, three of the four degrees of freedom in the loop momenta are removed. If the product of the three tree amplitudes is performed leaving the remaining degree of freedom as a free parameter, after subtracting the poles due to boxes, the result will be a power series in the free parameter. As will be shown later in this chapter, for the parametrisation used in this project, the coefficient to extract is the constant term in this expansion, as all other terms will vanish when integrated over the contour corresponding to the remaining degree of freedom from the momentum integration. For analytical methods this coefficient can be extracted by explicit rearrangement of the expression to isolate the needed term. For numerical methods this is not possible, instead, a discrete complex Fourier projection can be used to extract the coefficient, as long as the range of possible powers that can appear in the expansion is known. The continuous version of the projection for the term in the expansion of $f(z)$ of the form z^n is given by

$$P_n[f(z)] = \frac{1}{2\pi} \int_{\theta_0}^{\theta_0+2\pi} f(z_0 e^{it}) (z_0 e^{it})^{-n} dt, \quad (4.3)$$

where z_0 is an arbitrary complex value and θ_0 is an arbitrary real angle. For any power series, each term can be treated independently, as integration is linear and therefore, to prove the projection works, an expression of the form z^m can be substituted and then it can be shown that the result is 1 if $m = n$ and 0 otherwise. For $m = n$ the relation is given by

$$\begin{aligned}
P_n[n^n] &= \frac{1}{2\pi} \int_{\theta_0}^{\theta_0+2\pi} (z_0 e^{it})^n (z_0 e^{it})^{-n} dt \\
&= \frac{1}{2\pi} \int_{\theta_0}^{\theta_0+2\pi} (z_0 e^{it})^{n-n} dt \\
&= \frac{1}{2\pi} \int_{\theta_0}^{\theta_0+2\pi} 1 dt \\
&= \frac{1}{2\pi} [t]_{\theta_0}^{\theta_0+2\pi} \\
&= 1,
\end{aligned} \tag{4.4}$$

and for $m \neq n$ it is

$$\begin{aligned}
P_n[n^m] &= \frac{1}{2\pi} \int_{\theta_0}^{\theta_0+2\pi} (z_0 e^{it})^m (z_0 e^{it})^{-n} dt \\
&= \frac{1}{2\pi} \int_{\theta_0}^{\theta_0+2\pi} (z_0 e^{it})^{m-n} dt \\
&= \frac{1}{2\pi} \left[\frac{-i}{m-n} (z_0 e^{it})^{m-n} \right]_{\theta_0}^{\theta_0+2\pi} \\
&= \frac{1}{2\pi} \left(\left[\frac{-i}{m-n} (z_0 e^{i\theta_0})^{m-n} \right] - \left[\frac{-i}{m-n} (z_0 e^{i(\theta_0+2\pi)})^{m-n} \right] \right) \\
&= \frac{1}{2\pi} \frac{-i}{m-n} \left[(z_0 e^{i\theta_0})^{m-n} - (z_0 e^{i\theta_0})^{m-n} \right] \\
&= 0.
\end{aligned} \tag{4.5}$$

The discrete version to extract the coefficient of z^n using N evaluations is given by

$$D_{n,N}[f(z)] = \frac{1}{N} \sum_{j=0}^{N-1} f\left(z_0 e^{i2\pi \frac{j}{N}}\right) \left(z_0 e^{i2\pi \frac{j}{N}}\right)^{-n}, \tag{4.6}$$

The proof proceeds similar to before but now, if there are N terms in the projection, the result is 1 for $n = m \pmod N$ and 0 for $n \neq m \pmod N$. $m = n + kN$, with any integer k , results in

$$D_{n,N}[z^{n+kN}] = \frac{1}{N} \sum_{j=0}^{N-1} \left(z_0 e^{i2\pi \frac{j}{N}}\right)^{n+kN} \left(z_0 e^{i2\pi \frac{j}{N}}\right)^{-n}$$

$$\begin{aligned}
&= \frac{1}{N} \sum_{j=0}^{N-1} \left(z_0 e^{i2\pi \frac{j}{N}} \right)^{(n+kN)-n} \\
&= \frac{1}{N} \sum_{j=0}^{N-1} \left(z_0 e^{i2\pi \frac{j}{N}} \right)^{kN} \\
&= \frac{1}{N} z_0^{kN} \sum_{j=0}^{N-1} e^{i2\pi k j \frac{N}{N}} \\
&= \frac{1}{N} z_0^{kN} \sum_{j=0}^{N-1} e^{i2\pi k j} \\
&= \frac{1}{N} z_0^{kN} \sum_{j=0}^{N-1} 1^{kj} \\
&= \frac{1}{N} z_0^{kN} \sum_{j=0}^{N-1} 1 \\
&= \frac{1}{N} z_0^{kN} N \\
&= z_0^{kN} \\
&= 1 \text{ if } k = 0 ,
\end{aligned} \tag{4.7}$$

and for $m \neq n \pmod N$ it is given by

$$\begin{aligned}
D_{n,N}[z^k] &= \frac{1}{N} \sum_{j=0}^{N-1} \left(z_0 e^{i2\pi \frac{j}{N}} \right)^m \left(z_0 e^{i2\pi \frac{j}{N}} \right)^{-n} \\
&= \frac{1}{N} \sum_{j=0}^{N-1} \left(z_0 e^{i2\pi \frac{j}{N}} \right)^{m-n} \\
&= \frac{1}{N} z_0^{m-n} \sum_{j=0}^{N-1} e^{i2\pi j \frac{m-n}{N}} \\
&= \frac{1}{N} z_0^{m-n} \sum_{j=0}^{N-1} \left(e^{i2\pi \frac{m-n}{N}} \right)^j \\
&= \frac{1}{N} z_0^{m-n} \times 0 \\
&= 0 .
\end{aligned} \tag{4.8}$$

For $m - n \neq 0 \pmod N$, $e^{i2\pi \frac{m-n}{N}}$ is an N^{th} root of unity, and when raised to N successive integers, each root is given exactly once. Combining these with the fact that the sum of all the N^{th} roots of unity is exactly 0, allows the fifth line of [Equation 4.8](#) to be derived. From [Equations 4.7](#) and [4.8](#) it is clear that if the expansion contains terms from n_{\min} to n_{\max} then a projection with at least $n_{\max} - n_{\min} + 1$ terms is needed.

The powers of these remaining parameters will come from numerator terms left after the diagrams have been reduced by cancelling with propagators. For boxes, as there are four propagators and external momenta are four dimensional, any combination that contains loop momenta squared or higher powers will be cancellable against a propagator and therefore should be contained in triangles or bubbles instead. Therefore, the highest power of the loop momenta that can remain is a term linear in the loop momenta. For triangles there is now the possibility of combinations that no longer fully cancel against propagators and are in directions orthogonal to components that could be cancelled against loop momenta. The highest power can be calculated using power counting to derive the highest power that could come from any diagram and as any extra powers in the denominator must cancel to reduce the diagram to a combination of our basis, powers of the loop momenta appearing in the denominator can be cancelled against those from the numerator.

For a normal renormalisable theory, for example the case of pure gluons amplitudes, each propagator provides two powers of the loop momenta in the denominator and each propagator must be connected by either a three or four point vertex. As four point vertices do not contain any powers of loop momenta, they will not contribute to the highest power of the loop momenta. The highest, therefore, must come from terms containing three point vertexes which each contribute a power of the loop momenta. Therefore, the highest power would come from a term with the minimum number of propagators and a three point vertex on each corner which for triangles is three powers of the loop momenta. For the case of amplitudes with a complex scalar, ϕ , they contain one power higher of momenta and their Feynman rules include a two gluon vertex with two powers of the particle momenta. This term would contribute one higher power of the loop momenta and therefore would have a numerator with up to four powers of the loop momenta.

For bubbles there is one less propagator and therefore one less vertex in the cases that produce the highest powers of the loop momenta in the numerator and their highest power possible is exactly one less than is possible for the triangles in the same theory. Bubbles, therefore, have up to two powers of the loop momenta in the numerator in renormalisable theories and for amplitudes with a ϕ up to three powers of the loop momenta.

This, along with the ability to calculate the tree amplitudes, is the only change needed to extend this for amplitudes with ϕ , as all the properties used apply to

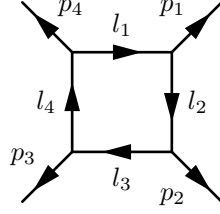


Figure 4.2: A box digram showing the labelling of the momenta and the direction each momenta is taken to be in.

any amplitude. To implement this efficiently and numerically, explicit formulas are required for the loop momenta solutions for each set of cuts. Firstly, for boxes, the relations that need to be solved are

$$\begin{aligned}
 l_1^2 &= \\
 (l_1 - p_1)^2 &= \\
 (l_1 - p_1 - p_2)^2 &= \\
 (l_1 - p_1 - p_2 - p_3)^2 &= 0, \tag{4.9}
 \end{aligned}$$

where $p_{1,2,3,4}$ are the external momenta outbound from each corner and l_1 is the loop momentum heading into corner 1 from corner 2 as shown in [Figure 4.2](#). If the momenta flowing out of at least one of the corners is massless, then the solutions take a particularly simple form given by [\[6\]](#)

$$\begin{aligned}
 l_1^\mu &= \frac{\langle 1|2|3|4|\gamma^\mu|1\rangle}{2\langle 1|2|4|1\rangle} & l_1'^\mu &= \frac{[1|2|3|4|\gamma^\mu|1]}{2[1|2|4|1]} \\
 l_2^\mu &= -\frac{\langle 1|\gamma^\mu|2|3|4|1\rangle}{2\langle 1|2|4|1\rangle} & l_2'^\mu &= -\frac{[1|\gamma^\mu|2|3|4|1]}{2[1|2|4|1]} \\
 l_3^\mu &= \frac{\langle 1|2|\gamma^\mu|3|4|1\rangle}{2\langle 1|2|4|1\rangle} & l_3'^\mu &= \frac{[1|2|\gamma^\mu|3|4|1]}{2[1|2|4|1]} \\
 l_4^\mu &= -\frac{\langle 1|2|3|\gamma^\mu|4|1\rangle}{2\langle 1|2|4|1\rangle} & l_4'^\mu &= -\frac{[1|2|3|\gamma^\mu|4|1]}{2[1|2|4|1]}, \tag{4.10}
 \end{aligned}$$

where without loss of generality corner 1 has been chosen to be a massless corner and $l_{1,2,3,4}$ and $l'_{1,2,3,4}$ are the two different solutions to the relations. These momenta satisfy the relations shown in [Equation 4.9](#) if it can be shown that all these vectors are massless and that the difference between neighbouring loop momenta is the appropriate corners momenta. That they are all massless is easy to see from the form of the momenta and the relations for spinors. For corner 1 the momentum difference

relation is given by

$$\begin{aligned}
l_1^\mu - l_2^\mu &= \frac{\langle 1|2|3|4|\gamma^\mu|1\rangle}{2\langle 1|2|4|1\rangle} - \frac{\langle 1|\gamma^\mu|2|3|4|1\rangle}{2\langle 1|2|4|1\rangle} \\
&= \frac{\langle 1|2|3|4|\gamma^\mu|1\rangle + \langle 1|\gamma^\mu|2|3|4|1\rangle}{2\langle 1|2|4|1\rangle} \\
&= \frac{p_4^\mu \langle 1|2|3|1\rangle - \langle 1|2|3|\gamma^\mu|4|1\rangle + \langle 1|\gamma^\mu|2|3|4|1\rangle}{2\langle 1|2|4|1\rangle} \\
&= \frac{2p_4^\mu \langle 1|2|3|1\rangle - 2p_3^\mu \langle 1|2|4|1\rangle + \langle 1|2|\gamma^\mu|3|4|1\rangle + \langle 1|\gamma^\mu|2|3|4|1\rangle}{2\langle 1|2|4|1\rangle} \\
&= \frac{2p_4^\mu \langle 1|2|3|1\rangle - 2p_3^\mu \langle 1|2|4|1\rangle + 2p_2^\mu \langle 1|2|3|4|1\rangle - \langle 1|\gamma^\mu|2|3|4|1\rangle + \langle 1|\gamma^\mu|2|3|4|1\rangle}{2\langle 1|2|4|1\rangle} \\
&= -p_3^\mu + \frac{p_4^\mu \langle 1|2|3|1\rangle + p_2^\mu \langle 1|3|4|1\rangle}{\langle 1|2|4|1\rangle} \\
&= -p_3^\mu + \frac{-p_4^\mu \langle 1|2|4|1\rangle - p_2^\mu \langle 1|2|4|1\rangle}{\langle 1|2|4|1\rangle} \\
&= -p_3^\mu - p_4^\mu - p_2^\mu \\
&= p_1^\mu, \tag{4.11}
\end{aligned}$$

where on the middle three lines commutation relations for slashed matrices have been used and on the last four lines momentum conservation between the external momenta has been used. The proofs for each of the other relations follow a similar method but are easier to show.

These relations only work if there is at least one momenta that is massless, which is true for pure gluon and quark amplitudes with seven or less external particles. For amplitudes with Higgs bosons, complex scalars, ϕ , or similarly any other external massive particles, these solutions are no longer enough for seven external particles, so a more general solution is required. To derive the general case, which is also the form of solution used for triangles and bubbles, a general basis of four massless independent vectors is needed. As there are in general no massless corners, the first step is to take two of the corners, which will be called corners 1 and 2 and their outbound momentum vectors K_1 and K_2 and project a pair of massless vectors by projecting them on each other. There are various definitions for how to perform the projection, but the one used here is a simplified version of that used by Berger and Forde[11] that leads to simpler formulas when working algebraically. This solution defines the two massless

projections, \tilde{K}_1 and \tilde{K}_2 as

$$\begin{aligned}\tilde{K}_1 &= K_1 - \frac{K_1^2}{\gamma} K_2 \\ \tilde{K}_2 &= K_2 - \frac{K_2^2}{\gamma} K_1 ,\end{aligned}\tag{4.12}$$

where γ is fixed such that the two projected vectors are massless and is given by

$$\begin{aligned}\gamma &= K_1 \cdot K_2 \left(1 \pm \sqrt{\frac{\Delta}{K_1 \cdot K_2^2}} \right) \\ \Delta &= K_1 \cdot K_2^2 - K_1^2 K_2^2 = \det \begin{pmatrix} K_1^2 & K_1 \cdot K_2 \\ K_1 \cdot K_2 & K_2^2 \end{pmatrix} ,\end{aligned}\tag{4.13}$$

where Δ is the determinant of the matrix of the products of any two of the three momentum conserving momenta, and is therefore independent of the labelling of the momenta. The two solutions for the momenta labeled by the choice of sign in γ are not independent. The solution with one sign is related to the solution with the opposite sign by the relations

$$\begin{aligned}\tilde{K}_1^\mp &= \frac{-\gamma^\pm}{K_2^2} \tilde{K}_2^\pm \\ \tilde{K}_2^\mp &= \frac{-\gamma^\pm}{K_1^2} \tilde{K}_1^\pm ,\end{aligned}\tag{4.14}$$

where the superscripts label which of the signs is chosen. From this it is clear that it is possible to choose either solution and ignore the other, as the results will be the same and using both would be a double counting. The solution with a positive sign is chosen so that the projected vectors are well behaved in the limit when K_1^2 and K_2^2 vanish, as for the negative sign solution, γ will vanish in this limit. The other two vectors used are built from the spinors for the two massless projected vectors and are given by

$$\begin{aligned}n_1 &= \frac{\langle \tilde{K}_1 | \gamma | \tilde{K}_2 \rangle}{2} \\ n_2 &= \frac{\langle \tilde{K}_2 | \gamma | \tilde{K}_1 \rangle}{2} .\end{aligned}\tag{4.15}$$

From these four vectors, any vector in 4 dimensional space can be built, as long as the four vectors are linearly independent, which they will be, if the two original vectors

are linearly independent. These four vectors are particularly convenient to use as the products of various combinations vanish or are very simple. The spinor products that are simple are given by

$$\begin{aligned}
n_1 \cdot n_2 &= -\tilde{K}_1 \cdot \tilde{K}_2 & (4.16) \\
0 &= \tilde{K}_1 \cdot n_1 = \tilde{K}_1 \cdot n_2 \\
&= \tilde{K}_2 \cdot n_1 = \tilde{K}_2 \cdot n_2 \\
&= K_1 \cdot n_1 = K_1 \cdot n_2 \\
&= K_2 \cdot n_1 = K_2 \cdot n_2 .
\end{aligned}$$

From this basis a general loop momenta can be written as

$$l_2 = \alpha \tilde{K}_1 + \beta \tilde{K}_2 + c_1 n_1 + c_2 n_2 , \quad (4.17)$$

where α , β , n_1 and n_2 are arbitrary complex constants. The first condition to apply will be to ensure that the loop momentum for the propagator between corners 1 and 2, l_2 , is on-shell; this will be needed for all cuts. This corresponds to the condition

$$\begin{aligned}
0 &= l_2^2 \\
&= \left(\alpha \tilde{K}_1 + \beta \tilde{K}_2 + c_1 n_1 + c_2 n_2 \right)^2 \\
&= 2 \left(\alpha \beta \tilde{K}_1 \cdot \tilde{K}_2 + c_1 \alpha \tilde{K}_1 \cdot n_1 + c_2 \alpha \tilde{K}_1 \cdot n_2 + c_1 \beta \tilde{K}_2 \cdot n_1 + c_2 \beta \tilde{K}_2 \cdot n_2 + c_1 c_2 n_1 \cdot n_2 \right) \\
&= 2 \left(\alpha \beta \tilde{K}_1 \cdot \tilde{K}_2 + c_1 c_2 n_1 \cdot n_2 \right) \\
&= 2 (\alpha \beta - c_1 c_2) \tilde{K}_1 \cdot \tilde{K}_2 , & (4.18)
\end{aligned}$$

which is used to fix $c_1 c_2 = \alpha \beta$. Which of c_1 and c_2 to fix in terms of the other variables is not specified. This is because, as will be shown below, α and β will be fixed by the next two cuts performed and one or both of them can be zero, in which case the solution for c_1 and c_2 will split into two branches, each with one of the constants vanishing and the other having a non zero value, ranging over all possible values depending on whether a fourth cut is applied. This form for the solution is particularly useful as the expression for the momenta, in terms of spinors, factorises

and is given by

$$\begin{aligned} l_2^\mu &= \left(\langle \tilde{K}_1 | + \frac{\beta}{c_1} \langle \tilde{K}_2 | \right) \gamma^\mu \left(\alpha | \tilde{K}_1] + c_1 | \tilde{K}_2] \right) \\ &= \left(\alpha \langle \tilde{K}_1 | + c_2 \langle \tilde{K}_2 | \right) \gamma^\mu \left(| \tilde{K}_1] + \frac{\beta}{c_2} | \tilde{K}_2] \right) , \end{aligned} \quad (4.19)$$

where the two solutions are equivalent, unless one of K_1^2 and K_2^2 vanishes, in which case they are each valid in one branch of the solution space, as in the other they evaluate to $\frac{0}{0}$.

The next two cuts to be applied, which are needed for both triangles and boxes, are for the propagators either side of the loop momenta, l_2 . These are given by the conditions

$$0 = (l_2 + K_1)^2 \qquad 0 = (l_2 - K_2)^2 , \quad (4.20)$$

for which solutions are given by

$$\begin{aligned} \alpha &= \frac{K_1 \cdot \tilde{K}_2 K_2 \cdot K_2 + K_1 \cdot K_1 K_2 \cdot \tilde{K}_2}{2(K_1 \cdot \tilde{K}_2 \tilde{K}_1 \cdot K_2 - K_1 \cdot \tilde{K}_1 K_2 \cdot \tilde{K}_2)} \\ &= \frac{K_2^2(K_1^2 + \gamma)}{4\Delta} \end{aligned} \quad (4.21)$$

$$\begin{aligned} \beta &= -\frac{K_1 \cdot K_1 \tilde{K}_1 \cdot K_2 + K_1 \cdot \tilde{K}_1 K_2 \cdot K_2}{2(K_1 \cdot \tilde{K}_2 \tilde{K}_1 \cdot K_2 - K_1 \cdot \tilde{K}_1 K_2 \cdot \tilde{K}_2)} \\ &= -\frac{K_1^2(K_2^2 + \gamma)}{4\Delta} . \end{aligned} \quad (4.22)$$

Using these two solutions produces simple spinor representations for the loop momenta, l_1 and l_3 , which are given by

$$\begin{aligned} l_1^\mu &= \left(\langle \tilde{K}_1 | + \frac{\beta(1 - C_{K_1})}{c_1} \langle \tilde{K}_2 | \right) \gamma^\mu \left(\frac{\alpha}{1 - C_{K_1}} | \tilde{K}_1] + c_1 | \tilde{K}_2] \right) \\ &= \left(\frac{\alpha}{1 - C_{K_1}} \langle \tilde{K}_1 | + c_2 \langle \tilde{K}_2 | \right) \gamma^\mu \left(| \tilde{K}_1] + \frac{\beta(1 - C_{K_1})}{c_2} | \tilde{K}_2] \right) \end{aligned} \quad (4.23)$$

$$\begin{aligned} l_3^\mu &= \left(\langle \tilde{K}_1 | + \frac{\beta}{c_1(1 - C_{K_3})} \langle \tilde{K}_2 | \right) \gamma^\mu \left(\alpha(1 - C_{K_3}) | \tilde{K}_1] + c_1 | \tilde{K}_2] \right) \\ &= \left(\alpha(1 - C_{K_3}) \langle \tilde{K}_1 | + c_2 \langle \tilde{K}_2 | \right) \gamma^\mu \left(| \tilde{K}_1] + \frac{\beta}{c_2(1 - C_{K_3})} | \tilde{K}_2] \right) , \end{aligned} \quad (4.24)$$

where C_{K_1} and C_{K_2} are constants given by

$$C_{K_1} = \frac{2(\gamma - K_1 \cdot K_2)}{\gamma + K_2^2} \quad (4.25)$$

$$C_{K_2} = \frac{2(\gamma - K_1 \cdot K_2)}{\gamma + K_1^2}. \quad (4.26)$$

As for l_2 , the two solutions are equivalent unless K_1^2 or $K_2^2 \rightarrow 0$, in which case they correspond to the two different branches of the solution.

The final cut is given by

$$0 = (l_2 - K_2 - K_3)^2, \quad (4.27)$$

and is solved to give

$$\begin{aligned} c_1 &= \frac{X \left(1 \pm \sqrt{1 - \frac{16\alpha\beta K_3 \cdot n_1 K_3 \cdot n_2}{X^2}} \right)}{4K_3 \cdot n_1} \\ &= \frac{X \left(1 \pm \sqrt{1 + \frac{K_1^2 K_2^2 \Delta_4 (K_1^2 + K_2^2 + 2K_1 \cdot K_2)}{X^2 \Delta^2}} \right)}{4K_3 \cdot n_1} \end{aligned} \quad (4.28)$$

$$\begin{aligned} c_2 &= \frac{X \left(1 \mp \sqrt{1 - \frac{16\alpha\beta K_3 \cdot n_1 K_3 \cdot n_2}{X^2}} \right)}{4K_3 \cdot n_2} \\ &= \frac{X \left(1 \mp \sqrt{1 + \frac{K_1^2 K_2^2 \Delta_4 (K_1^2 + K_2^2 + 2K_1 \cdot K_2)}{X^2 \Delta^2}} \right)}{4K_3 \cdot n_2}, \end{aligned} \quad (4.29)$$

where these forms are again chosen to simplify choices of signs in the limit of K_1^2 or $K_2^2 \rightarrow 0$ and the terms X and Δ_4 are given by

$$\begin{aligned} X &= (K_2 + K_3)^2 - 2\alpha\tilde{K}_1 \cdot (K_2 + K_3) - 2\beta\tilde{K}_2 \cdot (K_2 + K_3) \\ &= \frac{\left((K_1 + K_2)^2 + (K_1 - K_2) \cdot (K_4 - K_3) \right) K_1^2 K_2^2 + 2K_1 \cdot K_2 (K_1^2 K_2 \cdot K_3 + K_2^2 K_1 \cdot K_4)}{2\Delta} \\ &\quad + (K_2 + K_3)^2 \end{aligned} \quad (4.30)$$

$$\begin{aligned} \Delta_4 &= 2K_1 \cdot K_2 K_1 \cdot K_3 K_2 \cdot K_3 - (K_2 \cdot K_3)^2 K_1^2 - (K_1 \cdot K_3)^2 K_2^2 - (K_1 \cdot K_2)^2 K_3^2 + K_1^2 K_2^2 K_3^2 \\ &= \det \begin{pmatrix} K_1^2 & K_1 \cdot K_2 & K_1 \cdot K_3 \\ K_1 \cdot K_2 & K_2^2 & K_2 \cdot K_3 \\ K_1 \cdot K_3 & K_2 \cdot K_3 & K_3^2 \end{pmatrix}. \end{aligned} \quad (4.31)$$

These forms are not directly applicable to bubble diagrams as they only have one independent momentum vector and a basis must be built from at least two vectors. A solution of the same form would be very useful as it would again have a simple form in terms of spinors and would allow for efficient calculations using the same structures.

Therefore, an arbitrary on-shell reference vector, χ , is used as the second vector from which to build the basis. The exact value of this vector is irrelevant and will cancel in all calculations, but care is needed to ensure the vector will not have badly behaved products with any of the external momenta and as such should be chosen to be a complex vector that is not just a complex multiple of a real vector. The main example used in this project for numerical evaluation is $\chi = (1+2i, 1-2i, 1+4i, \sqrt{15})$. Using this vector and the method used above for triangle and boxes with the two cut conditions, equations are arrived at for the loop momenta of[12]

$$\begin{aligned} l_1 &= (1-y)\tilde{K}_1 + \frac{yK_1^2}{2K_1 \cdot \chi} \chi - tn_1 - \frac{y(1-y)K_1^2}{2tK_1 \cdot \chi} n_2 \\ l_2 &= -y\tilde{K}_1 - \frac{(1-y)K_1^2}{2K_1 \cdot \chi} \chi - tn_1 - \frac{y(1-y)K_1^2}{2tK_1 \cdot \chi} n_2, \end{aligned} \quad (4.32)$$

where t and y are the two degrees of freedom left after the two cuts. These momenta can again be written in a compact spinor form as

$$\begin{aligned} l_1^\mu &= \left(-t \langle \tilde{K}_1 | + \frac{yK_1^2}{2K_1 \cdot \chi} \langle \chi | \right) \gamma^\mu \left(-\frac{1-y}{t} | \tilde{K}_1] + | \chi] \right) \\ l_2^\mu &= \left(-t \langle \tilde{K}_1 | - \frac{(1-y)K_1^2}{2K_1 \cdot \chi} \langle \chi | \right) \gamma^\mu \left(\frac{y}{t} | \tilde{K}_1] + | \chi] \right). \end{aligned} \quad (4.33)$$

To perform the integrals, the Jacobian factors, J , and the integration contours for these choices of loop momenta representation are needed. The Jacobian factors result from both the change in integration variables from the loop momenta itself to the coefficients of our momenta definition and from performing the integrations over the delta functions. These can be given by the formula

$$J = \left| \frac{\sqrt{\det \left(\frac{\partial l}{\partial x_i} \cdot \frac{\partial l}{\partial x_j} \right)}}{\det \left(\frac{\partial d_k}{\partial x_l} \right)} \right|, \quad (4.34)$$

where x_i and x_j run over the set of new integration variables, d_k runs over the set of functions inside the delta functions and x_l runs over the variables integrated out with the delta functions. For example, for triangles this is due to changing from the loop momenta to the four variables α, β, c_1 and c_2 and then integrating out α, β and either c_1 or c_2 using the three delta functions. For this case x_i and x_j would run over all four of the variables α, β, c_1 and c_2 and x_l would run over α, β and whichever of c_1 and c_2 is integrated out.

For the case of the general box cut this gives a Jacobian factor, J_4 , of

$$\begin{aligned}
J_4 &= \left| \frac{\tilde{K}_1 \cdot \tilde{K}_2}{16 \left(K_1 \cdot \tilde{K}_1 K_2 \cdot \tilde{K}_2 - K_1 \cdot \tilde{K}_2 \tilde{K}_1 \cdot K_2 \right) (K_3 \cdot n_1 c_1 - K_3 \cdot n_2 c_2)} \right| \\
&= \frac{1}{8} \left| \sqrt{\frac{\Delta}{X^2 \Delta^2 + \Delta_4 K_1^2 K_2^2 (K_1 + K_2)^2}} \right|. \tag{4.35}
\end{aligned}$$

For the case of triangle solutions the Jacobian factor is

$$\begin{aligned}
J_3 &= \left| \frac{\tilde{K}_1 \cdot \tilde{K}_2}{8 \left(K_1 \cdot \tilde{K}_1 K_2 \cdot \tilde{K}_2 - K_1 \cdot \tilde{K}_2 \tilde{K}_1 \cdot K_2 \right) c} \right| \\
&= \frac{1}{8} \left| \frac{1}{c \sqrt{\Delta}} \right|, \tag{4.36}
\end{aligned}$$

where c is whichever of c_1 and c_2 is the remaining free parameter. For the case of bubble solutions the Jacobian factor is

$$\begin{aligned}
J_2 &= \left| \frac{\sqrt{(n_1 \cdot n_2)^2 (\tilde{K}_1 \cdot \chi)^2}}{4t K_1 \cdot \chi n_1 \cdot n_2} \right| \\
&= \frac{1}{4} \left| \frac{1}{t} \right|. \tag{4.37}
\end{aligned}$$

To calculate the conditions for the integration contour, it should be noted that the full integration space of the loop momenta is over all real momenta, so the contour will cover the set of parameters in the loop momenta that cause the loop momenta to be real. To evaluate these conditions the spinor representations are used along with the condition that for real vectors $\langle l |^\dagger \propto |l\rangle$. For triangles, this relation simplifies to

$$c^* c = \text{const}, \tag{4.38}$$

where c is whichever of c_1 and c_2 is the free parameter and const is a positive function of the external momenta that, as all poles other than those at $c = 0$ are subtracted leaving only a power series in c , is irrelevant by Cauchy residue theorem. For bubbles, as there are two parameters, the solution space is a surface rather than a line, but again a simple relation can be found which is

$$t^* t = \text{const } y(1 - y), \tag{4.39}$$

where again const is a positive function of external momenta and therefore, the solution space is that $0 \leq y \leq 1$ and t is integrated round a circle whose radius is a function of y but, as all poles other than those at $t = 0$ are removed, this simplifies to an integration round any circle centred on the origin.

The exact values of the Jacobian expressions are not relevant, as the same factor would appear when integrating both the scalar loop and the true amplitude terms. The relevant parts are the forms of these in terms of the free parameters, as the Jacobian factors must be well behaved at all relevant points and therefore not contribute poles and, when combined with the forms of the contours, will determine which terms in the integrand will be extracted by performing the remaining integrals. For triangles, as the Jacobian is of the form $1/c$ and the contour is a circle around 0, it is the constant terms in the expansion of the triangles in terms of c that will contribute when integrated and all others will vanish. They will still be needed for subtracting from bubbles, as the other coefficients in terms of c can contribute to the poles that need subtracting. For bubbles, in terms of t , it is the constant term that is needed, again using the same logic as for triangles. To numerically evaluate the integrals, it is necessary to know what ranges of powers there can be in each variable. For this momentum parametrisation the integrand must be a polynomial of order $p - 1$ of the monomials y , t and $\frac{y(1-y)}{t}$ and therefore, for the t^0 terms, the range of powers of y is from 0 to $p - 1$. From direct evaluation of the y integral it is clear that the relevant contribution is given by

$$B = \sum_n \frac{b_{0,n}}{n+1}, \quad (4.40)$$

where $b_{i,j}$ is the coefficient of $t^i y^j$ in the expansion of the bubble after removing poles due to boxes and triangles. To extract all these coefficients requires $n \times m$ evaluations, where n is one more than the difference between the lowest and highest power of t and m is one more than the highest power of m , as the lowest power of y is 0. However, this is not the most efficient that can be achieved as instead of extracting each coefficient and combining, a simpler expression can be produced that will directly produce this combination. If the highest power of y is 3, as it is for standard QCD amplitudes, then the formula

$$B = \frac{1}{5} \left(b_0(0) + 3b_0 \left(\frac{2}{3} \right) \right), \quad (4.41)$$

will extract the relevant contribution, where $b_i(y_0)$ is the coefficient of t^i evaluated for $y = y_0$. This relation will not work for amplitudes with a complex scalar, ϕ , as a

y^4 term is present, therefore a more general relation is needed that is correct for one more power in y . The formula needed is

$$B = \frac{1}{2} \left(b_0 \left(\frac{1}{2} \left(1 + \frac{1}{\sqrt{3}} \right) \right) + b_0 \left(\frac{1}{2} \left(1 - \frac{1}{\sqrt{3}} \right) \right) \right). \quad (4.42)$$

If even higher powers of y occurred then a relation with more evaluations would be needed but they are not required for this project.

With these terms the boxes can now be evaluated directly, but for the triangles and bubbles one more contribution is needed, the explicit forms of the subtractions. These poles are of the form of a numerator, which is the unintegrated numerator for the box or triangle diagram, divided by a denominator, which is the product of the propagators corresponding to the extra cuts in that diagram. For example, for the contribution of the box to the second triangle from the left in [Figure 4.3](#), the triangle has cuts for $(l_2 + K_1)^2 = 0$, $l_2^2 = 0$ and $(l_2 - K_2)^2 = 0$ and the box has an extra cut at $(l_2 - K_2 - K_3)^2 = 0$, so the pole contribution will be from “box numerator form”/ $(l_2 - K_2 - K_3)^2$. For this case the two reference vectors for the box are the same as the two reference vectors for the triangle, but the calculations are equivalent even if the choice of reference vectors is different, as long as care is made to use the correct signs on the box contributions and to use the vector for the extra pole evaluated relative to the appropriate loop momenta of the triangle.

The numerators of boxes can at most be linear in the free parameter in the triangle loop momentum parametrisation, as anything with higher powers would be cancellable with propagators in the box and therefore will be accounted for in triangle and bubble contributions. As its value at the location of the two box contributions are known, it must be given by

$$\text{“box numerator form”} = \frac{c(b^+ - b^-) + c^+b^- - c^-b^+}{c^+ - c^-}, \quad (4.43)$$

where c is whichever of c_1 and c_2 is the free parameter in the triangle parametrisation, $c^{+/-}$ are the two solutions to the extra box cut as given in [Equation 4.28](#) or [Equation 4.29](#) and $b^{+/-}$ are the two box coefficients corresponding to each of the solutions. Combining these relations, along with the $1/c$ from the Jacobian and simplifying

results in expressions for the poles of

$$\text{“pole”} = -\frac{1}{2K_3 \cdot n(c^+ - c^-)} \left(\frac{b^+}{c - c^+} - \frac{b^-}{c - c^-} \right), \quad (4.44)$$

where n is n_1 if c is c_1 and n is n_2 if c is c_2 .

Applying a triangle as a pole to a bubble, requires finding both the numerator structure and the simplified pole structure. The pole structure comes from evaluating the propagator corresponding to the extra cut at $l_2 - K_2$, using the momentum representation for l_2 given in Equation 4.32 which gives

$$\frac{1}{(l_2 - K_2)^2} = -\frac{tK_1 \cdot \chi}{K_2 \cdot n_2 K_1^2 (y^+ - y^-)} \left(\frac{1}{y - y^+} - \frac{1}{y - y^-} \right), \quad (4.45)$$

where $y^{+/-}$ are the two solutions of the propagator going on-shell for y which are given by

$$y^\pm = \left(\frac{1}{2} + t \frac{K_1 \cdot \chi K_1 \cdot K_2 - K_1^2 K_2 \cdot \chi}{K_2 \cdot n_2 K_1^2} \right) \left(1 \pm \sqrt{1 + \frac{t \frac{2tK_1 \cdot \chi K_2 \cdot n_1 + K_1^2 K_2 \cdot \chi + K_2^2 K_1 \cdot \chi}{K_1^2 K_2 \cdot n_2}}{\left(\frac{1}{2} + \frac{t(K_1 \cdot \chi K_1 \cdot K_2 - K_1^2 K_2 \cdot \chi)}{K_1^2 K_2 \cdot n_2} \right)^2}} \right). \quad (4.46)$$

Each c in the triangle form must come from a loop momentum multiplied with some combination of external momenta. Therefore, to expand out in terms of the space of the bubble momentum solutions, the values of the coefficients must be projected out of the full momentum. As the basis elements in the triangle momentum parametrisation have simple products with each other, an obvious choice for how to project out the coefficients, is to project using the basis vectors. The coefficients are found to be extracted by

$$c_1 = -\gamma \frac{l_2 \cdot n_2}{2\Delta} \quad c_2 = -\gamma \frac{l_2 \cdot n_1}{2\Delta}, \quad (4.47)$$

which if applied to the bubble momentum form will give

$$\begin{aligned} c_1(t, y) &= \frac{\gamma}{4\Delta} \left(\langle \tilde{K}_{t,2} | \tilde{K}_1 \rangle + \frac{(1-y) \langle \tilde{K}_{t,2} | \chi \rangle K_1^2}{2tK_1 \cdot \chi} \right) \left(t [\chi | \tilde{K}_{t,1}] + y [\tilde{K}_1 | \tilde{K}_{t,1}] \right) \\ c_2(t, y) &= \frac{\gamma}{4\Delta} \left(\langle \tilde{K}_{t,1} | \tilde{K}_1 \rangle + \frac{(1-y) \langle \tilde{K}_{t,1} | \chi \rangle K_1^2}{2tK_1 \cdot \chi} \right) \left(t [\chi | \tilde{K}_{t,2}] + y [\tilde{K}_1 | \tilde{K}_{t,2}] \right), \end{aligned} \quad (4.48)$$

where $\tilde{K}_{t,1/2}$ are the $\tilde{K}_{1/2}$ for the triangle by which this pole is caused. For triangles with a massless corner, whichever of these solutions corresponds to the variable that gives contributions for that triangle will be used. For the case where all corners are massive, both solutions will be used, as the solution which was not used to evaluate the triangle can be used to give an expression for $1/c$ that has no poles in terms of y . With these terms the all massive case can be evaluated, but in the massless triangle case an extra complication arises, as there are terms that are proportional to $\sqrt{1 + at + bt^2}$ where a and b are constants in terms of t and y . These terms integrate to zero on our contour, but would need an infinite number of terms in the Fourier projection to cause them to cancel correctly. For the massive case, due to symmetries, these terms cancel and so can be ignored, but for the massless case cancellations do not always occur and must be added in manually. These contributions come from partial fractioning of the pole form, which takes our numerator function $\text{num}(c(t, y))$ divided by the pole at $y = y^\pm$, and rearranges it to the form

$$\frac{C}{y^+ - y^-} \frac{\text{num}(c(t, y))}{y - y^\pm} = \frac{C}{y^+ - y^-} \frac{\text{num}(c(t, y^\pm))}{y - y^\pm} + \frac{Cf(t, y, y^\pm)}{y^+ - y^-}, \quad (4.49)$$

where C contains all the terms in the coefficient part of the expanded pole form that are free from y , y^+ and y^- and f is a rational function of its parameters, with only positive powers of y and y^\pm and both positive and negative powers of t . This function has terms that should vanish due to integrations over terms of the form $\sqrt{1 + at + bt^2}$, but which do not vanish in our numerical implementations and therefore must be subtracted. This expression will only give contributions that contain the dangerous term, when there are terms with an odd power of the square root. The simplest expression that would fully cancel this term is one which has exactly the same form, but with the opposite sign on the square root which corresponds to using the other solution for y . To directly evaluate f in terms of the other solution would require analytically calculating the form of f , which is a non trivial task, as such it is much easier to use the directly evaluated pole form to calculate the cancellation needed. This is a function of the form

$$\frac{C}{y^+ - y^-} \left(\frac{\text{num}(c(t, y))}{y - y^\pm} - \frac{\text{num}(c(t, y))}{y - y^\mp} + \frac{\text{num}(c(t, y^\mp))}{y - y^\mp} \right). \quad (4.50)$$

For poles from boxes contributing to bubbles, the product of the two extra prop-

agators and the known numerator form found above can be simplified using partial fractions to get a form for the poles that combines the box on triangle and triangle on bubble forms found above, giving

$$\begin{aligned} \frac{\text{“box numerator form”}}{(l_2 - K_2)^2 (l_2 - K_3)^2} &= -\frac{tK_1 \cdot \chi}{K_1^2} \left(\right. \\ &+ \frac{1}{K_2 \cdot n_2 (y_{2+} - y_{2-})} \left(\frac{\text{“pole”}(K_3, c(t, y_{2+}))}{y - y_{2+}} - \frac{\text{“pole”}(K_3, c(t, y_{2-}))}{y - y_{2-}} \right) \\ &+ \left. \frac{1}{K_3 \cdot n_2 (y_{3+} - y_{3-})} \left(\frac{\text{“pole”}(K_2, c(t, y_{3+}))}{y - y_{3+}} - \frac{\text{“pole”}(K_2, c(t, y_{3-}))}{y - y_{3-}} \right) \right), \quad (4.51) \end{aligned}$$

where $y_{2/3\pm}$ are the two solutions for y for the poles with K_2 or K_3 respectively and “pole”(K, c) is the box on triangle pole from Equation 4.44 for the extra cut at K evaluated for the value of c given. From this it is clear that the box on bubble pole gives a pair of contributions that each look like the contribution of a triangle and as such for evaluation they can be combined with the triangle pole that has the same pole structure.

For efficiency reasons it is useful to skip the calculation of terms that are known to not contribute. The more terms that can be skipped, the less work there is to do and the less terms there are at higher numbers of loop propagators, the less poles there are that need subtracting from lower loop propagator calculations. The simplest set of terms to ignore is any for which one of the corners is an amplitude with helicity choices that vanish, which applies irrespective of the particles involved, but has more cases to consider, as which helicity combinations vanish depends on the particles involved. The next few simplifications depend on properties of the loop momenta when corners are on-shell. Firstly, when a single corner goes on-shell, the loop momenta on each side of it can be simplified to

$$\begin{aligned} l_1^\mu &= \langle K_1 | \gamma^\mu \left(\left(1 + \frac{K_2^2}{2K_1 \cdot K_2} \right) |K_1\rangle + c_1 |\tilde{K}_2\rangle \right) \\ &= \left(\left(1 + \frac{K_2^2}{2K_1 \cdot K_2} \right) \langle K_1 | + c_2 \langle \tilde{K}_2 | \right) \gamma^\mu |K_1\rangle \\ l_2^\mu &= \langle K_1 | \gamma^\mu \left(\frac{K_2^2}{2K_1 \cdot K_2} |K_1\rangle + c_1 |\tilde{K}_2\rangle \right) \\ &= \left(\frac{K_2^2}{2K_1 \cdot K_2} \langle K_1 | + c_2 \langle \tilde{K}_2 | \right) \gamma^\mu |K_1\rangle, \quad (4.52) \end{aligned}$$

where the massless corner has been labelled as 1. From these expressions it is clear that whether c_1 or c_2 is non-zero, determines which type of spinor is proportional

for all the momenta in that corner. For amplitudes with three on-shell particles, only two cases are non vanishing and each of these must have two particles with one helicity and one with the opposite helicity and will be labelled by the helicity of which they have two. These amplitudes will each depend on only one type of spinor and therefore will vanish when the spinors they depend on are proportional. As a result, if a corner is massless, only one branch of the momentum solutions need be considered. Therefore, for numerical calculations, if there is at least one massless corner then one of them will be chosen to be corner 1 and only evaluated with the correct branch of the solution for the helicities in that corner.

If two neighbouring corners are on-shell then the loop momenta between them and either side of them are related in spinor forms and are given by

$$\begin{aligned}
l_1^\mu &= \langle K_1 | \gamma^\mu (|K_1] + c_1 |K_2]) \\
&= (\langle K_1 | + c_2 \langle K_2 |) \gamma^\mu |K_1] \\
l_2^\mu &= c_1 \langle K_1 | \gamma^\mu |K_2] \\
&= c_2 \langle K_2 | \gamma^\mu |K_1] \\
l_3^\mu &= (c_1 \langle K_1 | - \langle K_2 |) \gamma^\mu |K_2] \\
&= \langle K_2 | \gamma^\mu (c_2 |K_1] - |K_2]) ,
\end{aligned} \tag{4.53}$$

where spinors can now be written for K_1 and K_2 , as they are on-shell and so the \tilde{K}_1 and \tilde{K}_2 projected vectors are just the same as the unprojected vectors and so are no longer needed. If two adjacent corners are massless, then they must have opposite helicities, as otherwise they each vanish on opposite solutions and so neither solution will contribute for that diagram. For boxes, if two opposite corners are both on-shell, the solutions will have the same type of spinor proportional in each of the two corners and therefore must in both corners depend only on the other type of spinor and must be of the same helicity. These relations serve to greatly reduce the numbers of diagrams that are relevant and which need to be computed. For bubbles, an even stricter condition can be found as scalar loop diagrams vanish if there are no scales in the diagram and as there are only two corners, their momentum vectors must be equal and opposite, so the only scale available is the mass of this momentum. If it vanishes then the scalar loop function vanishes, so diagrams with only one massless particle on one side of the bubble will vanish and do not need to be computed. All

these relations apply equally well to amplitudes with a complex scalar, ϕ , as to pure QCD amplitudes, although care must be taken as a corner with contains only a ϕ is not a massless corner.

Contributions that are not relevant have been removed from the set of terms shown in [Figure 4.3](#), [Figure 4.4](#) and [Figure 4.5](#) using all these relations.

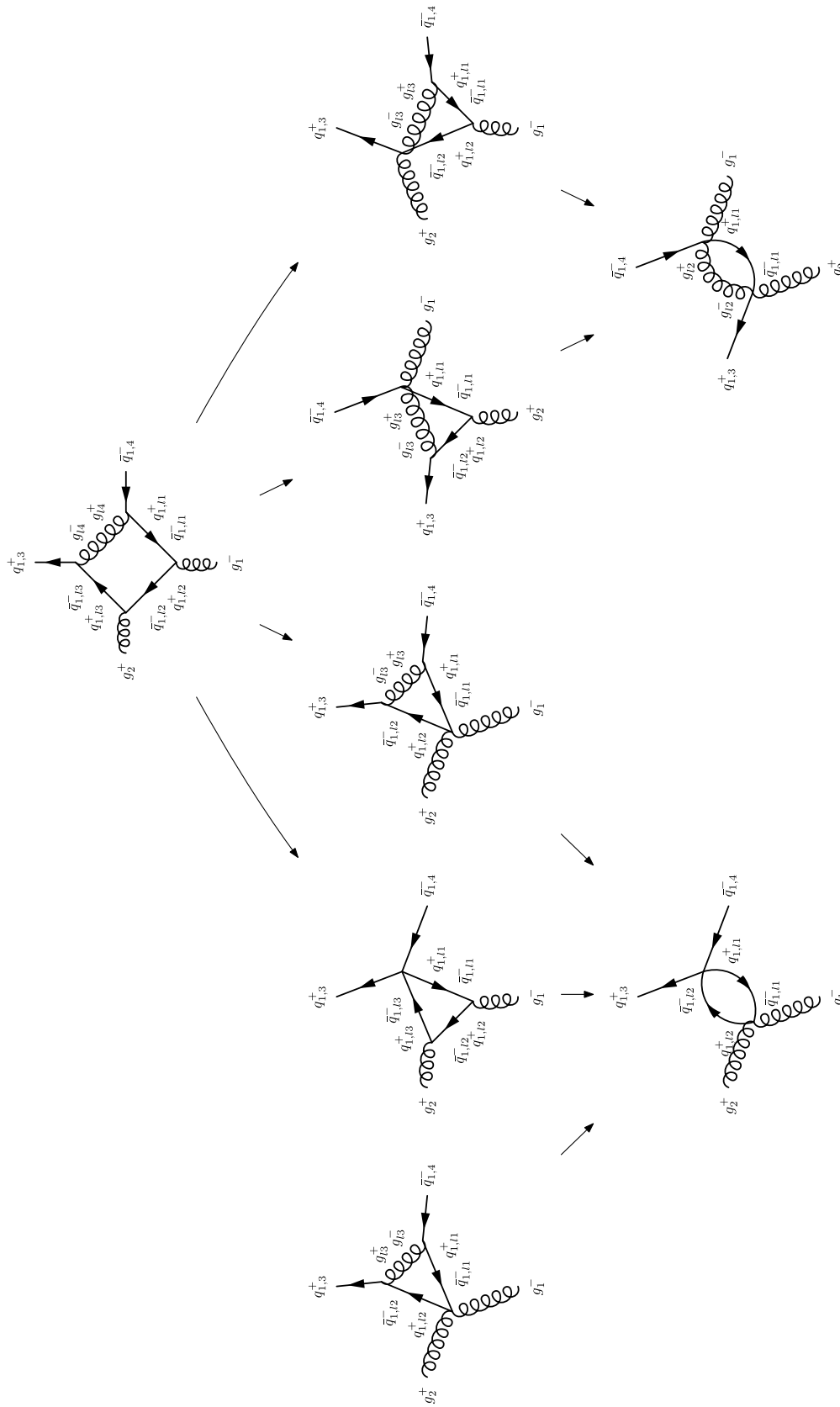
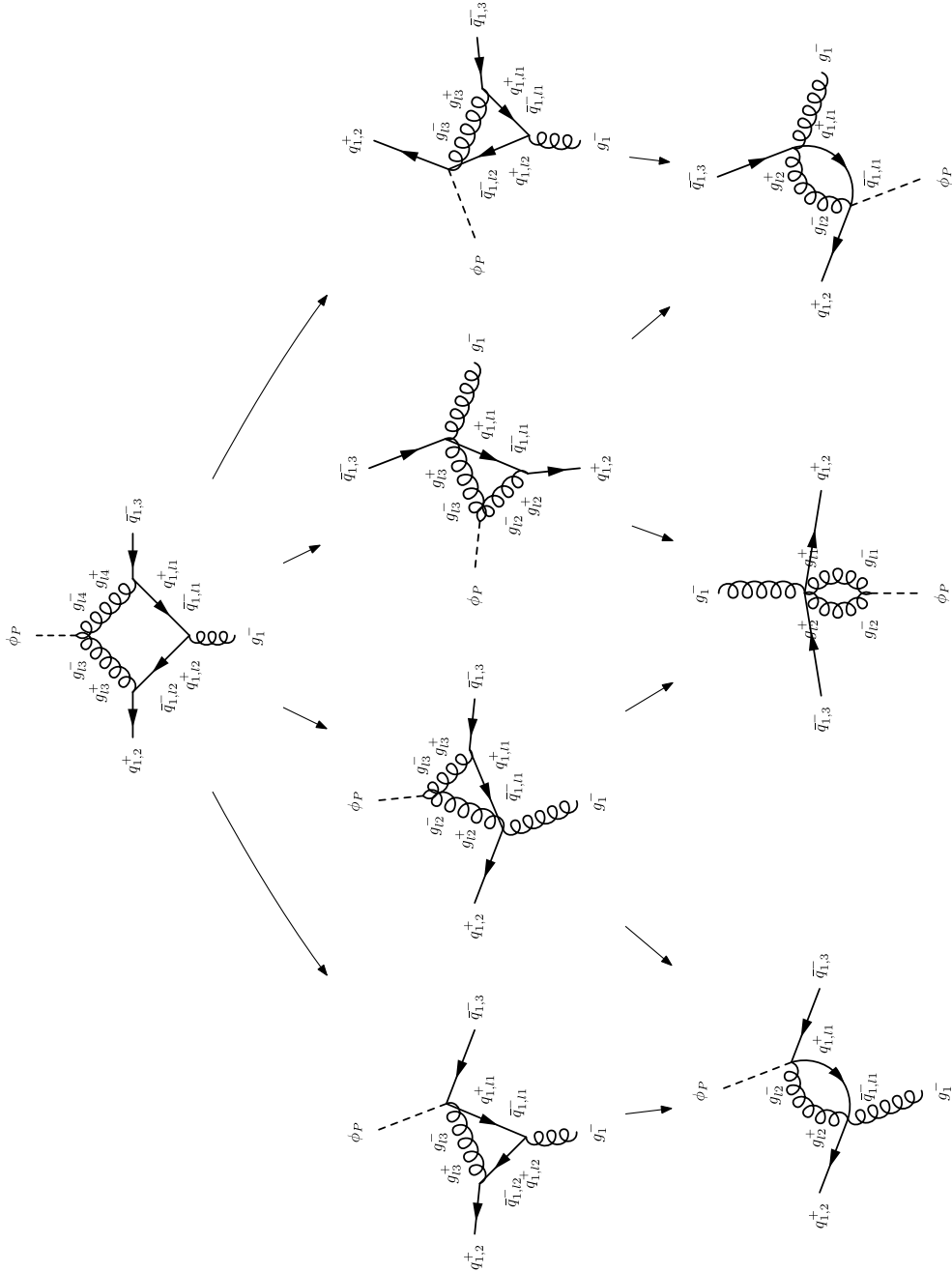
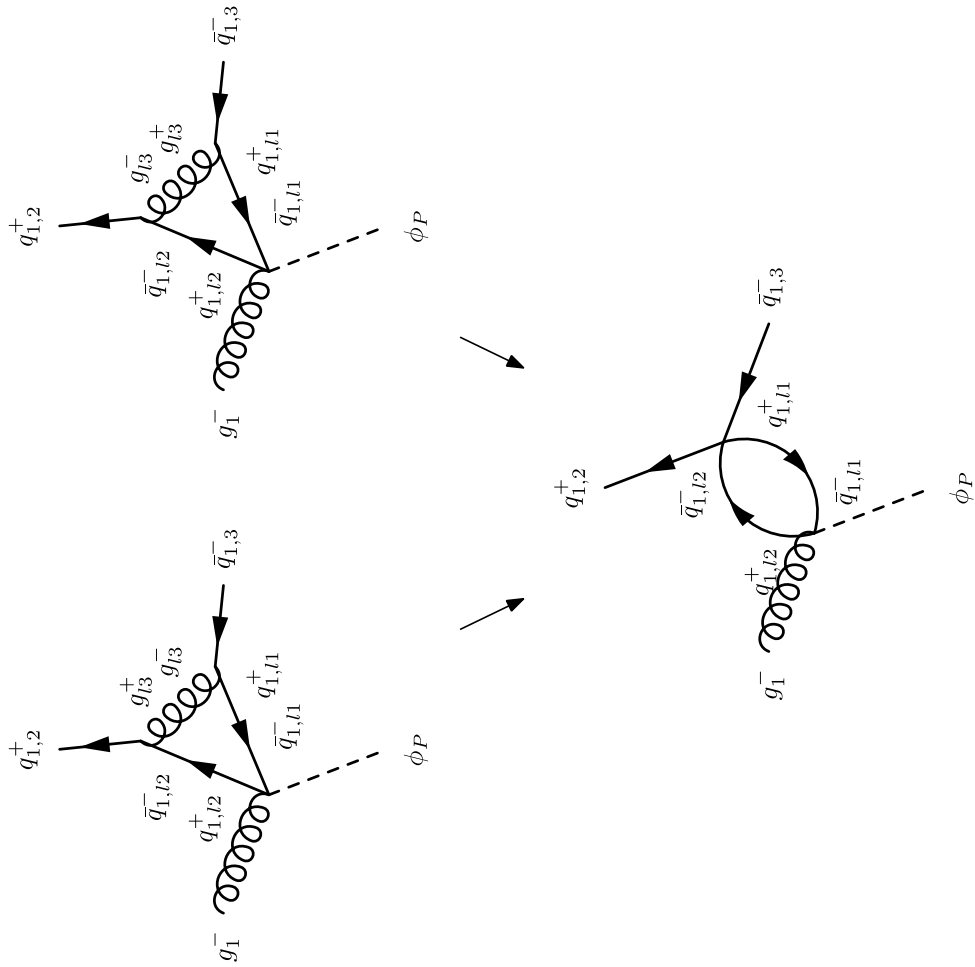


Figure 4.3: A graph showing the different terms that contribute to the amplitude for $g_1^- g_2^+ q_{1,3}^+ \bar{q}_{1,4}^-$ with the quark travelling left. The arrows indicate other diagrams to which a diagram contributes a pole.



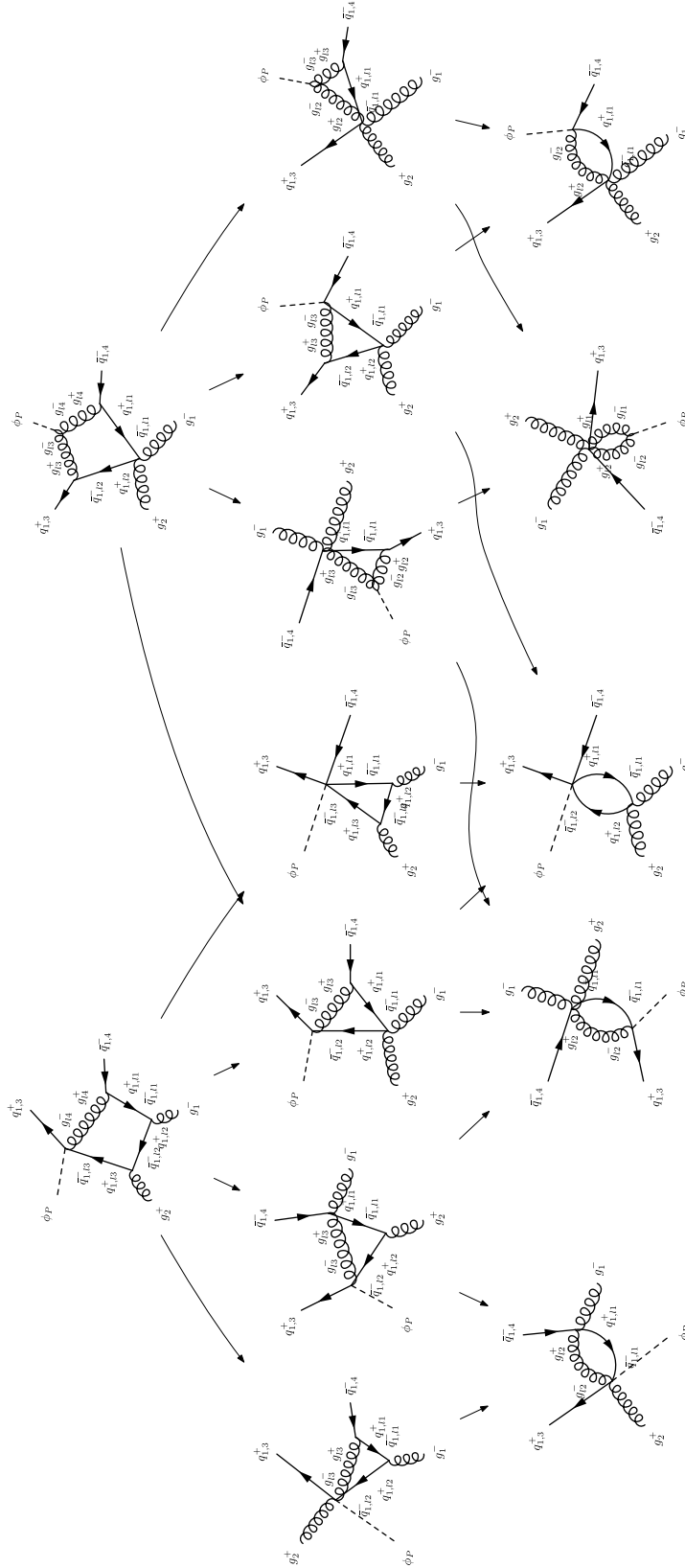
(4.4.a) One of the two fragments of the diagram of terms. The other is in [Figure 4.4.b](#).

Figure 4.4: A graph showing the different terms that contribute to the amplitude for $g_1^- q_{1,2}^+ \bar{q}_{1,3} \Phi_P$ with the quark travelling left. The arrows indicate other diagrams to which a diagram contributes a pole.



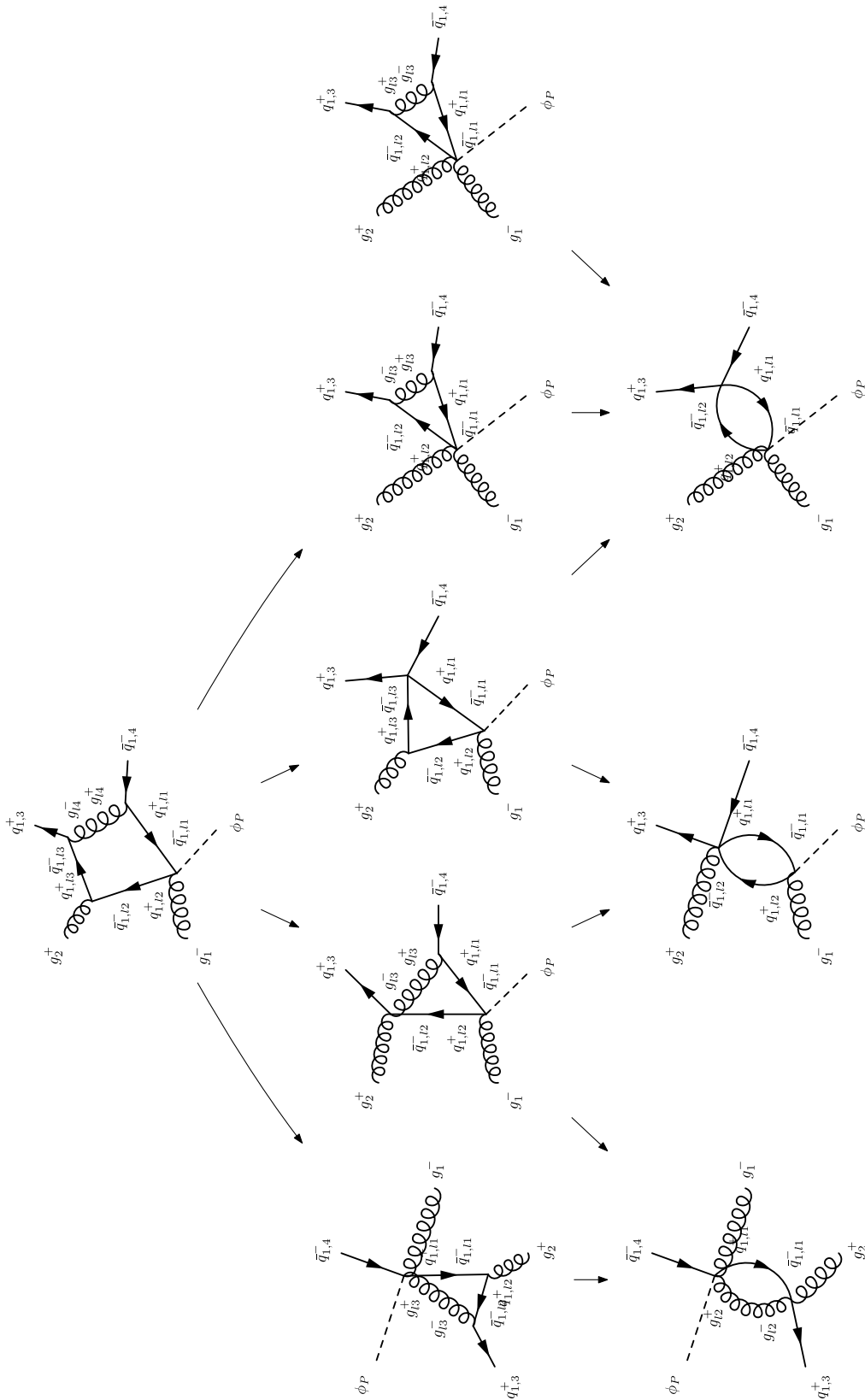
(4.4.b) One of the two fragments of the diagram of terms. The other is in [Figure 4.4.a](#).

Figure 4.4: A graph showing the different terms that contribute to the amplitude for $g_1^- q_{1,2}^+ \bar{q}_{1,3}^- \Phi_P$ with the quark travelling left. The arrows indicate other diagrams to which a diagram contributes a pole.



(4.5.a) One of the two fragments of the diagram of terms. The other is in Figure 4.5.b.

Figure 4.5: A graph showing the different terms that contribute to the amplitude for $g_1^- g_2^+ q_{1,3}^+ \bar{q}_{1,4}^- \Phi_P$ with the quark travelling left. The arrows indicate other diagrams to which a diagram contributes a pole.



(4.5.b) One of the two fragments of the diagram of terms. The other is in Figure 4.5.a.

Figure 4.5: A graph showing the different terms that contribute to the amplitude for $g_1^- g_2^+ q_{1,3}^+ \bar{q}_{1,4}^- \Phi_P$ with the quark travelling left. The arrows indicate other diagrams to which a diagram contributes a pole.

Chapter 5

Rational Terms

In the previous chapters, methods have been developed that will construct the majority of each amplitude for processes with a Higgs boson and many jets at one loop, however there are terms that cannot be extracted using just these techniques. These terms are the terms that do not have any poles or cuts in 4 dimensions and therefore cannot be constructed by techniques that require poles to build the amplitudes. These terms are called rational terms. One of the common ways to calculate rational terms is to use relations involving supersymmetric amplitudes[13]. These amplitudes contain combinations of the physical amplitudes and extra amplitudes. The most common combination to use is a combination of $\mathcal{N} = 1$ chiral and $\mathcal{N} = 4$ supersymmetric theories, where \mathcal{N} counts how many supersymmetries there are in the theory. For external gluon amplitudes these are given by[13]

$$\begin{aligned} A_n^{\mathcal{N}=4} &\equiv A_n^g + 4A_n^f + 3A_n^S \\ A_n^{\mathcal{N}=1 \text{ chiral}} &\equiv A_n^f + A_n^S, \end{aligned} \quad (5.1)$$

where A^g , A^f and A^S are the different versions of an amplitude, with gluons, Weyl fermions and complex scalars respectively in the loop. Using these building blocks pure gluon amplitudes and quark amplitudes can be built up as[13]

$$\begin{aligned} A_n^g &= A_n^{\mathcal{N}=4} - 4A_n^{\mathcal{N}=1 \text{ chiral}} + A_n^S \\ A_n^f &= A_n^{\mathcal{N}=1 \text{ chiral}} - A_n^S. \end{aligned} \quad (5.2)$$

As all supersymmetric amplitudes are cut constructable in the four dimensional

helicity scheme, they have no rational terms. Therefore, the rational terms of their component amplitudes are related to each other by[13]

$$A_n^g|_{\text{rational}} = A_n^S|_{\text{rational}} \qquad A_n^f|_{\text{rational}} = -A_n^S|_{\text{rational}} \ , \qquad (5.3)$$

where $|_{\text{rational}}$ means take only the rational part of the amplitude. From these relations it can be seen that the rational part of quark and gluon amplitudes is the same as the rational part of their related scalar amplitudes, but with a minus sign for quark amplitudes. These scalar amplitudes are much simpler and so can be extracted using generalised unitarity in 6 dimensions or recursion relations. Unfortunately this is not possible with Higgs boson amplitudes, as forming a supersymmetric amplitude would require using the supersymmetric extension for the Higgs boson, rather than the standard model Higgs boson.

Another type of method that has been used to extract these terms is recursion methods using ideas similar to BCFW as done by Berger, Del Duca and Dixon[15]. These methods use properties of the loop amplitudes under shifts of the external momenta to relate the rational terms to the cut terms and therefore rely on the pole structure of the cut terms when shifts are applied. These methods are therefore not as easy to apply numerically as they rely on complete knowledge of the pole structure of the amplitudes.

The method used here is to use generalised unitarity applied directly in more than 4 dimensions, as in more than 4 dimensions the rational terms have cuts and poles. If calculating in any specific dimensions greater than four, then the dependence on the dimensionality of space time can be extracted by using the dependence on the magnitude of the momenta pointing into the extra dimensions, as long as there is no explicit dependence on the direction or complex argument of any components. In the four dimensional helicity scheme (FDH)[3], which is being used here, all internal spinor and gluon states are kept in 4 dimensions. Therefore, the dependence of amplitudes on the dimension of space time, due to loop momenta and due to extra spinor and gluon states, must be separated. To derive the relations used to separate these two dependencies, it is useful to separate the dimension of the loop momenta, which will be called D_l , from the dimension of the spinor and gluon polarisation states, which

will be called D_s . Using this split, a loop amplitude is given by[14]

$$A_{D_l, D_s}(\{p_i\}) = \int \frac{d^{D_l} l}{i(\pi)^{\frac{D_l}{2}}} \frac{\mathcal{N}_{D_s}(\{p_i\}, l)}{d_1 d_2 \dots d_N} . \quad (5.4)$$

where D_l must be less than or equal to D_s .

If there is a closed quark loop in an amplitude, the extra contributions due to an increase of two in dimensions will be factors of two, due to there being twice as many states and the gamma matrices being twice as big. If chiral quarks are used, the quark space will split into two. If it can be shown that using either of the chiral quark subspaces is a valid representation of the 4 dimensional states and that amplitudes do not mix the two spaces and are identical in either case, then the state reduction can be performed by using only one of the two spaces in the calculations. For the case of pure gluon amplitudes, it can be shown that the dependence of the loop amplitude on D_s must be at most linear, as it arises purely from terms that form a closed loop of metric tensors. Therefore the amplitude in any dimensionality can be given by[14]

$$A_{D_l, D_s}(\{p_i\}) = A_{D_l}^0(\{p_i\}) + (D_s - 4) A_{D_l}^1(\{p_i\}) , \quad (5.5)$$

where $A_{D_l}^0(\{p_i\})$ and $A_{D_l}^1(\{p_i\})$ is independent of D_s , but will still depend on D_l . Combining two evaluations at two different values for D_s , the values of the two different components can be extracted which will allow a continuation to different values of D_s . Therefore, the value in the four dimensional helicity scheme is given by[14, 16]

$$A_{FDH} = \frac{(D_2 - 4)A_{D_l, D_s=D_1} - (D_1 - 4)A_{D_l, D_s=D_2}}{D_2 - D_1} . \quad (5.6)$$

To evaluate these expressions explicitly requires calculating in two different, even, integer dimensions, both of which must be greater than or equal to D_l , which is itself greater than 4. This would require working in both 6 and 8 dimensions although with the restriction that the loop momenta remains in either 5 or 6 dimensions. In both of these sets of dimensions a generalised unitarity calculation would need to be performed to extract the coefficients of the integrals and then once combined they can be limited back down to $d = 4 - 2\epsilon$ dimensions. Again the basis of loop integrals is needed and now the basis is extended to include pentagon scalar loop integrals. The coefficients now contain extra tensor components dependant on the loop momenta and there are now multiple elements that do not vanish when integrated. Some of these

$$A = \sum_i e_i \text{ (pentagon) } + \sum_i d_i \text{ (diamond) } + \sum_i c_i \text{ (triangle) } + \sum_i b_i \text{ (circle) } + \sum_i a_i \text{ (circle) }$$

Figure 5.1: The expansion of an amplitude in terms of the scalar basis functions in D dimensions. The sums are over each possible scalar loop function of that form, defined by the momenta at each corner. The coefficients contain loop momentum tensor structures and are therefore inside the loop momentum integral.

tensor terms give rise to D_l dependence which when combined with poles in the loop integrals results in the rational terms. In terms of this basis the amplitude is expanded as shown in [Figure 5.1](#). The forms that can appear in the coefficients are limited as any term containing external momenta can be replaced by a combination of inverse propagators, loop momenta squared and rational functions of external momenta as

$$l \cdot P = \frac{1}{2} (l^2 + P^2 - (l - P)^2) . \quad (5.7)$$

A further restriction is that there is no dependence on the direction in the extra $D - 4$ dimensions, as there are no external vectors in this subspace, only on the total magnitude of these components which will be labelled by μ^2 which is defined by

$$l^2 = \bar{l}^2 - \mu^2 , \quad (5.8)$$

where \bar{l} contains the 4 dimensional components of l . Therefore, the terms that can appear are all possible terms built from μ^2 and $t_i = l \cdot n_i$, where n_i are unit vectors that form an orthonormal basis of the subspace of 4 dimensional space that is transverse to all external momenta. A final restriction is on the highest power of loop momenta that can occur in the integrals, which for amplitudes involving a Higgs boson is one higher than the number of propagators in the loop momenta. Therefore, the general forms of the coefficients are

$$e_i = \mu^2 \tilde{e}_i \quad (5.9)$$

$$d_i = \bar{d}_i^0 + t_1 \bar{d}_i^1 + \mu^2 (\tilde{d}_i^0 + t_1 \tilde{d}_i^1) + \mu^4 (\tilde{d}_i^2 + t_1 \tilde{d}_i^3) \quad (5.10)$$

$$\begin{aligned}
c_i &= \bar{c}_i^0 + t_1 \bar{c}_i^1 + t_2 \bar{c}_i^2 + (t_1^2 - t_2^2) \bar{c}_i^3 + t_1 t_2 \bar{c}_i^4 + (t_1^2 - 3t_2^2) t_1 \bar{c}_i^5 + (t_2^2 - 3t_1^2) t_2 \bar{c}_i^6 \\
&+ (t_1^4 + t_2^4 - 6t_1^2 t_2^2) \bar{c}_i^7 + (t_1^2 - t_2^2) t_1 t_2 \bar{c}_i^8 \\
&+ \mu^2 (\tilde{c}_i^0 + t_1 \tilde{c}_i^1 + t_2 \tilde{c}_i^2 + (t_1^2 - t_2^2) \tilde{c}_i^3 + t_1 t_2 \tilde{c}_i^4) + \mu^4 \tilde{c}_i^5
\end{aligned} \tag{5.11}$$

$$\begin{aligned}
b_i &= \bar{b}_i^0 + t_1 \bar{b}_i^1 + t_2 \bar{b}_i^2 + t_3 \bar{b}_i^3 + (t_1^2 - t_3^2) \bar{b}_i^4 + (t_2^2 - t_3^2) \bar{b}_i^5 + t_1 t_2 \bar{b}_i^6 + t_1 t_3 \bar{b}_i^7 + t_2 t_3 \bar{b}_i^8 \\
&+ t_1 (t_1^2 - 3t_2^2) \bar{b}_i^9 + t_2 (t_2^2 - 3t_3^2) \bar{b}_i^{10} + t_3 (t_3^2 - 3t_1^2) \bar{b}_i^{11} + t_1 (t_2^2 - t_3^2) \bar{b}_i^{12} \\
&+ t_2 (t_3^2 - t_1^2) \bar{b}_i^{13} + t_3 (t_1^2 - t_2^2) \bar{b}_i^{14} + t_1 t_2 t_3 \bar{b}_i^{15} + \mu^2 (\tilde{b}_i^0 + t_1 \tilde{b}_i^1 + t_2 \tilde{b}_i^2 + t_3 \tilde{b}_i^3) ,
\end{aligned} \tag{5.12}$$

where the dependence on t_i encodes all the 4 dimensional loop momentum dependence, coefficients with a bar over them appear in 4 dimensional generalised unitarity calculations, coefficients with a $\tilde{}$ over them are new coefficients for D dimensional calculations and all coefficients are independent of the dimension of the loop momenta, D_l . It is now possible to expand each of these coefficients out into their own integrals. Most of the integrals vanish and as such only a small subset are needed directly, but for numerical calculations, in theory, all coefficients are needed as subtractions. The only integrals, other than the normal 4 dimensional basis, that do not vanish in the limit of $D = 4 - 2\epsilon \rightarrow 4$ are integrals whose numerators can only be powers of μ^2 . However the box and pentagon with μ^2 inserted do vanish in this limit. In the limit of $D = 4 - 2\epsilon \rightarrow 4$, dropping terms of order ϵ , the extra non-vanishing integrals are given by[17]

$$I_4[\mu^4] = \int \frac{d^D l}{i\pi^{\frac{D}{2}}} \frac{\mu^4}{D_1 D_2 D_3 D_4} = -\frac{1}{6} \tag{5.13}$$

$$I_3[\mu^4] = \int \frac{d^D l}{i\pi^{\frac{D}{2}}} \frac{\mu^4}{D_1 D_2 D_3} = \frac{1}{24} (P_1^2 + P_2^2 + P_3^2) \tag{5.14}$$

$$I_3[\mu^2] = \int \frac{d^D l}{i\pi^{\frac{D}{2}}} \frac{\mu^2}{D_1 D_2 D_3} = \frac{1}{2} \tag{5.15}$$

$$I_2[\mu^2] = \int \frac{d^D l}{i\pi^{\frac{D}{2}}} \frac{\mu^2}{D_1 D_2} = -\frac{1}{6} P_1^2 = -\frac{1}{6} P_2^2 , \tag{5.16}$$

where the $I_n[X]$ is the scalar integral with n propagators and a numerator term of X , D_i is the i^{th} propagator and P_i is the momenta flowing out of the i^{th} corner. Combining these with the coefficients above, the rational term is given by

$$R = -\frac{1}{6} \sum_i \bar{a}_i^2 + \frac{1}{2} \sum_i \left(\frac{P_{i,1}^2 + P_{i,2}^2 + P_{i,3}^2}{12} \tilde{c}_i^5 + \tilde{c}_i^0 \right) - \frac{1}{6} \sum_i P_{i[1]}^2 \tilde{b}_i^0 , \tag{5.17}$$

where $P_{i,j}$ is the momenta in the j^{th} corner of the cut i .

To calculate the amplitudes in the four dimensional helicity scheme, the relevant coefficients must be extracted using the methods from the previous chapter and combined to form the rational terms in both 6 and 8 dimensions. The rational terms in each dimension are then combined using [Equation 5.6](#) to form the rational term in the four dimensional helicity scheme. Alternatively, as the relations are linear, the extrapolation of the spinor and polarisation state dimension dependence to 4 dimensions can be done at the coefficient level and then only combined to give the rational term once the FDH versions of each coefficient are found.

Working in higher numbers of dimensions increases the complexity of the calculations, as such it is always simpler to work in fewer dimensions and a lower number of different dimensions. It is possible to avoid working in 8 dimensions by simplifying the calculation so only a single dimension greater than 4 is needed by making use of the form of the Feynman rules, as is done by Giele, Kunszt and Melnikov[14] and Davies[16]. The single dimension now needed will be taken to be 6 dimensions. This is done by noticing that if D_l is taken as being 5, then D_1 and D_2 can be taken to be 5 and 6 respectively. Furthermore, if the amplitude is purely gluons then the difference between working in D_1 and D_2 , will be the contribution due to gluons with their polarisation vectors pointing in the 6th dimension, which, if the Feynman rules can be extracted consistently, is equivalent to a complex scalar particle[14, 16]. The new Feynman rules needed for pure gluon amplitudes are just the normal three and four gluon vertex factors, but with one of the gluons forced to be in the 6th dimension, along with the propagator factor for gluons in the 6th dimension and gluon polarisation vectors in the 6th dimension. The polarisation vector must be given by

$$\epsilon_6^\mu = n_6^\mu , \quad (5.18)$$

where n_6^μ is the unit vector in the 6th dimension. On contracting this with a three gluon vertex it is clear that the combination will vanish, unless exactly one of the other gluons is also polarised in the direction of the 6th dimension. Therefore, taking two of the gluons, with momenta k_1 and k_2 , as being polarised in the 6th dimension, the vertex factor is given by

$$-\frac{1}{\sqrt{2}}(k_1 - k_2)^\sigma n_6^\mu n_6^\nu , \quad (5.19)$$

where the σ index is contracted with the remaining gluon line and the indices μ and ν are contracted with gluons pointing in the 6th dimension. The same logic applies to the four gluon vertex, which requires that if any gluons are polarised in the 6th dimension, either two or four are. If two neighbouring gluons are polarised in the 6th dimension the four gluon vertex is given by

$$-\frac{i}{2}g^{\sigma\rho}n_6^\mu n_6^\nu, \quad (5.20)$$

where the indices μ and ν are contracted with the gluons polarised in the 6th dimension and the indices σ and ρ are contracted with the two remaining gluons which are next to each other. It is clear from the form of the gluon propagator that if the gluon at one end is polarised in the 6th dimension, the other end will also be, which allows it to be simplified to

$$\frac{in_6^\mu n_6^\nu}{k^2 + i0}. \quad (5.21)$$

For the Higgs boson calculations the same extraction must be applied to the Higgs boson vertex rules. These two vertex rules obey the same relations as above and therefore will be given for the case of two gluons being in the direction of the 6th dimension. The vertex rule from the 2 gluon vertex is given by

$$-2ik_1 \cdot k_2 n_6^\mu n_6^\nu, \quad (5.22)$$

where k_1 and k_2 are the momenta of the gluons polarised in the 6th dimension which are contracted with the indices μ and ν . The vertex rule derived from the 3 gluon vertex is given by

$$-i\sqrt{2}(k_1 - k_2)^\sigma n_6^\mu n_6^\nu, \quad (5.23)$$

where k_1 and k_2 are the momenta of the gluons polarised in the 6th dimension which are contracted with the indices μ and ν and the remaining index, σ , is contracted with the remaining gluon.

From these relations it is clear that for the tree amplitudes in any corner of a generalised unitarity calculation, as there will be exactly two external gluons polarised in the 6th dimension, there must be a single line through the diagrams of gluons polarised in this direction and these gluons can all be treated as scalars using the new Feynman rules calculated above with the factors of n_6^μ stripped. The Feynman rules

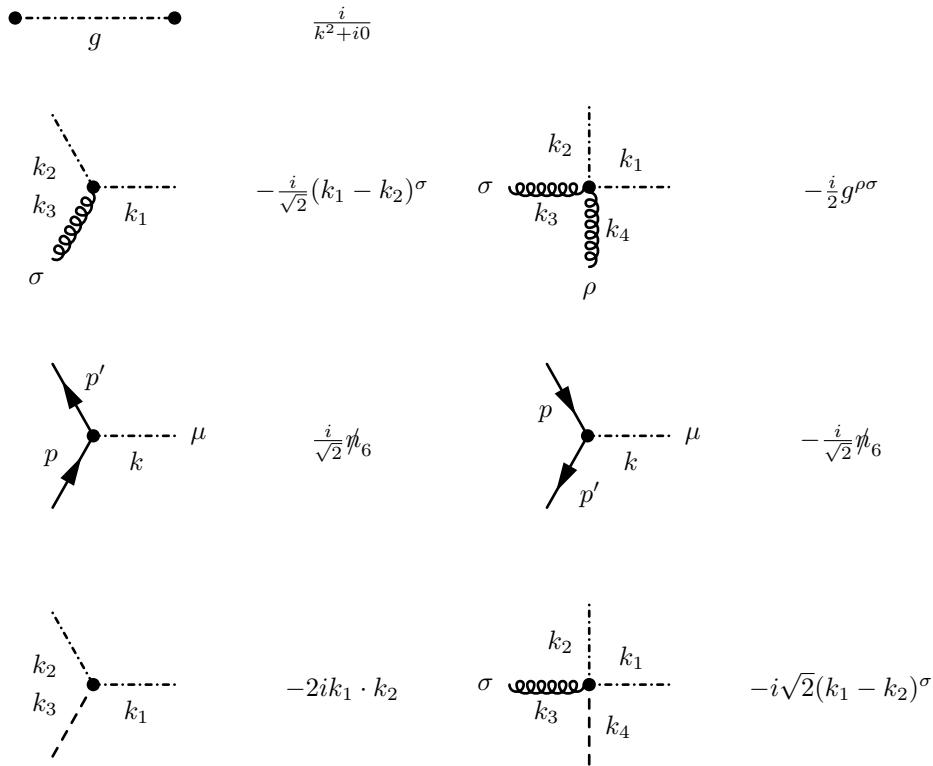


Table 5.1: Colour ordered Feynman rules in Faddeev-Popov gauge for the scalar particle equivalent to a gluon polarised in the 6th dimension. The dash-dotted lines are the new scalar and all momenta are inbound.

for the new complex scalar are given in [Table 5.1](#).

The last case to handle is amplitudes with external quarks which enter the loop, which can have a mixture of quarks and gluons as the cut particles. This case can be handled by working with only one of the two copies of 4 dimensional spinors, to reduce the quark state dimension dependence, and then using scalar 6th dimensional gluons to subtract the extra gluon polarisation states. This is possible, as after reducing to only the 4 dimensional set of spinors, the dependence on D_s is due to terms with a closed loop of metric tensors and pairs of gamma matrices in a quark line, which again can result in at most linear dependence on D_s . The extra Feynman rules needed for this case are also shown in [Table 5.1](#).

Therefore, if the calculations for the full 6 dimensional amplitudes and, where there are gluons in the loop, the subtraction terms using the scalar equivalent for the gluon, can be performed, the coefficients of scalar loop integrals in the amplitude can be extracted and then the rational terms can be calculated using only 6 dimensions.

To perform these calculations, an explicit numerical implementation must be derived and must be checked to ensure the properties used above are true. These aspects will be investigated in the next few sections.

5.1 6 Dimensional Spinor Helicity Formalism

The 6 dimensional extension of the spinor helicity formalism will enable efficient calculation of amplitudes in 6 dimensions by making use of the many cancellations and vanishing products, especially as many of the vectors are still purely 4 dimensional. To derive 6 dimensional spinors, a 6 dimensional generalisation of the gamma matrices is needed. The conditions these have to obey, as for the 4 dimensional case, are

$$\{\gamma^\mu, \gamma^\nu\} = g^{\mu\nu} , \quad (5.24)$$

where $g^{\mu\nu} = \text{diag}(1, -1, -1, \dots)$ is the metric of space time. There are many solutions to these equations but some are more convenient to use than others. This project uses a recursive definition, so that the different dimensions can be compared directly. In addition, a basis is used where as many elements as possible are zero, so that as many elements as possible in both spinors and products are zero and different elements mix as little as possible in products. The basis is also chosen to be based on the Weyl (chiral) basis as helicity spinors will be used and therefore a diagonal chirality operator, the γ_5 equivalent, will ensure different helicities do not mix. The chirality operator from now on will be referred to as γ_C , as γ_5 is the name of a normal gamma matrix in 6 or more dimensions. The basis chosen is[18]

$$\begin{aligned} \gamma_d^\mu &= \gamma_{d-2}^\mu \otimes \sigma_3 & \gamma_d^{d-2} &= iI_{d-2} \otimes \sigma_1 & \gamma_d^{d-1} &= -iI_{d-2} \otimes \sigma_2 \\ \gamma_2^0 &= \sigma_1 & \gamma_2^1 &= -i\sigma_2 , \end{aligned} \quad (5.25)$$

where I_d is the identity matrix in d dimensions which is given by

$$I_d = I_2 \otimes I_{d-2} \quad I_2 = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix} , \quad (5.26)$$

and the Pauli matrices, σ_i , are given by

$$\sigma_1 = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \quad \sigma_2 = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix} \quad \sigma_3 = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}. \quad (5.27)$$

Using these definitions the 6 dimensional gamma matrices are given by

$$\gamma_6^0 = \begin{pmatrix} 0 & 1 & & & & & & \\ & 1 & 0 & & & & & \\ & & 0 & -1 & & & & \\ & & -1 & 0 & & & & \\ & & & & 0 & -1 & & \\ & & & & -1 & 0 & & \\ & & & & & & 0 & 1 \\ & & & & & & 1 & 0 \end{pmatrix} \quad \gamma_6^1 = \begin{pmatrix} 0 & -1 & & & & & & \\ 1 & 0 & & & & & & \\ & & 0 & 1 & & & & \\ & & 0 & -1 & 0 & & & \\ & & & & 0 & 1 & & \\ & & & & -1 & 0 & & \\ & & & & & & 0 & -1 \\ & & & & & & 1 & 0 \end{pmatrix}$$

$$\gamma_6^2 = \begin{pmatrix} 0 & i & & & & & & \\ i & 0 & & & & & & \\ & & 0 & -i & & & & \\ & & -i & 0 & & & & \\ & & & & 0 & 1 & & \\ & & & & 1 & 0 & & \end{pmatrix} \quad \gamma_6^3 = \begin{pmatrix} 0 & 1 & & & & & & \\ -1 & 0 & & & & & & \\ & & 0 & -1 & & & & \\ & & -1 & 0 & & & & \\ & & & & 0 & 1 & & \\ & & & & 1 & 0 & & \end{pmatrix} \quad \gamma_6^4 = \begin{pmatrix} 0 & i \\ i & 0 \end{pmatrix} \quad \gamma_6^5 = \begin{pmatrix} 0 & 1 \\ -1 & 0 \end{pmatrix}, \quad (5.28)$$

where the matrices are given in block form and are all 8×8 matrices. Using this basis the γ_C is given by

$$\gamma_C = \begin{pmatrix} 1 & 0 & & & & & & \\ & 0 & -1 & & & & & \\ & & & -1 & 0 & & & \\ & & & 0 & 1 & & & \\ & & & & & -1 & 0 & \\ & & & & & 0 & 1 & \\ & & & & & & 1 & 0 \\ & & & & & & 0 & -1 \end{pmatrix}. \quad (5.29)$$

In hindsight, a basis that produced a chirality operator that was diagonal, with the top four values being one and the bottom four values being minus one, would have made the calculations simpler, as splitting the space into two using the helicity projector would have split spinors and matrices in half, top from bottom, whereas with this basis the space is split in a more complex pattern of rows, 1, 4, 6, 7 and 2, 3, 5, 8. However, as this only makes a difference to readability and simplicity of algebraic implementations, the basis defined in [Equation 5.25](#) is used from here on.

Using this basis, spinors can be derived by using the defining equation

$$\not{p}u(p) = 0 , \quad (5.30)$$

where $u(p)$ is a spinor corresponding to the momentum p . After using this definition there are still four degrees of freedom left. One is removed by the normalisation condition of

$$\sum u(p)\bar{u}(p) = \not{p} , \quad (5.31)$$

where \bar{u} is the conjugate spinor for u and for real momenta is defined by $\bar{u} = u^\dagger\gamma_0$. The others are used to split the general spinor into multiple spinor states. The first split is into the eigen spaces of the chirality operator, γ_C , which splits the spinor space into two subspaces, each with one remaining degree of freedom. So that the external particles, which are still 4 dimensional, have simple spinors, the last split should reduce to the eigen states of the 4 dimensional chirality operator, $\gamma_{6,4C}$, when the momenta is 4 dimensional. Unfortunately it is not possible to split the remaining spinor space by the 4 dimensional helicity operator for full 6 dimensional momenta. For this project, one of the simplest extensions away from the 4 dimensional limit was chosen, which is to keep the components that are already non-zero the same and only allow one extra component to become non-zero. Any choice for both gamma matrices and spinors should produce the same answers, other than an overall phase change on amplitudes, but this has not been tested directly as it would require reproducing the complete calculations for multiple choices of spinor basis.

Using these choices the spinors are given by

$$\begin{aligned}
u_{++} &= \begin{pmatrix} \sqrt{p_0 + p_1} \\ 0 \\ 0 \\ \frac{p_3 - ip_2}{\sqrt{p_0 + p_1}} \\ 0 \\ \frac{p_5 - ip_4}{\sqrt{p_0 + p_1}} \\ 0 \\ 0 \end{pmatrix} & u_{-+} &= \begin{pmatrix} 0 \\ \frac{p_5 + ip_4}{\sqrt{p_0 + p_1}} \\ 0 \\ 0 \\ \sqrt{p_0 + p_1} \\ 0 \\ 0 \\ \frac{p_3 - ip_2}{\sqrt{p_0 + p_1}} \end{pmatrix} & u_{+-} &= \begin{pmatrix} 0 \\ 0 \\ 0 \\ -\frac{p_5 + ip_4}{\sqrt{p_0 + p_1}} \\ 0 \\ \frac{p_3 + ip_2}{\sqrt{p_0 + p_1}} \\ \sqrt{p_0 + p_1} \\ 0 \end{pmatrix} & u_{--} &= \begin{pmatrix} 0 \\ \frac{p_3 + ip_2}{\sqrt{p_0 + p_1}} \\ \sqrt{p_0 + p_1} \\ 0 \\ 0 \\ 0 \\ 0 \\ -\frac{p_5 - ip_4}{\sqrt{p_0 + p_1}} \end{pmatrix} .
\end{aligned} \tag{5.32}$$

As in 4 dimensions, it is possible to write the conjugate spinors in terms of a combination of the transposed spinors. For this choice of spinors, the conjugates are given by

$$\bar{u}_{h-l} = i u_{hl}^T \sigma_1 \otimes \sigma_2 \otimes \sigma_1 \equiv \underline{u}_{hl} , \tag{5.33}$$

where the underlined spinor, \underline{u}_{hl} , is the conjugate spinor, but is labelled by the helicity of the spinor it can be written in terms of and $i\sigma_1 \otimes \sigma_2 \otimes \sigma_1$ is a real anti-diagonal matrix. For example \bar{u}_{++} is given by

$$\begin{aligned}
\bar{u}_{++} &= u_{++}^\dagger \gamma_0 \\
&= \begin{pmatrix} \sqrt{p_0 + p_1} & 0 & 0 & \frac{p_3 + ip_2}{\sqrt{p_0 + p_1}} & 0 & \frac{p_5 + ip_4}{\sqrt{p_0 + p_1}} & 0 & 0 \end{pmatrix} \gamma_0 \\
&= \begin{pmatrix} 0 & \sqrt{p_0 + p_1} - \frac{p_3 + ip_2}{\sqrt{p_0 + p_1}} & 0 & -\frac{p_5 + ip_4}{\sqrt{p_0 + p_1}} & 0 & 0 & 0 & 0 \end{pmatrix} ,
\end{aligned} \tag{5.34}$$

and \underline{u}_{+-} is given by

$$\begin{aligned}
\underline{u}_{+-} &= -i u_{+-}^T \sigma_1 \otimes \sigma_2 \otimes \sigma_1 \\
&= -i \begin{pmatrix} 0 & 0 & 0 & -\frac{p_5 + ip_4}{\sqrt{p_0 + p_1}} & 0 & \frac{p_3 + ip_2}{\sqrt{p_0 + p_1}} & \sqrt{p_0 + p_1} & 0 \end{pmatrix} \sigma_1 \otimes \sigma_2 \otimes \sigma_1 \\
&= \begin{pmatrix} 0 & \sqrt{p_0 + p_1} - \frac{p_3 + ip_2}{\sqrt{p_0 + p_1}} & 0 & -\frac{p_5 + ip_4}{\sqrt{p_0 + p_1}} & 0 & 0 & 0 & 0 \end{pmatrix} .
\end{aligned} \tag{5.35}$$

This shows that $\underline{u}_{+-} = \bar{u}_{++}$ as expected.

The spinor products for each helicity combination can be worked out, both in general from the relations and in the specific case of these spinors. The spinor products

that are exactly zero are $\underline{u}_{+h_1}(p)u_{+h_2}(q)$ and $\underline{u}_{-h_1}(p)u_{-h_2}(q)$, where p and q are two general on-shell 6 dimensional momenta and h_1 and h_2 are arbitrary signs, which can be shown using the definition of the spinors. As the spinors are eigenstates of the 6 dimensional helicity operator, the appropriate sign projection operator,

$$P_{\pm} = \frac{1 \pm \gamma_C}{2}, \quad (5.36)$$

can be applied to each spinor. When applied to spinors and conjugate spinors the projection operators satisfy the relations

$$\begin{aligned} P_{\pm}u_{\pm,h} &= u_{\pm,h} & P_{\pm}u_{\mp,h} &= 0 \\ \underline{u}_{\mp,h}P_{\pm} &= \underline{u}_{\mp,h} & \underline{u}_{\pm,h}P_{\pm} &= 0. \end{aligned} \quad (5.37)$$

If the projection operators introduced for the two spinors in a spinor product have opposite signs then they will cancel and the product will vanish.

In general, the other spinor products are non-zero, however a similar logic to that used above, but using the 4 dimensional chirality operator, shows that in the limit of both momenta being 4 dimensional $\underline{u}_{h_1h_2}(p)u_{h_1h_2}(q)$ vanishes. The remaining spinor products, in this limit, must be given by the standard 4 dimensional spinor products, apart from possible changes in sign. For the choice of spinors used in this project, various spinor products can be related to each other. The full set of relations and their limits in the case of 4 dimensional momenta are given by,

$$\begin{aligned} \underline{u}_{++}(p)u_{--}(q) &= -\underline{u}_{--}(q)u_{++}(p) = \underline{u}_{--}(p)u_{++}(q) = -\underline{u}_{++}(q)u_{--}(p) \xrightarrow{p,q_{4,5} \rightarrow 0} 0 \\ \underline{u}_{-+}(p)u_{+-}(q) &= -\underline{u}_{+-}(q)u_{-+}(p) = \underline{u}_{+-}(p)u_{-+}(q) = -\underline{u}_{-+}(q)u_{+-}(p) \xrightarrow{p,q_{4,5} \rightarrow 0} 0 \\ \underline{u}_{++}(p)u_{-+}(q) &= \underline{u}_{-+}(q)u_{++}(p) = -\underline{u}_{-+}(p)u_{++}(q) = -\underline{u}_{++}(q)u_{-+}(p) = -\underline{u}_{-+}(p)u_{+}(q) \\ \underline{u}_{--}(p)u_{+-}(q) &= \underline{u}_{+-}(q)u_{--}(p) = -\underline{u}_{+-}(p)u_{--}(q) = -\underline{u}_{--}(q)u_{+-}(p) = -\underline{u}_{-}(p)u_{-}(q), \end{aligned} \quad (5.38)$$

where the spinors with only one sign are 4 dimensional spinors as given in [Equation 2.19](#). In [Section 5.2](#) even simpler explicit expressions for these products in terms of massless 4 dimensional projections will be derived.

The other main contribution needed for calculating amplitudes in 6 dimensions is a representation for polarisation vectors of gluons. As for the spinors, a representa-

tion that reduces to their 4 dimensional versions would be helpful as it would allow easy comparison to 4 dimensional amplitudes and calculations. Unfortunately, the simplest option of using the same expression as is used in 4 dimensions, as shown in Equation 2.21, does not obey the relations expected for polarisation vectors when used in 6 dimensions. It is possible to find a simple extension to the 4 dimensional case which does obey all the required relations, which is given by

$$\epsilon_{h,l,-h,m}^\mu(p; k) = \frac{\underline{u}_{hl}(p)\gamma^\mu \not{k} \underline{u}_{-hm}}{2\sqrt{2}p \cdot k} , \quad (5.39)$$

where k is again a arbitrary on-shell vector that fixes the gauge for the gluon. This representation though does have the issue of there being eight different states where as physically there are known to be four gluon states. This issue is resolved by noticing that the states form pairs that are equivalent, up to a choice of sign, related by

$$\epsilon_{+,h,-,l}(p; k) = -hl\epsilon_{-,l,+,h}(p; k) . \quad (5.40)$$

Therefore, where convenient, the polarisation vector form will be simplified to

$$\epsilon_{h,l}(p; k) = \epsilon_{+,h,-,l}(p; k) . \quad (5.41)$$

The explicit mention of the arbitrary reference vector will also be dropped where there is only one polarisation vector and will be implicitly labelled as k .

The other expression needed when calculating amplitudes and expressions involving gluons is a formula for the contraction of a polarisation vector with a gamma matrix, either in another polarisation vector or in a spinor chain. This can be shown to be given by

$$\not{\epsilon}_{hl}(p) = \frac{\not{k} \underline{u}_{-,l}(p) \underline{u}_{+,h}(p) - \underline{u}_{-,l}(p) \underline{u}_{+,h}(p) \not{k} + hl (\not{k} \underline{u}_{+,h}(p) \underline{u}_{-,l}(p) - \underline{u}_{+,h}(p) \underline{u}_{-,l}(p) \not{k})}{\sqrt{2}p \cdot k} , \quad (5.42)$$

where, when combined with any spinor chain or polarisation vector, at most two of the four terms will contribute.

5.2 Reducing 6 Dimensional Spinors to 4 Dimensions

A 6 dimensional vector can be decomposed into a 4 dimensional vector and a vector purely in the 5th and 6th dimensions,

$$p^\mu = p_{4D,M}^\mu + p_{6D}^\mu, \quad (5.43)$$

where p is the full 6 dimensional vector, $p_{4D,M}$ is the 4 dimensional part and p_{6D} is the part purely in the 5th and 6th dimensions. In components the vectors are given by

$$p = \begin{pmatrix} p_0 \\ p_1 \\ p_2 \\ p_3 \\ p_4 \\ p_5 \end{pmatrix} \quad p_{4D,M} = \begin{pmatrix} p_0 \\ p_1 \\ p_2 \\ p_3 \\ 0 \\ 0 \end{pmatrix} \quad p_{6D} = \begin{pmatrix} 0 \\ 0 \\ 0 \\ 0 \\ p_4 \\ p_5 \end{pmatrix}. \quad (5.44)$$

Both $p_{4D,M}$ and p_{6D} are in general massive vectors and their dot product is zero as they have no components in common directions. Using an extra arbitrary massless 4 dimensional vector, a , it is possible to write the 4 dimensional part in terms of a massless 4 dimensional vector as

$$p_{4D}^\mu = p_{4D,M}^\mu - a^\mu \frac{p_{4D,M}^2}{2p_{4D,M} \cdot a}. \quad (5.45)$$

The vector p_{6D} can be written in terms of any pair of 4 dimensional vectors and the values of the 5th and 6th dimensional components. If the two vectors are chosen to be the vectors a and p_{4D} , p_{6D} can be written as

$$p_{6D}^\nu = hl \frac{\mu_p^{hl} \underline{u}_{hl}(p_{4D}) \gamma^\nu \not{p}_{-h-l}(p_{4D}) - \mu_p^{-hl} \underline{u}_{h-l}(p_{4D}) \gamma^\nu \not{p}_{hl}(p_{4D})}{4p_{4D} \cdot a}, \quad (5.46)$$

where h and l are arbitrary signs and $\mu_p^\pm = p_5 \pm ip_4$ contains the values of the 5th and 6th dimensional components of the vector p .

The spinors for the full 6 dimensional vector, p , can also be written in terms of the spinors for their massless 4 dimensional projection and the arbitrary vector

used in the massless projection by inserting the fraction $\frac{\not{p}_{4D} + \not{p}_{4D} \not{a}}{2p_{4D} \cdot a}$, which due to the commutation properties of gamma matrices is equivalent to the identity, and gives

$$\begin{aligned}
u_{h,l}(p) &= \frac{\not{p}_{4D} + \not{p}_{4D} \not{a}}{2p_{4D} \cdot a} u_{hl}(p) \\
&= u_{h-l}(a) \frac{\not{u}_{hl}(a) \not{p}_{4D} u_{hl}(p)}{2p_{4D} \cdot a} + u_{hl}(a) \frac{\not{u}_{h-l}(a) \not{p}_{4D} u_{hl}(p)}{2p_{4D} \cdot a} + \\
&\quad u_{h-l}(p_{4D}) \frac{\not{u}_{hl}(p_{4D}) \not{p}_{4D} u_{hl}(p)}{2p_{4D} \cdot a} + u_{hl}(p_{4D}) \frac{\not{u}_{h-l}(p_{4D}) \not{p}_{4D} u_{hl}(p)}{2p_{4D} \cdot a} . \quad (5.47)
\end{aligned}$$

The remaining dependence on the vector p is in the coefficients of each spinor, which can be evaluated in terms of p_a , a and μ_p^\pm , once a choice is given for a .

For the rest of this project a is chosen to be

$$a^\mu = \frac{1}{2} \begin{pmatrix} 1 \\ -1 \\ 0 \\ 0 \\ 0 \\ 0 \end{pmatrix} , \quad (5.48)$$

as this gives one of the simplest forms for all the expressions for the spinor definition used. With this choice of reference vector the spinor can be written as

$$u_{h,l}(p) \rightarrow u_{h,l}(p_{4D}) - hl \frac{\not{u}_{h-l}(p_{4D})}{2p_{4D} \cdot a} \mu_p^{-hl} . \quad (5.49)$$

The reason for this choice is to ensure that only two spinors appear in the replacement for the 6 dimensional spinor, rather than the four that can appear in general, and that they have different helicities. If this is ensured then in an expansion of the product of a spinor for a 6 dimensional vector and a spinor for any other 4 dimensional vector, at most one term will exist and in the expansion of a spinor product for two 6 dimensional vectors, at most two terms will exist. To ensure that only two spinors appear in the replacement for the 6 dimensional spinor and that they have different helicities requires that the products $\not{u}_{hl}(p_{4D}) \not{p}_{4D} u_{hl}(p)$ and $\not{u}_{h-l}(a) \not{p}_{4D} u_{hl}(p)$ vanish. This choice of vector is the unique real vector that causes both of these products to vanish. It is also the vector that when contracted with a momenta gives the expression that will appear in the square roots within its spinors.

By applying this replacement to the definition of the conjugate spinor as given in Equation 5.33, the replacement for the conjugate spinor can be derived, which gives

$$\underline{u}_{h,l}(p) \rightarrow \underline{u}_{h,l}(p_{4D}) + hl \frac{\underline{u}_{-h,l}(p_{4D}) \not{a}}{2p_{4D} \cdot a} \mu_p^{-hl} . \quad (5.50)$$

Combining this replacement with the replacement for spinors, a replacement for a slashed matrix can also be derived which is given by

$$\not{p} = \not{p}_{4D} + \frac{\not{a}}{2p_{4D} \cdot a} \mu_p^+ \mu_p^- + \sum_{h,l=\pm} \frac{hl (\not{a} u_{-h,-l}(p_{4D}) \underline{u}_{h,l}(p_{4D}) - u_{h,l}(p_{4D}) \underline{u}_{-h,-l}(p_{4D}) \not{a}) \mu_p^{hl}}{2p_{4D} \cdot a} . \quad (5.51)$$

Using these replacements on spinor products gives the relations

$$\underline{u}_{h,l}(p) u_{-h,l}(q) = \underline{u}_{h,l}(p_{4D}) u_{-h,l}(q_{4D}) = -hl \underline{u}_l(p_{4D}) u_l(q_{4D}) \quad (5.52)$$

$$\underline{u}_{h,l}(p) u_{-h,-l}(q) = hl \frac{1}{2} \left(\frac{\underline{u}_{-h,l}(p_{4D}) \not{a} u_{-h,-l}(q_{4D})}{p_{4D} \cdot a} \mu_p^{-hl} + \frac{\underline{u}_{h,l}(p_{4D}) \not{a} u_{h,-l}(q_{4D})}{q_{4D} \cdot a} \mu_q^{-hl} \right) , \quad (5.53)$$

where the right hand sides depend only on 4 dimensional vectors and can therefore be written as 4 dimensional spinor products.

5.3 Calculating 6 Dimensional Amplitudes

The amplitudes calculated in this project will always have physical external particles in 4 dimensions. This limits where the extra dimensions can be introduced into the tree and loop amplitudes being calculated. Therefore many of the amplitudes can be greatly simplified. It is also necessary to check the validity of the properties needed for 6 dimensional generalised unitarity calculations and for the state sum reduction using scalar particles.

The corners of the cut loops require tree amplitudes. These tree amplitudes will have external momenta for all but the two legs that are cut loop propagators. Therefore these are the only momenta that can have 6 dimensional components. As there are only two particles that have 6 dimensional components, by conservation of momentum, the 5th and 6th components of their momenta must be equal and opposite to the values in the other particles' momenta. A simple logic would also imply that as there are no other 6 dimensional vectors, the direction in the 5th and 6th components should be irrelevant and a rotation between the components should not change

anything, but this needs to be proven explicitly and it must be shown precisely which types of amplitudes are guaranteed to show this behaviour. There are also restrictions on the helicity states that can appear for external and internal particles. In general, quarks and gluons both have twice as many states as they do in 4 dimensions, however the Higgs boson has one state, which is the same as in 4 dimensions.

To show that loop amplitudes do not depend on the direction of the 5th and 6th components of the momenta, the dependence of tree amplitudes on them must be shown. It must then be shown that any dependence cancels when used in generalised unitarity calculations. The simplest case is an n gluon amplitude which will therefore be handled first. Without loss of generality the first two particles are chosen to be the ones from the cut loop propagators and will be called p and q respectively. Using Feynman rules, the possible terms that can appear in any given amplitude can be calculated without evaluating the amplitude in full. For a pure gluon amplitude the Feynman rules are given in [Table 2.1](#). To prove exactly what terms can be contributed for each diagram, one vertex can be chosen and starting with its vertex factor, new vertex factors and propagators connecting them can be added. These can then be split into separate terms. If this process is repeated, at each stage any term will have the form

$$C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j^{\mu_j} \prod_{k=1}^n g^{\nu_k \sigma_k}}{\prod_{l=1}^{N_g+m-n-2} S_l^2}, \quad (5.54)$$

where C is a complex number, P , Q , R and S are momentum vectors containing some combination of neighbouring momentum from the total amplitude, m and n count how many of their respective types of factors are in the term and obey the relations

$$0 \leq m \leq \left\lfloor \frac{N_g}{2} - 1 \right\rfloor \quad m + 1 \leq n \leq \left\lfloor \frac{N_g}{2} \right\rfloor, \quad (5.55)$$

and N_g is the number of external indices or equivalently the number of gluon polarisation vectors the term needs to be contracted with to form an amplitude. The number of vertex factors of each type can be calculated as

$$\begin{aligned} N_3 &= N_g + 2m - 2n \\ N_4 &= n - m - 1, \end{aligned} \quad (5.56)$$

where N_3 is the number of three gluon vertices and N_4 is the number of four gluon vertices. This form will be proved in [Appendix A](#).

When these terms are contracted with the gluon polarisation vectors, each polarisation vector will either contract with a free momenta, R , or with a metric, g , that will contract a pair of polarisation vectors together. From this, every term in an amplitude is of the form

$$C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j \cdot \epsilon_j \prod_{k=1}^n \epsilon_k \cdot \epsilon'_k}{\prod_{l=1}^{N_g+m-n-2} S_l^2}, \quad (5.57)$$

where $\epsilon_{j/k}^{(l)}$ are gluon polarisation vectors.

Extending the general form of a term to an amplitude with a quark line, can be done by tracing along the quark line, contracting each gamma matrix from a vertex factor with a gluon term of the form in [Equation 5.54](#). Each time a gluon term is added, either a free vector, R , or a metric, g , must be contracted with the gamma matrix in the vertex term. If a vector is contracted then it will introduce a slashed matrix into the quark line. If a metric is contracted then a free gamma matrix is left that will be contracted with a gluon polarisation vector and so introduce a slashed gluon polarisation vector to the quark line. In general therefore, the terms in an amplitude with one quark pair and any number of gluons are of the form

$$C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-N_{qg}-2n} R_j \cdot \epsilon_j \prod_{k=1}^n \epsilon_k \cdot \epsilon'_k}{\prod_{l=1}^{N_g+m-n-N_{qR}-1} S_l^2} \bar{u} \left(\prod_a^{N_{qR}+N_{qg}-1} \psi_a T_a \right) \psi_{(N_{qR}+N_{qg})} u, \quad (5.58)$$

where N_{qg} of the U_a s are gluon polarisation vectors and the remaining N_{qR} of the U_a s, along with T_a s, are momentum vectors containing some combination of neighbouring momenta from the total amplitude and the summation limits now obey the conditions

$$\begin{aligned} 0 \leq m &\leq \left\lfloor \frac{N_g + N_{qR} + N_{qg}}{2} \right\rfloor - N_{qR} - N_{qg} \\ m + N_{qR} \leq n &\leq \left\lfloor \frac{N_g + N_{qg} + N_{qR}}{2} \right\rfloor - N_{qg}. \end{aligned} \quad (5.59)$$

To show how the amplitudes depend on the direction of the extra components of

vectors, it is required to show how the various types of term depend on the extra dimensions. Three of the types of term needed are products of two momentum vectors, products of two polarisation vectors and products of polarisation vectors and momentum vectors. The final type of term is spinor chains with an odd number of gamma matrices, each contracted with either a momenta or a polarisation vector. Firstly, products of two momenta will be considered. They can only depend on the 4th and 5th dimensional components if one or both of them contains one of the loop propagator momenta, p or q . Without loss of generality, both of the momenta in a product can be expanded as the sum of a possibly massive momentum vector, which contains all contributions from any other momenta and one, both or none of the loop momenta. It is then possible to expand the products into multiple terms. The terms that this can produce and their expansions in terms of their 4 dimensional equivalent vectors are

$$R \cdot S = R \cdot S \quad (5.60)$$

$$R \cdot p = \frac{\mu_p^- \mu_p^+ R \cdot a}{2p_{4D} \cdot a} + R \cdot p_{4D} = p_{4D,M} \cdot R \quad (5.61)$$

$$R \cdot q = \frac{\mu_q^- \mu_q^+ R \cdot a}{2q_{4D} \cdot a} + R \cdot q_{4D} = q_{4D,M} \cdot R \quad (5.62)$$

$$p \cdot p = q \cdot q = 0 \quad (5.63)$$

$$\begin{aligned} p \cdot q &= q_{4D} \cdot p_{4D} + \frac{(\mu_q^+ p_{4D} \cdot a - \mu_p^+ q_{4D} \cdot a)(\mu_q^- p_{4D} \cdot a - \mu_p^- q_{4D} \cdot a)}{p_{4D} \cdot a q_{4D} \cdot a} \\ &= q_{4D,M} \cdot p_{4D,M} - \frac{1}{2} (\mu_q^- \mu_p^+ + \mu_p^- \mu_q^+) , \end{aligned} \quad (5.64)$$

where R and S are the massive vectors containing only 4 dimensional momenta. All of these terms only depend on $\mu_{p/q}^\pm$ through the combinations $\mu_p^+ \mu_p^- = \mu_q^+ \mu_q^-$ and $\mu_p^+ \mu_q^- = \mu_q^+ \mu_p^- = -\mu_p^+ \mu_p^-$. These forms are the only ways that $\mu_{p/q}^\pm$ can occur in Minkowski products of momenta.

The next form that occurs is $R_j \cdot \epsilon_j$. As for Minkowski products of momenta, the momenta, R , can be expanded as a 4 dimensional massive vector and possibly one or both of the 6 dimensional vectors. The resulting terms can be expanded in terms of 4 dimensional equivalent vectors as

$$\epsilon_{hh}(r) \cdot S = \frac{\underline{u}_{+h}(r) \not{k} \underline{u}_{-h}(r)}{2\sqrt{2}r \cdot k} \quad (5.65)$$

$$\begin{aligned}
\epsilon_{hh}(r) \cdot p &= \frac{u_{+h}(r) \not{p} \not{k} u_{-h}(r)}{2\sqrt{2}r \cdot k} \\
&= \frac{u_{+h}(r) \not{p} \not{k} u_{-h}(r) \mu_p^+ \mu_p^-}{4\sqrt{2}r \cdot k p_{4D} \cdot a} + \frac{u_{+h}(r) \not{p}_{4D} \not{k} u_{-h}(r)}{2\sqrt{2}r \cdot k} \\
&= \frac{u_{+h}(r) \not{p}_{4D,M} \not{k} u_{-h}(r)}{2\sqrt{2}r \cdot k} \tag{5.66}
\end{aligned}$$

$$\begin{aligned}
\epsilon_{hh}(p) \cdot S &= \frac{u_{+h}(p) \not{S} \not{k} u_{-h}(p)}{2\sqrt{2}p \cdot k} \\
&= \frac{u_{+h}(p_{4D}) \not{p} \not{S} \not{k} u_{-h}(p_{4D}) \mu_p^+ \mu_p^-}{8\sqrt{2} \left(p_{4D} \cdot k + \frac{k \cdot a}{2p_{4D} \cdot a} \mu_p^+ \mu_p^- \right) (p_{4D} \cdot a)^2} + \frac{u_{+h}(p_{4D}) \not{S} \not{k} u_{-h}(p_{4D})}{2\sqrt{2} \left(p_{4D} \cdot k + \frac{k \cdot a}{2p_{4D} \cdot a} \mu_p^+ \mu_p^- \right)} \\
&= \frac{u_{+h}(p_{4D}) \not{p} \not{S} \not{k} u_{-h}(p_{4D}) \mu_p^+ \mu_p^-}{8\sqrt{2} p_{4D,M} \cdot k (p_M \cdot a)^2} + \frac{u_{+h}(p_{4D}) \not{S} \not{k} u_{-h}(p_{4D})}{2\sqrt{2} p_{4D,M} \cdot k} \tag{5.67}
\end{aligned}$$

$$\begin{aligned}
\epsilon_{h-h}(p) \cdot S &= \frac{u_{+h}(p) \not{S} \not{k} u_{-h}(p)}{2\sqrt{2}p \cdot k} \\
&= h \frac{S \cdot a p_{4D} \cdot k - k \cdot a p_{4D} \cdot S}{\sqrt{2} p_{4D} \cdot a \left(p_{4D} \cdot k + \frac{k \cdot a}{2p_{4D} \cdot a} \mu_p^+ \mu_p^- \right)} \mu_p^{-h} \\
&= h \frac{S \cdot a p_{4D,M} \cdot k - k \cdot a p_{4D,M} \cdot S}{\sqrt{2} p_{4D,M} \cdot a p_{4D,M} \cdot k} \mu_p^{-h}, \tag{5.68}
\end{aligned}$$

where only terms involving p are shown, as it is easy to produce the equivalent terms involving q from the versions shown. For these cases it can be seen that factors of $\mu_p^+ \mu_p^-$ appear in many places. However, lone factors of μ_p^+ or μ_p^- only appear in terms from gluons with opposite helicities and always have the sign of the μ as the second sign of the helicity. The other case, involving contracting the polarisation vector for one of the 6 dimensional momenta with the other 6 dimensional momenta, show the same pattern as for the ones shown above but are too complex to show here.

The same can be done for the contraction of two polarisation vectors and again the same relations apply. Now, as there are two polarisation vectors, it is possible for both gluons to have the same pair of helicities and for the two helicities on each gluon to be opposite, i.e. $\epsilon_{h-h} \cdot \epsilon_{h-h}$. In this situation the terms will have an overall factor of $\mu_p^{\pm 2}$. If the two gluons have exactly opposite polarisations and both have their two signs opposite, i.e. $\epsilon_{h-h} \cdot \epsilon_{-hh}$, then there will be no overall factor of $\mu_p^+ \mu_p^-$ as factors of $\mu_p^+ \mu_p^-$ can be converted to Minkowski products of vectors. Overall, this means that each term in a pure gluon amplitude can contain many factors of $\mu_p^+ \mu_p^-$, but any lone factors of μ_p^{\pm} always come from polarisation vectors with their two helicities opposite. The sign of the factors depends on the helicities of the most common type of gluon with opposite signs. If the most common type of gluon is ϵ_{h-h} then the amplitude will have overall factors of μ_p^{-h} and the number of factors will be $\#\epsilon_{h-h} - \#\epsilon_{-hh}$,

where $\#\epsilon_{hl}$ means the number of polarisation vectors of type ϵ_{hl} .

For all pure gluon amplitudes needed to calculate the rational parts, there can only be two gluons that have these helicities, so at most there will be two overall factors of μ and that will only be the case if both the gluons have the same helicities. As the signs of the gluons either side of a cut are related, the overall factors of μ^\pm for the product of the amplitudes either side of the cut must be the same as for a single amplitude without the two gluons from the cut. Therefore, the full combination of tree amplitudes used to build a cut contribution must have no overall dependence on μ^\pm and must only depend on μ^\pm via the combination $\mu^+\mu^-$.

To extend this to amplitudes with quarks, first requires proving the forms of tree amplitudes with quarks in and then again combining them to show that any overall factors of μ^\pm cancel. The only extra contribution needed for quark amplitudes is knowing the dependence of one or more quark lines on μ^\pm . Firstly the cases of only a single spinor chain will be considered. It is clear from the vanishing products and the expansion of slashed matrices in terms of spinors that any quark line will vanish if the quarks at either end have opposite values for their first helicity signs. This separates the sets of amplitudes into two spaces as is required. It still remains to show that the two spaces are both equivalent but this just requires the amplitudes to not depend on the value of the first helicity sign. To simplify the relations all slashed momenta will be expanded as above, but this time also replacing any instances of the loop momenta q with minus the sum of all the other momenta using conservation of momenta. This leaves spinor chains with an odd number of elements which are either \not{p} , slashed polarisation vectors for p or q , a possibly massive 4 dimensional slashed momenta or slashed polarisation vectors for 4 dimensional external gluons. To further simplify the cases that need to be handled, any \not{p} can be commuted to the start of the spinor line, followed by the polarisation vector for p if present, then the polarisation vector for q and finally followed by all the purely 4 dimensional slashed vectors and polarisation vectors. Each of these swaps will also introduce an extra term of the form of a spinor chain with the two swapped elements removed, multiplied by the Minkowski product of the two items removed. These extra terms will also always have an odd number of slashed matrices and be of the same form as the original spinor chain but with less elements in them. The simplest case to handle is an amplitude with the two cut particles being the quarks and no other quarks in the amplitude. Then the form of the spinor lines can always be converted to the spinor for the quark p followed by a chain

of an odd number of slashed momenta or polarisation vectors which are all purely 4 dimensional. Any spinor chain terms that have a \not{p} in them will result in terms with no \not{p} in them once the commutations are applied as any \not{p} will be commuted to the start and then cause the term to vanish as $\underline{u}(p)\not{p} = 0$. There now remain two forms to calculate for this case which depend on the helicities of the quarks and are given by

$$\underline{u}_{hl}(p) \overbrace{\not{X}}^{n \times} u_{h-l}(q) = \underline{u}_{hl}(p_{4D}) \overbrace{\not{X}}^{n \times} u_{h-l}(q_{4D}) + \frac{\underline{u}_{-hl}(p_{4D}) \overbrace{\not{X}}^{n \times} \not{a} u_{-h-l}(q_{4D})}{4p_{4D} \cdot a q_{4D} \cdot a} \mu^+ \mu^- \quad (5.69)$$

$$\underline{u}_{hl}(p) \overbrace{\not{X}}^{n \times} u_{hl}(q) = \frac{hl}{2} \left(\frac{\underline{u}_{hl}(p_{4D}) \overbrace{\not{X}}^{n \times} u_{-hl}(q_{4D})}{p_{4D} \cdot a} + \frac{\underline{u}_{-hl}(p_{4D}) \overbrace{\not{X}}^{n \times} \not{a} u_{hl}(q_{4D})}{q_{4D} \cdot a} \right) \mu^{-hl}, \quad (5.70)$$

where \not{X} represents the n slashed matrices in the chain and n is odd. From these forms it can be seen that amplitudes with a single quark line, where the two quarks are the cut loop propagators, have an overall factor of μ^{-hl} if the quarks are both of helicity hl and have no overall factors of μ^\pm otherwise. All other dependence on μ^\pm is via the combination $\mu^+ \mu^-$.

The next simplest case is an amplitude with one of the cuts being a quark and the other being a gluon. Without loss of generality the cut quark will be chosen to be p and the cut gluon will be chosen to be q . After applying all the simplifications and relations as above, two types of spinor chain remain. Firstly are spinor chains with only 4 dimensional slashed matrices and secondly are chains with the slashed polarisation vector, $\not{\epsilon}(q)$, at the start. The different cases for these terms are given by

$$\underline{u}_{hl}(p) \overbrace{\not{X}}^{n \times} u_{h-l}(r) = \underline{u}_{hl}(p_{4D}) \overbrace{\not{X}}^{n \times} u_{h-l}(r) \quad (5.71)$$

$$\underline{u}_{hl}(p) \overbrace{\not{X}}^{n \times} u_{hl}(r) = hl \frac{\underline{u}_{-hl}(p_{4D}) \overbrace{\not{X}}^{n \times} u_{hl}(r)}{2p_{4D} \cdot a} \mu^{-hl} \quad (5.72)$$

$$\underline{u}_{hl}(p) \not{\epsilon}_{m,m}(q) \overbrace{\not{X}}^{n-1 \times} u_{h-l}(r) \propto 1 \quad (5.73)$$

$$\underline{u}_{hl}(p) \not{\epsilon}_{m,-m}(q) \overbrace{\not{X}}^{n-1 \times} u_{h-l}(r) \propto \mu^{-m} \quad (5.74)$$

$$\underline{u}_{hl}(p)\not{\epsilon}_{m,m}(q)\overbrace{\not{X}}^{n-1\times}u_{hl}(r)\propto\mu^{-hl}\quad (5.75)$$

$$\underline{u}_{hl}(p)\not{\epsilon}_{h,-l,-h,l}(q)\overbrace{\not{X}}^{n-1\times}u_{h-l}(r)\propto 1\quad (5.76)$$

$$\underline{u}_{hl}(p)\not{\epsilon}_{h,l,-h,-l}(q)\overbrace{\not{X}}^{n-1\times}u_{h-l}(r)\propto(\mu^{-hl})^2,\quad (5.77)$$

where in the later terms, factors that do not depend on μ^\pm , or only depend on μ^\pm via the combination $\mu^+\mu^-$, have been removed as the expressions are too complex to show in full here.

The last type of term that can contribute to an amplitude with a single quark line is the case where the cut particles are both gluons. If the connection between the two 6 dimensional gluons does not include the quark line, then the quark line is independent of p and q and the dependence on μ^\pm must come from pure gluon factors. The remaining cases are too complex to show here but have been calculated and are found to obey the same relations as above, namely that each gluon of helicity $\epsilon_{h,-h}$ and each quark line where the quarks have helicities \underline{u}_{hl} and u_{hl} , contributes a factor of μ^{-hl} or removes a factor of μ^{hl} . These relations can also be shown to work for multiple quark lines by examining the new cases contributed. For pure gluon amplitudes it has already been shown above that these factors of μ^\pm cancel between amplitudes when combined for a generalised unitarity contribution. Using the same logic, combined with the formulas derived above for the dependence of quark lines, it is clear that the overall factors of μ^\pm for any product of tree amplitudes is again the same as a single amplitude without the cut quarks. Therefore, for amplitudes with quark lines in, there again must be no overall dependence on μ^\pm for the combination of tree amplitudes needed for a generalised unitarity term. In addition, for any generalised unitarity contribution for an amplitude which includes a combination of quarks, gluons and optionally a Higgs boson, the only dependence on μ^\pm is via the combination $\mu^+\mu^-$.

Therefore, it has been shown that the exact value of the components p_4 and p_5 of the loop momenta are irrelevant and can be chosen freely without affecting the results of the generalised unitarity calculations and that μ^+ and μ^- can be chosen to have the same value and to always have a positive real part or be pure imaginary with a positive imaginary part. This reduces the two extra complex components of the momenta to one extra complex value for $\mu = \mu^+ = \mu^-$.

To show that the quark state reduction works, requires showing that amplitudes

do not depend on which of the two quark spaces they are from, other than an overall sign shared by all terms. The effect from the quark helicities is only able to affect the quark lines in each term. For the cases with no cut quarks, the two spaces give exactly the same value as they must have 4 dimensional type helicities. If the quarks are cut then there are more options available and the tree amplitudes, in general, will depend on which quark space is used, due to the factors of hl and μ^{-hl} that appear, but these always cancel to terms of the form $\mu^+\mu^-$, while also cancelling the dependence on which space they are from. Therefore, both spaces will give equal contributions and the state sum reduction works.

It now remains to show that the extra amplitudes needed for the gluon state reduction also simplify and obey the relations expected. For scalar subtraction amplitudes for pure gluon amplitudes, the terms that can contribute are related to those for the pure gluon amplitude, by the requirement that for any term of the gluon amplitude where the two cut gluons would have been contracted with each other using a metric tensor, there is a term of the same form but with that metric tensor removed. Each corner in these amplitudes must depend on μ^+ and μ^- only via the combination $\mu^+\mu^-$, as all contributions that give overall factors of μ^\pm , were shown earlier to be contractions of the polarisation vectors for the cut gluons with either each other, other polarisation vectors or momentum vectors and these factors can not appear in these terms. Therefore, these subtraction terms are again independent of the direction of the extra components and could even be calculated with the loop momenta taken in a direction that includes the 6th dimension, which would not otherwise have been permitted.

For amplitudes with external quarks, as well as some gluons in the loop, if any corners do not contain the quarks, then by the same logic as above, those corners must only depend on μ^+ and μ^- via the combination $\mu^+\mu^-$. If all the cut particles in a contribution are gluons and each quark pair is isolated to a single corner, then the only case to consider is amplitudes like those shown on the left of [Figure 5.2](#). Many diagrams in these amplitudes will not have the scalars enter the quark line and will therefore have the two lines connected by a gluon. For this case, the same logic as for pure gluon amplitudes can be used to show that again the term must have no overall factors of μ^+ or μ^- and can only depend on μ^\pm via the combination $\mu^+\mu^-$. The remaining type of diagram for this type of corner is where the scalar lines both enter the quark line and is shown in the rightmost diagram in [Figure 5.2](#). For this

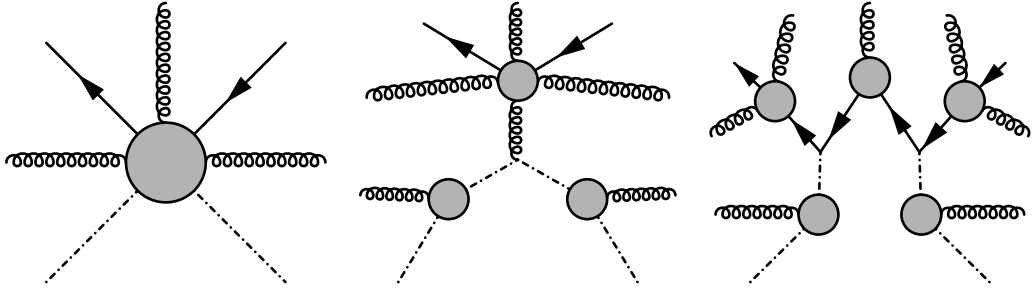


Figure 5.2: The general form of the scalar subtraction amplitude for a corner with a quark pair that does not enter the loop, along with the two types of term that contribute to it. The dash-dotted lines represent the scalar corresponding to gluons in the 6th dimension. An external gluon shown in the diagrams can represent any number of external gluons, including none. The cut particles are always the bottom two particles in each diagram and shaded circles represent all possible colour ordered diagrams that contribute to the amplitude or type of term.

case the term that is new and could potentially break the assumptions, is the quark line with the scalar lines entering it, as these introduce explicit factors of \not{n}_6 , which is not a massless physical momentum vector. Again these terms will be handled by commuting these factors to the very end of the quark chain, which will leave one term with both of these slashed matrices at the end and many terms with one or both of these vectors removed from the chain and contracted with another contribution from the line. The term which still has both factors of \not{n}_6 present and has them adjacent to each other can then be simplified by removing these two slashed matrices, as

$$\not{n}_6 \not{n}_6 = n_6^2 I = -I , \quad (5.78)$$

where I is the identity matrix for the space of gamma matrices. This term therefore gives a contribution without overall factors of μ^+ or μ^- and only depends on μ^\pm via the contribution $\mu^+ \mu^-$. The other terms produced by this commutation will contain chains with one or zero factors of \not{n}_6 , multiplied by one or two factors, each of the form of n_6 contracted with either a momentum vector or an external polarisation vector. These extra factors will all vanish as the loop momenta is restricted to only be in 5 dimensions and the polarisation vectors are for external gluons with 4 dimensional polarisations and momenta and either of these will vanish when contracted with n_6 .

The remaining type of corner that needs investigation is those that have a quark for one of the cut particles and a gluon for the other as shown in [Figure 5.3](#). Again each

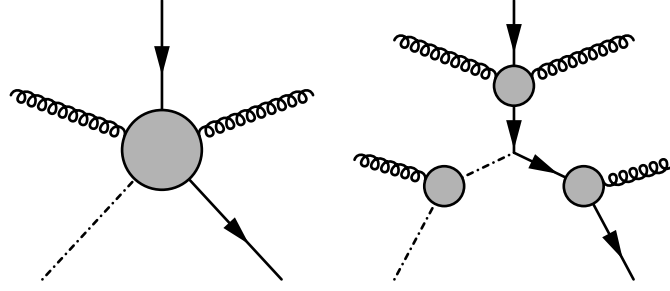


Figure 5.3: The general form of the scalar subtraction amplitude for a corner with a single quark that enters the loop, along with the form of the only type of term that contributes to it. The dash-dotted lines represent the scalar corresponding to gluons in the 6th dimension. An external gluon shown in the diagrams can represent any number of external gluons, including none. The cut particles are always the bottom two particles in each diagram and shaded circles represent all possible colour ordered diagrams that contribute to the amplitude or type of term.

term will be commuted so that a fixed order of the slashed matrices is achieved. Here, as for the other quark corner above, any extra terms produced by the commutation of \not{n}_6 must vanish as they will introduce factors of n_6 contracted with either a loop momenta or an external polarisation vector. The quark line will contain exactly one factor of \not{n}_6 and a combination of slashed external polarisation vectors and slashed momenta, which will contain a combination of the loop momenta, l , and external momenta. As for the normal amplitudes, if there are multiple factors of the loop momenta, they can be commuted to be next to each other and will then vanish. Therefore, there can be at most one loop momenta left in the amplitude after this. The remaining factor of the loop momenta will also vanish if it is commuted to the end of the quark line that is the cut quark, leaving just terms with a single factor of \not{n}_6 next to the spinor for the cut quark. This quark line is potentially problematic as it may introduce factors of μ^\pm that do not combine with a factor of the opposite sign. Using the properties of quark lines derived above, the dependence of these terms on μ can be shown to be

$$\underline{u}_{hl}(p)\not{n}_6 \overbrace{\not{X}}^{n-1\times} u_{hl}(r) = l\underline{u}_{hl}(p_{4D}) (u_{-hl}(n_{6,4D})\underline{u}_{-hl}(a) - u_{-hl}(a)\underline{u}_{-hl}(n_{6,4D})) \overbrace{\not{X}}^{n-1\times} u_{hl}(r) \quad (5.79)$$

$$\underline{u}_{hl}(p)\not{n}_6 \overbrace{\not{X}}^{n-1\times} u_{h-l}(r) = -h\underline{u}_{hl}(p_{4D})\not{a}u_{-h,-l}(n_{6,4D})\underline{u}_{-h-l}(a) \overbrace{\not{X}}^{n-1\times} u_{h-l}(r) \frac{\mu_p^{-hl}}{2p_{4D} \cdot a}, \quad (5.80)$$

where $n_{6,a}$ is a 4 dimensional vector introduced as the 4 dimensional equivalent to n_6 . $n_{6,a}$ is derived by splitting $2n_6$ into the difference between two massless 6 dimensional vectors, n_6^+ and n_6^- , which are given by

$$n_6^+ = \begin{pmatrix} 1 \\ 0 \\ 0 \\ 0 \\ 0 \\ 1 \end{pmatrix} \quad n_6^- = \begin{pmatrix} 1 \\ 0 \\ 0 \\ 0 \\ 0 \\ -1 \end{pmatrix}, \quad (5.81)$$

and then applying the normal reduction to 4 dimensional vectors. Both of these vectors, when reduced, result in the same 4 dimensional massless vector, $n_{6,a}$ which is given by

$$n_{6,4D} = \frac{1}{2} \begin{pmatrix} 1 \\ 1 \\ 0 \\ 0 \end{pmatrix}. \quad (5.82)$$

As all terms in these amplitudes must have a factor of this form, the overall dependence on μ^\pm is consistent, but it is no longer as it was for normal amplitudes. This is potentially problematic as now factors of μ^\pm will no longer cancel and there could be overall factors of μ^\pm that depend on the helicity of the cut terms. If the two corners of this form are neighbours, then there are two cases for the internal dependencies, which are given in [Table 5.2](#). If there are other corners between these two corners, then there are four cases, as all that matters is the helicities of the external and cut quarks next to the corners where the quarks exit the diagram because, as found above, multiple neighbouring corners with both cut particles being quarks have the same overall factors of μ^\pm as one corner without the internal cuts. These four cases are also shown in [Table 5.2](#). These terms, as shown in the table, are also well behaved and depend on μ^\pm only via the combination $\mu^+\mu^-$ and as such all of the calculations are independent of the direction of the extra components, other than the requirement that $l \cdot n_6 = 0$ when calculating these extra cut terms.

As all the amplitudes are independent of the value of the components of the loop momenta in the extra dimensions, it would greatly simplify implementing calculations

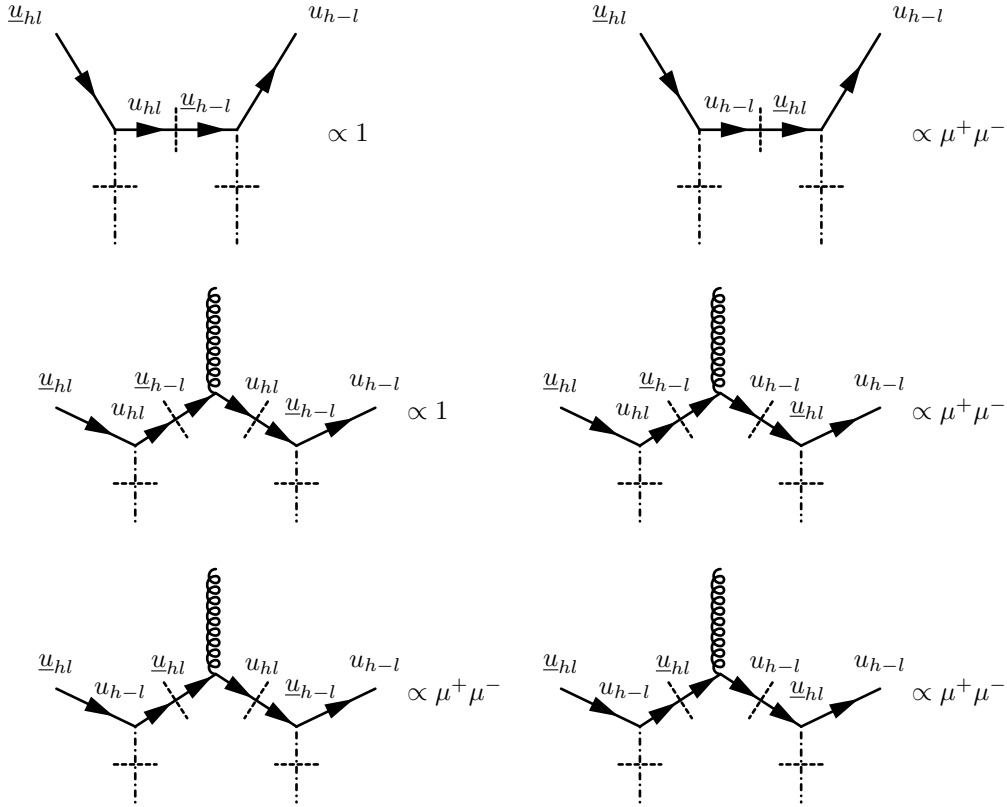


Table 5.2: The different types of quark line dependence in generalised unitarity terms for the scalar gluon subtraction terms. The top row are the contributions if the corners where the quark line exits the diagram are next to each other and the other rows are for if there are any number of extra corners between them. Only one extra corner is shown, as the contribution from a series of neighbouring corners with both cut particles being quarks, combines to give the same contribution as only a single corner with the same edge cut quarks. Any extra external gluons in any corner are not shown as the contribution depends only on the quark line and scalar gluons.

if all amplitudes were calculated in terms of the massless 4 dimensional projected momenta and the value of μ^2 . This will require new effective particle types that label the particles that are 6 dimensional. For these particles, their true momenta must be calculated by combining the massless projection with the given value for μ^2 . To actually calculate these amplitudes will still require using the full 6 dimensional Feynman rules, but all external momenta will be purely 4 dimensional and all factors and terms in the amplitude will be written in terms of purely 4 dimensional spinor products, Minkowski products, constants and the value of μ^2 . It is important to note that conservation of momenta in amplitudes involving external particles that are 6 dimensional is more complex, as it is the total 6 dimensional momenta that is conserved not the reduced momenta. The exact relation between the projected

on-shell momenta of two cut on-shell 6 dimensional particles is

$$l_{i+1,4D}^\mu = l_{i,4D}^\mu - P^\mu - a^\mu \frac{P \cdot a}{2l_{i,4D} \cdot a (l_{i,4D} \cdot a - P \cdot a)} \mu^+ \mu^- , \quad (5.83)$$

where the full 6 dimensional momenta satisfy the relation $l_{i+1}^\mu = l_i^\mu - P^\mu$ and $l_{i+1,4D}$ will be massless if and only if the full 6 dimensional vector would have been massless.

After a complete set of the lowest multiplicity and simplest amplitudes is calculated explicitly from the Feynman rules, the BCFW recursion relation can be used to calculate the higher multiplicity amplitudes, by applying it to the purely 4 dimensional momenta. However, care must be taken to use the modified conservation of momenta rule given above and to sum over all possible states including the extra 6 dimensional ones, if the two 6 dimensional momenta are on opposite sides of a cut. The exact formula that combines the two amplitudes on either side of the cut to give the contribution will also need to be calculated. It is also important to check which choices of shifted particles will be valid. These conditions are not yet known and are likely to be complex if one of the 6 dimensional particles is chosen.

5.4 Calculating the Rational Terms using 6 Dimensional Spinors

Once tree amplitudes can be calculated, the next step is to combine them to produce the coefficients of the scalar basis integrals and then combine the coefficients with the values of the basis integrals to give the rational terms. For numeric calculations an explicit numerical expression for the loop momenta is again required. As before, it would be convenient if the loop momenta could be written in terms of spinors and vectors for the external momenta. If the cut conditions are evaluated in terms of the massive 4 dimensional loop momenta, then the conditions extend in a relatively simple way and give the loop momentum for three or more cuts of

$$l_{2,4D,M}^2 = c_1 n_1^\mu + c_2 n_2^\mu + \tilde{K}_1^\mu \frac{K_2^2 (K_1^2 + \gamma)}{4\Delta} - \tilde{K}_2^\mu \frac{K_1^2 (K_2^2 + \gamma)}{4\Delta} , \quad (5.84)$$

where c_1 and c_2 now satisfy the conditions

$$c_1 c_2 = - \frac{\gamma \left(\mu^2 + \frac{K_1^2 K_2^2 (K_1^2 + K_2^2 + 2K_1 \cdot K_2)}{4\Delta} \right)}{4\Delta} , \quad (5.85)$$

and if a fourth cut is performed they are given by

$$c_1 = \frac{X \left(1 \pm \sqrt{1 + \frac{\Delta_4}{X^2 \Delta} \left(\frac{K_1^2 K_2^2 (K_1^2 + K_2^2 + 2K_1 \cdot K_2)}{\Delta} + 4\mu^2 \right)} \right)}{4K_3 \cdot n_1} \quad (5.86)$$

$$c_2 = \frac{X \left(1 \mp \sqrt{1 + \frac{\Delta_4}{X^2 \Delta} \left(\frac{K_1^2 K_2^2 (K_1^2 + K_2^2 + 2K_1 \cdot K_2)}{\Delta} + 4\mu^2 \right)} \right)}{4K_3 \cdot n_2}, \quad (5.87)$$

where the conventions for the directions of momenta and definitions for \tilde{K}_1 , \tilde{K}_2 , n_1 , n_2 , γ , Δ , Δ_4 and X are the same as for Equation 4.28. As this definition is massive it does not have a spinor representation, but its massless projection does. To project down into a spinor form, the vector a must be expanded in terms of the basis of \tilde{K}_1 , \tilde{K}_2 , n_1 and n_2 , so that extra linearly dependant vectors and spinors are not introduced. In terms of this basis, a is given by

$$a^\mu = \frac{\tilde{K}_1^\mu (K_2^2 K_1 \cdot a - \gamma K_2 \cdot a) + \tilde{K}_2^\mu (K_1^2 K_2 \cdot a - \gamma K_1 \cdot a) + n_1^\mu \gamma n_2 \cdot a + n_2^\mu \gamma n_1 \cdot a}{2K_1^2 K_2^2 - 2(K_1 \cdot K_2)^2}, \quad (5.88)$$

where $n_1 \cdot a$ and $n_2 \cdot a$ are related by

$$n_1 \cdot a \ n_2 \cdot a = \frac{(\gamma - 2K_1 \cdot K_2) \left(K_2^2 (K_1 \cdot a)^2 + K_1^2 (K_2 \cdot a)^2 - 2K_1 \cdot a K_2 \cdot a K_1 \cdot K_2 \right)}{K_1^2 K_2^2}. \quad (5.89)$$

Using this representation for a , the spinor form for the massless 4 dimensional loop momentum is given by

$$\begin{aligned} l_{2,4D}^\mu &= \left(\langle \tilde{K}_1 | + \frac{\beta + \mu^2 \frac{K_1^2 K_2 \cdot a - \gamma K_1 \cdot a}{C(c_1, c_2)}}{c_1 + \mu^2 \frac{\gamma n_2 \cdot a}{C(c_1, c_2)}} \langle \tilde{K}_2 | \right) \gamma^\mu \\ &= \left(\left(\alpha + \mu^2 \frac{K_2^2 K_1 \cdot a - \gamma K_2 \cdot a}{C(c_1, c_2)} \right) | \tilde{K}_1 \right] + \left(c_1 + \mu^2 \frac{\gamma n_2 \cdot a}{C(c_1, c_2)} \right) | \tilde{K}_2 \right] \\ &= \left(\left(\alpha + \mu^2 \frac{K_2^2 K_1 \cdot a - \gamma K_2 \cdot a}{C(c_1, c_2)} \right) \langle \tilde{K}_1 | + \left(c_2 + \mu^2 \frac{\gamma n_1 \cdot a}{C(c_1, c_2)} \right) \langle \tilde{K}_2 | \right) \\ &\quad \gamma^\mu \left(| \tilde{K}_1 \right] + \frac{\beta + \mu^2 \frac{K_1^2 K_2 \cdot a - \gamma K_1 \cdot a}{C(c_1, c_2)}}{c_2 + \mu^2 \frac{\gamma n_1 \cdot a}{C(c_1, c_2)}} | \tilde{K}_2 \right] \end{aligned}, \quad (5.90)$$

where the common function $C(c_1, c_2)$ is extracted to simplify the representation and is given by

$$\begin{aligned}
C(c_1, c_2) &= 2(K_2^2 K_1 \cdot a(K_1^2 + K_1 \cdot K_2) - K_1^2 K_2 \cdot a(K_2^2 + K_1 \cdot K_2) + 2\Delta(c_1 n_1 \cdot a + c_2 n_2 \cdot a)) \\
&= 4c_1 \Delta n_1 \cdot a + 2(K_2^2 K_1 \cdot a(K_1^2 + K_1 \cdot K_2) - K_1^2 K_2 \cdot a(K_2^2 + K_1 \cdot K_2)) \\
&\quad + \frac{(K_2^2 K_1 \cdot a^2 + K_1^2 K_2 \cdot a^2 - 2K_1 \cdot a K_2 \cdot a K_1 \cdot K_2) \left(\mu^2 + \frac{K_1^2 K_2^2 (K_1^2 + K_2^2 + 2K_1 \cdot K_2)}{4\Delta} \right)}{c_1 n_1 \cdot a} \\
&= 4c_2 \Delta n_2 \cdot a + 2(K_2^2 K_1 \cdot a(K_1^2 + K_1 \cdot K_2) - K_1^2 K_2 \cdot a(K_2^2 + K_1 \cdot K_2)) \\
&\quad + \frac{(K_2^2 K_1 \cdot a^2 + K_1^2 K_2 \cdot a^2 - 2K_1 \cdot a K_2 \cdot a K_1 \cdot K_2) \left(\mu^2 + \frac{K_1^2 K_2^2 (K_1^2 + K_2^2 + 2K_1 \cdot K_2)}{4\Delta} \right)}{c_2 n_2 \cdot a}.
\end{aligned} \tag{5.91}$$

As in 4 dimensions the bubble can not directly use the above form as it only has one independent external momenta so an extra arbitrary vector is required to parametrise the set of momenta. By the same method as above, the loop momenta cut conditions can be solved and the loop momentum given by

$$\begin{aligned}
l_{1,4D,M} &= (1-y)\tilde{K}_1 + \frac{yK_1^2}{2K_1 \cdot \chi} \chi - tn_1 - \frac{y(1-y)K_1^2 - \mu^2}{2tK_1 \cdot \chi} n_2 \\
l_{2,4D,M} &= -y\tilde{K}_1 - \frac{(1-y)K_1^2}{2K_1 \cdot \chi} \chi - tn_1 - \frac{y(1-y)K_1^2 - \mu^2}{2tK_1 \cdot \chi} n_2,
\end{aligned} \tag{5.92}$$

where again this solution gives a massive vector. As for the loop momentum parametrisation for pentagons, boxes and triangles, the reduced 4 dimensional loop momentum will have a spinor representation which will factorise. As in 4 dimensions no extra variants of the loop momentum expression are needed to handle the case where $K_1^2 \rightarrow 0$ as this will not contribute due to the presence of K_1^2 in the scalar integral's value.

It is necessary to know exactly what form the combinations of tree amplitudes will have in terms of the remaining degrees of freedom and which of these coefficients are needed to form the rational terms for each cut in order to evaluate the coefficients. This is calculated by taking the form of the loop momentum solution, which for three or more cuts is given above, and substituting it into the form of the integral numerators as given in Equations 5.9 to 5.12. The directions for the vectors n_i can be chosen, subject to the conditions that defined them, for each box, triangle and bubble diagram, so that the resulting forms are as simple as possible, without changing the form that results because of the symmetry of the numerator forms. The result will be

a power series in the remaining degrees of freedom, with a relatively complex shape in the space of the different degrees of freedom.

The reduced momenta forms given above include extra poles, but these must cancel when the different amplitudes and coefficients are combined to produce the full generalised unitarity contribution. Care must be taken if the vectors become 4 dimensional, as then $\mu^2 \rightarrow 0$ and the loop momenta should collapse back to the standard 4 dimensional form, but may numerically diverge from this solution. It is also important to ensure that the product of the loop momenta with a does not vanish, as then the projection would fail. This should be unlikely to happen as long as the products of external momenta with a do not vanish and complex values for the extra components are used. If the product does vanish then there is likely to be a factorisation of the badly behaved components that allows them to cancel between the coefficients. No extra cases will be needed to handle the situation where the momenta of one or both of the external corners used to construct the momenta representation is massless, unlike in 4 dimensions, as long as μ^2 is not zero.

The final component needed to calculate the rational terms, are the subtraction terms resulting from contributions that have extra cuts. These again will be functions of the coefficients for the higher terms, though there are now significantly more terms to handle. It is hoped that some of the coefficients can be ignored if it can be shown that both their contribution to the terms that are actually needed and their contribution to the subtractions for lower diagrams will always vanish when the relevant terms are projected out. As for the pure 4 dimensional calculation, each coefficient can be projected out using complex Fourier projections. Combining all these steps, the coefficients for each amplitude and any subtraction amplitude can be calculated and then combined using [Equation 5.17](#) and [Equation 5.6](#), to give the rational terms in the four dimensional helicity scheme. The full details of how to implement this calculation have not been worked out in this project due to time constraints.

Chapter 6

Implementing the Calculations

Throughout this project Mathematica 8[19] has been used to implement the calculations and derive the formulas given in this thesis. For the 4 dimensional calculations in Chapters 3 and 4, the Mathematica package Spinors@Mathematica (S@M)[20] was used, which provides an implementation of the spinor helicity formalism in 4 dimensions for both analytical and numeric calculations. This Mathematica package uses the same definition for spinors as is used in BlackHat. This enables direct numerical comparisons of the values produced by BlackHat and those produced by the Mathematica implementation, as no overall external momentum dependent phases need to be combined with the values before comparison. The automatic numerical calculations for tree amplitudes and the cut coefficients are implemented in terms of the basic spinor and 4-vector objects provided by this package, as a collection of Mathematica source files. This implementation can calculate the tree amplitudes and coefficients of the one loop scalar basis integrals for any amplitude containing one or zero Higgs bosons with any number of quark pairs and gluons. The Mathematica code for the 4 dimensional calculations is divided into three files: common shared code in `Common.txt`; the tree level, colour ordered, helicity amplitudes in `HelAmplN.txt` and the one loop cut part calculations in `Loop-Cuts.txt`. Types of particles are labelled by objects of the form `Type[Properties, ...]`, where `Type` is the type of particle, either `Higgs`, `Phi`, `Gluon` or `Quark`, and `Properties, ...` includes the helicity of the particle and, for quarks, its direction (1 for a quark and -1 for an anti-quark which is the

same as a quark travelling in the opposite direction) and flavour. For example, a positive helicity gluon is labelled as `Gluon[1]`, a negative helicity quark of flavour 2 is labelled as `Quark[1,-1,2]` and a negative helicity anti quark of flavour 3 is labelled as `Quark[-1,-1,3]`. Various properties of the types of particles that are used later in the calculations are defined in `Common.txt`. This file also loads `S@M`, defines a few extra methods for it and improves the functions that calculate spinors from momenta to handle cases where elements are infinite or indeterminate (of the form $0/0$).

The tree level amplitudes are calculated using the function `HelAmplN`, which is called using expressions of the form `HelAmplN[{Momenta,...},{ParticleTypes,...}]`, where the `Momenta` are either lists containing the elements of the momenta, or names that have been defined in `S@M` as four vectors or spinors and have their momenta defined. This function's arguments are separated based on whether or not a particle is colour ordered and then the particles are sorted into a canonical order so that when declaring amplitudes, the minimum number of combinations needs to be checked for. In `HelAmplN.txt`, explicit formulas for the MHV and anti-MHV amplitudes are declared, along with some amplitudes that vanish. These amplitudes are used by the BCFW implementation to build up higher multiplicity amplitudes. It was discovered that some theoretically valid shifts had terms of the form $0/0$ which caused numerical problems and were not handled well in this implementation, therefore multiple pairs of shift particles are tried until one is found that provides a valid amplitude. Many cases where this could happen were handled by explicitly checking for certain forms of terms and declaring that they would vanish, even though often they are badly behaved numerically. Care was needed, while combining the two amplitudes to form a BCFW term, to ensure that the correct factors were introduced by the contribution for the particles either side of the cut, so that they form a valid propagator, given that they were calculated with momenta in the opposite directions which are not guaranteed to have a simple relationship. A simpler method was used in the cut part calculations, where the spinors for the reverse particle are explicitly defined as i times the non-reversed particle's spinors, as this does not require as much care on these factors.

The coefficients of the scalar loop integrals were also calculated in Mathematica using the code in `Loop-Cuts.txt`. For this section of calculations the processes and set of momenta must be declared using the functions `DeclareProcess` and `DeclareMomConf`. Both of these functions take as their first argument a list containing

a pair of lists, where the first list is the colour ordered particles and the second is the non-colour ordered particles. `DeclareProcess` also has a second argument which is the quarks that turn left after entering the loop. For each quark pair either the quark or the anti-quark is left turning. It is also possible to have a closed quark loop which is represented by the quark flavour -1 . It is not important whether the closed loop quark is a quark or an anti-quark, as the differences always cancel in closed quark loops. The helicity property of left turning quarks is also not specified as it is implicit from the helicity of the external quark. Both these declaration functions return a new unique symbol that labels the process or momentum configuration.

The Mathematica code generates the set of all possible non-vanishing cuts and the sets of cut propagators for each cut. Cuts are described by a pair of lists, the first of which lists particles which are the first particle not in the previous corner, which is the same as the first particle in that corner unless the corner is empty. The second list shows which corner contains each non-colour ordered particle. For example, $\{\{2,4,4,5\},\{2\}\}$ in an amplitude with five ordered particles and one unordered particle, means the first corner has particles 2 and 3, the second corner contains the first and only unordered particle, the third corner contains just particle 4 and the last corner contains particles 5 and 1. This split is shown in [Figure 6.1](#). The list of propagator particles is given by the particles heading out of each corner, towards the next corner, in corner number order. This labelling is also shown in [Figure 6.1](#). The list of possible cuts is generated using `SplitOptions[process,n]`, where `process` is the symbol returned by `DeclareProcess` and `n` is the number of cuts wanted and is 2 for bubbles, 3 for triangles and 4 for boxes. The list of possible sets of propagators for a given cut is calculated using `SplitValidPropOptions[process,split]`, where `split` is a split as returned by `SplitOptions`. It is possible for a split to be returned by `SplitOptions`, but for it to have no valid sets of propagators and therefore for `SplitValidPropOptions` to return an empty list.

The numeric value of the coefficient of each scalar function is calculated using the functions `CalculateBoxContribution`, `CalculateTriContribution` and `CalculateBubbleContribution`, all of which take the process, momentum configuration, cut and cut propagators as arguments, in that order.

For tree level amplitudes the values were tested against the values produced by `BlackHat` and found to agree to within the expected error, which is of the order of the square root of the error in the input, due to square roots in the spinor definitions. This

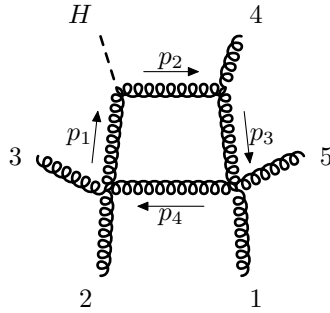


Figure 6.1: The cut labelled $\{\{2, 4, 4, 5\}, \{2\}\}$ for a five gluon, one Higgs amplitude. The propagators are labelled by p_i , where i is the index in the list of propagators for the cut. The direction that the propagator particle is viewed as travelling in is shown with arrows. All external particles are taken as outbound. This shape of diagram and cut label is equally valid for any set of five ordered and one non-ordered external particles.

is performed with the python script [TreeTests.py](#) which links to BlackHat, generates momenta and processes and calculates them, then generates Mathematica code to calculate each amplitude and compare it to the value calculated with BlackHat. The cut part integral coefficients were also tested to ensure that both implementations produced the same set of non vanishing amplitudes and the same numeric values for all cuts for all types of amplitude and again the values were found to agree. This is tested using the python script [OneLoopTests.py](#). The only issue was with amplitudes with a closed quark loop which were out by a factor of -1 , which is a factor that, in BlackHat, has been absorbed into the coefficients rather than the integral. There were also a few cuts that one version produced but the other did not produce, but these always had a value of zero as is required. This inconsistency is not necessarily an issue as time spent calculating coefficients that can be shown to be zero is traded against the effort needed to work out which terms are going to vanish. The main class of amplitudes that vanish but are still produced by BlackHat, was cuts for Higgs boson amplitudes where the corner containing the Higgs boson had helicities that vanish. It would improve the efficiency of BlackHat's calculation to not calculate these contributions that are known to always vanish. The methods of implementation for the cut part calculations for amplitudes with a Higgs boson and for pure quark and gluon amplitudes were compared in BlackHat and showed the same differences as were found in this project. This is another confirmation that the existing implementation in BlackHat is correct for calculating the cut parts of one loop amplitudes with a Higgs boson.

Care was required when comparing cuts between BlackHat and the Mathematica

implementation as different labelling conventions are used. If the amplitude contains only colour ordered particles, then there is only a difference if the first particle in the first corner is not particle 1, in which case the last element from the label is moved to the front to form the BlackHat label. If there are non-colour ordered particles then it is even more complex to convert the split labels, as BlackHat inserts the non-colour ordered particles into the colour ordered particles in different positions for each cut, so it can label the cut as if all the particles were ordered. This conversion along with the many other conversions needed to convert between BlackHat labels and Mathematica labels are in [BHMathLink.py](#).

There is also Mathematica code in [DrawCuts.txt](#) to draw graphs showing the possible cuts for an amplitude along with which diagrams contribute poles to which other diagrams. This was used to produce Figures 4.3–4.5. These diagrams are produced using pdfLaTeX[21], feynmp[22], pdfcrop[23] and dot[24] to draw the individual Feynman diagrams, process them and then combine them to produce the overall figure. Similar functionality is also available in BlackHat and can be accessed using the `--plotgraph` option of [OneLoopTests.py](#) or its short version.

For the rational terms, a Mathematica implementation of spinors in 6 dimensions has been developed. It has been used to derive and check the relations between 6 dimensional and 4 dimensional spinors and between 6 dimensional momenta and spinor expressions. It is implemented in [6DSpinorHelicity.txt](#) and uses the spinors as defined in [Section 5.1](#). Both algebraic and numeric calculations can be performed. There are many rearrangements and simplifications available. Spinors and conjugate spinors are represented by the objects `spinor[name,helicity,...]` and `spinorbar[name,helicity,...]`, where `name` is the label for the particle and `helicity,...` represents the one or two helicity signs for the spinor depending on if the spinor is 4 or 6 dimensional respectively. Slashed matrices are represented by `Sm[name,helicity]` for massless momenta and `SmM[name,helicity]` for massive momenta, where `helicity` is the helicity of the spinor space that the slashed matrix is in. These are combined in products using the function `Sp[elements,...]` where each element is a spinor, conjugate spinor or slashed matrix. It is also possible to insert linear functions of these elements as elements in larger spinor products. All products are expanded as fully as possible and completely closed spinor products are extracted and separated into individual factors. Closed spinor products will also reverse direction to produce a canonical form. As was shown in [Section 5.1](#) many spinor products vanish

```

Sp[spinorbar[p, 1, 1], Sm[r], Sm[s], Sm[t], spinor[q, -1, -1]]
Sp[spinorbar[p, 1, 1], Sm[r], Sm[s], Sm[t], spinor[q, 1, -1]]
Sp[spinorbar[q, 1, -1], Sm[t], Sm[s], Sm[r], spinor[p, 1, 1]]
% == %%
%% // TraditionalForm
Sp[spinorbar[q, 1, -1] + spinorbar[p, 1, -1],
  (Sp[Sm[r], Sm[s]] + Sp[Sm[s], Sm[r]]), spinor[p, -1, -1]]
0
Sp[u_{1,1}[p], Sm[r, 1], Sm[s, -1], Sm[t, 1], u_{1,-1}[q]]
Sp[u_{1,1}[p], Sm[r, 1], Sm[s, -1], Sm[t, 1], u_{1,-1}[q]]
True
(u_{1,1}(p), Sm(r, 1), Sm(s, -1), Sm(t, 1), u_{1,-1}(q))
Sp[u_{1,-1}[p], Sm[r, 1], Sm[s, -1], u_{1,-1}[p]] + Sp[u_{1,-1}[p], Sm[s, 1], Sm[r, -1], u_{1,-1}[p]] +
  Sp[u_{1,-1}[q], Sm[r, 1], Sm[s, -1], u_{1,-1}[p]] + Sp[u_{1,-1}[q], Sm[s, 1], Sm[r, -1], u_{1,-1}[p]]
2 Mp[Mom[s], Mom[r]] Sp[u_{1,-1}[q], u_{1,-1}[p]]
DeclareVectorDimension[r, 4]
DeclareVectorDimension[s, 4]
Sp[spinorbar[p, 1, 1], spinor[q, -1, 1]]
Sp[spinorbar[p, 1, 1], spinor[q, -1, -1]]
Sp[spinorbar[r, 1, 1], spinor[s, -1, 1]]
Sp[spinorbar[r, 1, 1], spinor[s, -1, -1]]
4
4
Sp[u_{1,1}[p], u_{1,1}[q]]
Sp[u_{1,1}[p], u_{1,-1}[q]]
Sp[u_{1,1}[r], u_{1,1}[s]]
0

```

Figure 6.2: Examples of the spinor and slashed matrix objects and their products using the implementation in `6DSpinorHelicity.txt`. The first block of equations show various of the simplifications that are performed automatically by the implementation for any momenta while the second block shows simplifications that only happen if the momenta are 4 dimensional.

and these conditions are implemented here and cause the relevant terms to simplify to zero. Using the function `DeclareVectorDimension[momenta, dimensions]`, it is possible to declare that vectors are in a lower number of dimensions, for example, a 4 dimensional vector while working in 6 dimensions, where `momenta` is the name of the momenta and `dimensions` is the number of dimensions that the vector is in. Declaring that vectors are 4 dimensional while working with 6 dimensional spinors and momenta allows the extra properties given in [Section 5.1](#) to be used and many more products to be simplified away. These spinor and slashed matrix objects and some of the automatic simplifications performed are shown in [Figure 6.2](#).

Momenta are represented by `Mom[name]` for massless momenta and `MomM[name]` for massive momenta, where `name` is the name of the vector. These are combined to give their Minkowski product using the function `Mp[mom1, mom2]`, where `mom1` and `mom2` are both momenta and examples are shown in [Figure 6.3](#). Again, as for spinor products,

```

Mp[Mom[p], Mom[q]]
% // TraditionalForm
Mp[Mom[q], Mom[p]]
% == %%
Mp[Mom[p], Mom[p]]
Mp[MomM[p], MomM[p]]
% // TraditionalForm
Mp[Mom[q], Mom[p]]

Pq · Pp
Mp[Mom[q], Mom[p]]

True
0
Mp[MomM[p], MomM[p]]

Pp · Pp

```

Figure 6.3: Examples of momenta objects and their Minkowski products using the implementation in `6DSpinorHelicity.txt`.

```

Mp[Sp[spinorbar[p, 1, 1], γ, Sm[q], spinor[r, -1, -1]], Mom[s]]
% // TraditionalForm
% // DoMToSm // TraditionalForm
Mp[Sp[u_{1,1}[p], σ[1], Sm[q, -1], u_{-1,-1}[r]], Mom[s]]
⟨u_{1,1}(p), σ^μ, Sm(q, -1), u_{-1,-1}(r)⟩ · p_{s μ}
⟨u_{1,1}(p), Sm(s, 1), Sm(q, -1), u_{-1,-1}(r)⟩

Mp[Sp[spinorbar[p, 1, 1], γ, Sm[q], spinor[r, -1, -1]],
Sp[spinorbar[s, 1, -1], γ, spinor[t, 1, 1]]]
% // TraditionalForm
% // DoMToSpinorChains // Simplify // TraditionalForm
Mp[Sp[u_{1,1}[p], σ[1], Sm[q, -1], u_{-1,-1}[r]], Sp[u_{1,1}[t], σ[1], u_{1,-1}[s]]]
⟨u_{1,1}(p), σ^μ, Sm(q, -1), u_{-1,-1}(r)⟩ · ⟨u_{1,1}(t), σ_μ, u_{1,-1}(s)⟩
- 2 ⟨u_{1,1}(t), u_{-1,-1}(r)⟩ ⟨u_{1,1}(p), Sm(q, 1), u_{1,-1}(s)⟩ +
2 ⟨u_{1,-1}(s), u_{-1,-1}(r)⟩ ⟨u_{1,1}(t), Sm(q, 1), u_{1,1}(p)⟩ +
2 ⟨u_{1,1}(p), u_{-1,-1}(r)⟩ ⟨u_{1,1}(t), Sm(q, 1), u_{1,-1}(s)⟩

```

Figure 6.4: Examples of spinor chains containing gamma matrices and the simplification of the products with momenta and each other using the implementation in `6DSpinorHelicity.txt`.

products are fully expanded and if needed, reversed, to give a canonical representation. This canonicalisation is very important as otherwise it is possible to have a complex expression that should cancel, but does not, as the different terms are using different forms of their spinor chains or Minkowski products. It is not easy to find these manually in large expressions. Unfortunately there are still equivalent expressions that cannot be automatically canonicalised, as they would involve commuting matrices around in spinor chains, which if it does not cause a cancellation could drastically increase the number of terms in an expression, causing it to be too complex to handle and work with.

Complex, momenta like, expressions can also be written involving spinor chains with gamma matrices, `\[Gamma]` or `\[Sigma][h]`, inserted, which can be combined

using `Mp`; these and simplification of their products are shown in [Figure 6.4](#). If the other vector in the Minkowski product is a normal momenta rather than a spinor chain, then the product can be simplified by applying `DoMpToSm` to it. Simplifications for the Minkowski product of two spinor chains containing gamma matrices are harder to perform in general, but for most cases can be performed by applying `DoMpToSpinorChains`. The main case this function cannot replace is expressions of the form $\underline{u}_{\pm,h}(p)\gamma^\mu\underline{u}_{\pm,l}(q)\underline{u}_{\pm,m}(r)\gamma_\mu\underline{u}_{\pm,n}(s)$, as it is not possible to write them purely in terms of the spinors they contain. If this type of term is encountered, the simplest solution is to insert the fraction $\frac{\not{a}\not{b}+\not{b}\not{a}}{2a\cdot b}$ into one of the spinor chains, where a and b are two vectors which should be chosen so that the expression simplifies. As this requires knowledge of what the other vectors are, it cannot be automated and as such is not provided as a function.

Expressions involving chains of slashed matrices can often be simplified by commuting matrices so that objects for the same momenta are neighbouring, or so that different terms are rearranged to have the same form. The function `CommuteMatrices[element1,element2]`, where `element1` and `element2` are both either explicit gamma matrix objects of the form `\[Sigma][h]` or the name of a momenta, can be applied to an expression to commute `element1` and `element2` whenever they occur in that order in the expression. Basic automatic commutations are also possible where it is simple to see that the commutation will produce simpler expressions, as it will cause there to be two neighbouring slashed matrices for the same momenta, or cause a momenta to be next to its corresponding spinor, and as such these terms will vanish. This is done using the functions `CommuteMatricesAway` and `CommuteSigmaMatricesAway`, depending on whether it is a gamma matrix or only slashed matrices that are in between the matching items that will be commuted away. Examples of these functions for commuting slashed matrices are shown in [Figure 6.5](#).

These spinor and Minkowski products can also be evaluated numerically using functions defined in [6DSpinorHelicity.txt](#). There can be multiple sets of values defined for any momenta as the definitions are attached to a `tag`. To define a value for a momentum for a specific `tag` the function `DeclareVectorMomentum` is used, which takes arguments of the form `DeclareVectorMomentum[tag,momentum,{p0,p1,...}]` or `DeclareVectorMomentum[tag,momentum,{helicities}->spinor,...]`, where `helicities` is the helicities of the spinor and all possible helicities must be included in

```

Sp[spinorbar[p, 1, 1], Sm[r], Sm[s], Sm[t], spinor[q, 1, -1]] // CommuteMatrices[r, s]
Sp[spinorbar[p, 1, 1], Sm[r], Sm[q], Sm[r], spinor[q, 1, -1]] // CommuteMatricesAway
Sp[spinorbar[q, 1, -1] + spinorbar[p, 1, -1],
  (Sp[Sm[r], Sm[s]] + Sp[Sm[s], Sm[r]]), spinor[p, -1, -1]]
% // CommuteMatrices[r, s]

2 Mp[Mom[s], Mom[r]] Sp[u_{1,-1}[p], Sm[t, 1], u_{1,-1}[q]] -
  Sp[u_{1,-1}[p], Sm[s, 1], Sm[r, -1], Sm[t, 1], u_{1,-1}[q]]

2 Mp[Mom[r], Mom[q]] Sp[u_{1,-1}[p], Sm[r, 1], u_{1,-1}[q]]

Sp[u_{1,-1}[p], Sm[r, 1], Sm[s, -1], u_{1,-1}[p]] + Sp[u_{1,-1}[p], Sm[s, 1], Sm[r, -1], u_{1,-1}[p]] +
  Sp[u_{1,-1}[q], Sm[r, 1], Sm[s, -1], u_{1,-1}[p]] + Sp[u_{1,-1}[q], Sm[s, 1], Sm[r, -1], u_{1,-1}[p]]

2 Mp[Mom[s], Mom[r]] Sp[u_{1,-1}[q], u_{-1,-1}[p]]

```

Figure 6.5: Examples of commuting spinor chains using the implementation in `6DSpinorHelicity.txt`.

the list, otherwise missing options will not be usable for evaluating expressions. Once momenta are declared, then expressions using them can be evaluated by applying `Ev[dimensions, tag]`, where `dimensions` is the number of dimensions the expression should be evaluated for, which is required as some expressions can be used algebraically in any number of dimensions, but numerical evaluation requires the number of dimensions to be stated. Numerically evaluating various expressions along with the different ways of declaring momenta values are shown in [Figure 6.6](#).

The 4 dimensional spinors used in `6DSpinorHelicity.txt` as part of the recursive definition unfortunately do not have the same form as those used in `BlackHat` or `S@M`, but differ by a rotation of vectors in space. The rotation swaps the 2nd and 4th components and negates the 3rd component and is given by the Lorentz transformation

$$\Lambda_{\nu}^{\mu} = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 \\ 0 & 0 & -1 & 0 \\ 0 & 1 & 0 & 0 \end{pmatrix}. \quad (6.1)$$

This is especially important for the reduction of 6 dimensional momenta to massless 4 dimensional momenta, as the vector a used in the reduction was chosen for that particular spinor representation and needs to be converted if a different representation is used. For the `BlackHat` and `S@M` definition of spinors it is clear that a , as given

```

DeclareVectorMomentum [mom1, p, {sqrt[474], 3, 12, -4, -4, 17}];
DeclareVectorMomentum [mom1, q, {-sqrt[583], 11, -13, 2, 15, 8}];
DeclareVectorMomentum [mom2, p, {1 - 5 i, 3 + 2 i, 11 - 3 i, -7 i, 2 + 11 i, 5}];
DeclareVectorMomentum [mom2, q, {{1, 1} -> {{4 - 3 i}, {-3 - 18 i}, {16 - 2 i}, {0}},
    {1, -1} -> {{0}, {6/5 + 2 i/5}, {i}, {1}}, {-1, 1} -> {{-6/5 - 2 i/5}, {0}, {1}, {42/25 - 81 i/25}},
    {-1, -1} -> {{3 + 4 i}, {4 - 3 i}, {0}, {-16 + 2 i}}];
DeclareVectorMomentum [mom2, r, {5, 0, 3, 0, 0, 0, 0, 4}];

Mcm [p] // Ev [6, mom2]
Mcm [q] // Ev [6, mom2]
% == %
{1 - 5 i, 3 + 2 i, 11 - 3 i, -7 i, 2 + 11 i, 5}
{1 - 5 i, 3 + 2 i, 11 - 3 i, -7 i, 2 + 11 i, 5}
True

{Sp [spinorbar [p, 1, 1], spinor [r, -1, 1]],
 Sp [spinorbar [p, 1, -1], spinor [r, -1, 1]] // Ev [6, mom2] // Simplify
 {Sp [spinorbar [q, 1, 1], spinor [r, -1, 1]], Sp [spinorbar [q, 1, -1], spinor [r, -1, 1]] //
 Ev [6, mom2] // Simplify
 % == %
 %% [[1]] %% [[2]] == %% [[1]] %% [[2]]

{(-42/5 + 66 i/5) sqrt[4/5 - 3 i/5], (6 + 2 i) sqrt[4/5 - 3 i/5]}
{6 + 78 i/sqrt[5], 6 + 2 i/sqrt[5]}
False
True

Sm [p, h] // Ev [6, mom1] // TraditionalForm
Sm [p, h] // Ev [6, mom2] // TraditionalForm
(
sqrt[474] - 3 h    4 - 12 i    -17 + 4 i    0
-4 - 12 i    -3 h - sqrt[474]    0    -17 + 4 i
17 + 4 i    0    -3 h - sqrt[474]    -4 + 12 i
0    17 + 4 i    4 + 12 i    sqrt[474] - 3 h
)
(
(1 - 5 i) - (3 + 2 i) h    -3 - 4 i    6 - 2 i    0
-3 - 18 i    (-3 - 2 i) h - (1 - 5 i)    0    6 - 2 i
16 - 2 i    0    (-3 - 2 i) h - (1 - 5 i)    3 + 4 i
0    16 - 2 i    3 + 18 i    (1 - 5 i) - (3 + 2 i) h
)

```

Figure 6.6: Examples of declaring numeric values for momenta and evaluating spinors, momenta and slashed matrix objects in terms of them using the implementation in `6DSpinorHelicity.txt`.

in Equation 5.48, will be given by

$$a_{\text{BH}}^\mu = \frac{1}{2} \begin{pmatrix} 1 \\ 0 \\ 0 \\ -1 \end{pmatrix}. \quad (6.2)$$

These replacements for 6 dimensional spinors in terms of massless 4 dimensional spinors are provided in the variables `rep` and `reph[h]`, where in the latter the replacement uses `h` and its negative for the summed over helicities needed in the replacement, which can greatly help with producing a simple result. The reverse replacement is also provided in the variable `unrep`. These leave the expression defined in terms of the explicitly named extra component of the momenta p_4 and p_5 , which can then be replaced with the μ expressions defined in Equation 5.46. This is achieved by using the replacements stored in the variables `tomu`, `tomuh[h]` and `tomuhp[hfn]`, where the first replacement uses `+` and `-` as the signs in the μ , the second uses `+h` and `-h` as the signs and the third version's argument is a function that takes the name of the momenta and returns the expression to base the signs for the μ s on. Again the reverse replacement is provided and is stored in the variable `frommu`. To avoid replacing all momenta in an expression these replacements will currently only replace an explicitly specified list of vectors; the replacements will need to be manually updated if extra vectors need to be replaced, by changing the pattern named as `pp` in the replacement rules.

Due to time constraints the full calculation for 6 dimensional tree and one loop amplitudes has not been performed, but could be developed using the objects and tools defined in `6DSpinorHelicity.txt` along with the Feynman rules defined in Tables 2.1 and 5.1. Unfortunately the conversion to 4 dimensional spinors cannot be performed using replacement rules, as slashed matrices have the same form in any number of dimensions, but when converting from 6 to 4 dimensions should convert to the sum of both helicities, for which it is impossible to declare a rule. There are also issues with simplifications for spinor products applying during the replacement causing incorrect results and terms to vanish that should not have, since while in the middle of applying the replacement some objects are 6 dimensional and some are already 4 dimensional. As such the final conversion to 4 dimensional expressions will have to be done manually while writing the 4 dimensional implementation.

Chapter 7

Conclusions

In this thesis progress has been made towards adding Higgs boson amplitudes to BlackHat. This requires calculating tree amplitudes with any number of quarks and gluons both with and without a Higgs boson. These tree amplitudes have been calculated and are shown to agree with BlackHat, apart from a few differences in sign conventions for quarks. The cut parts have also been calculated and again agree with BlackHat, after taking account of the differences in quark signs, apart from a factor of -1 for closed quark loop contributions. BlackHat was seen to produce cuts that were known to vanish. The main case where this happened was amplitudes containing a Higgs boson where the corner with the Higgs boson causes the coefficient to vanish unless it is restricted to have at least two negative helicities and this is not currently done in BlackHat. Adding this restriction would save calculating a large class of cuts that can easily be seen to vanish. Apart from this possible improvement in speed the code was found to reliably and efficiently calculate the cut parts. The various restrictions and assumptions used in deriving the calculations and therefore the limitations of the method as currently implemented in this project are discussed. The main restriction on the method as discussed in this thesis is that only massless particles are allowed in loops. This restriction could be lifted and will have to be lifted when it is used in calculating the rational terms.

The method to calculate the rational terms was unfortunately not completed in this project due to time constraints so could not be implemented into BlackHat. The basic method is discussed in [Chapter 5](#) and a Mathematica implementation of 6 dimensional spinor helicity formalism has been developed, along with a method to reduce the calculations back down to 4 dimensions by introducing effective massive particles

which is discussed in [Chapter 6](#). This implementation is included in [Appendix C](#). The general form of the dependence of tree amplitudes needed for generalised unitarity calculations on the extra dimensions has been investigated and there is no dependence found on the direction of the extra components, apart from in the amplitudes needed to subtract the extra gluon polarisation states. For those amplitudes it is required that the product of the loop momenta with the polarisation vector in the 6th dimension is taken to be zero. If this is ensured, even if the direction of the polarisation vector has to be effectively chosen based on the direction of the loop momenta, then again the amplitudes are independent of the direction of the extra components. With these restrictions the rational terms can be calculated in terms of only 4 dimensional momenta and amplitudes with new effective massive particles, however the internal degrees of freedom in the Feynman rules still require full 6 dimensional momenta and states.

Once a set of the simplest amplitudes has been calculated, higher multiplicity amplitudes can again be calculated with the BCFW recursion relations, but this time using massive 4 dimensional momenta. The general form has been investigated, but the details of the implementation, including which particles are valid to use as the shifted particles and what other factors need to be introduced into each term other than the pair of tree amplitudes, has not been investigated. Once all tree amplitudes can be calculated the scalar loop integral coefficients needed for the rational terms can be extracted using generalised unitarity. The forms of the numerator terms in terms of the degrees of freedom in the loop momenta definitions needs calculating, along with the subtraction terms needed to subtract the higher order terms from the lower order terms, which must be formulated in a way that can be evaluated numerically. Once these elements of the calculation are completed, the rational terms can be calculated. Finally, these calculations will need to be implemented into BlackHat, which will also require converting the calculation to use the conventions used in BlackHat and implementing it in terms of BlackHat's choice of spinor conventions. These combined with the already implemented cut part calculations will provide an efficient and automatic calculation for the loop amplitude in a way that will make BlackHat an even more useful addition to the set of tools available for NLO calculations at the LHC. This thesis provides an important set of building blocks towards this aim.

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Appendix A

The Form of a Generic Term in an Amplitude.

The first case to prove is the generic term in a pure gluon amplitude which is given by

$$C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j^{\mu_j} \prod_{k=1}^n g^{\nu_k \sigma_k}}{\prod_{l=1}^{N_g+m-n-2} S_l^2} . \quad (\text{A.1})$$

This can be proven to be the correct form using proof by induction. Firstly the base cases are a single three or four gluon vertex on its own which corresponds to the cases where $m = 0$, $n = 1$ and $N_g = 3$ or $m = 0$, $n = 2$ and $N_g = 4$ respectively. To prove the recursion this form must be combined with each possible form of propagator and extra vertex combination and show that all terms in the result are also of this form. Firstly the case of combining a momentum index with the metric tensor from a three

gluon vertex is considered. The relation is then given by

$$\begin{aligned}
& C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j^{\mu_j} \prod_{k=1}^n g^{\nu_k \sigma_k}}{N_g+m-n-2 \prod_{l=1} S_l^2} \frac{i}{k^2+i0} \frac{-i}{\sqrt{2}} g^{\mu_1 \mu'_1} (k_1 - k_2)^{\mu(N_g-2n+1)} \\
&= \frac{C}{\sqrt{2}} \frac{\prod_{i=1}^m P_i \cdot Q_i \left(\prod_{j=2}^{N_g-2n} R_j^{\mu_j} \right) R_1^{\mu'_1} (k_1 - k_2)^{\mu(N_g-2n+1)} \prod_{k=1}^n g^{\nu_k \sigma_k}}{k^2 \prod_{l=1}^{N_g+m-n-2} S_l^2} \\
&= C' \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N'_g-2n} R'_j{}^{\mu'_j} \prod_{k=1}^n g^{\nu_k \sigma_k}}{N'_g+m-n-2 \prod_{l=1} S_l'^2}, \tag{A.2}
\end{aligned}$$

where in the last line the new factors have been relabelled to match the terms they combine with and N_g has been replaced with $N'_g = N_g + 1$ which is the correct increase in external gluons. The next case is when a metric tensor in the term is contracted with a metric tensor of the extra three gluon vertex and is given by

$$\begin{aligned}
& C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j^{\mu_j} \prod_{k=1}^n g^{\nu_k \sigma_k}}{N_g+m-n-2 \prod_{l=1} S_l^2} \frac{i}{k^2+i0} \frac{-i}{\sqrt{2}} g^{\sigma_1 \sigma'_1} (k_1 - k_2)^{\mu(N_g-2n+1)} \\
&= \frac{C}{\sqrt{2}} \frac{\prod_{i=1}^m P_i \cdot Q_i \left(\prod_{j=1}^{N_g-2n} R_j^{\mu_j} \right) (k_1 - k_2)^{\mu(N_g-2n+1)} \left(\prod_{k=2}^n g^{\nu_k \sigma_k} \right) g^{\nu_1 \sigma'_1}}{k^2 \prod_{l=1}^{N_g+m-n-2} S_l^2} \\
&= C' \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N'_g-2n} R'_j{}^{\mu'_j} \prod_{k=1}^n g^{\nu_k \sigma'_k}}{N'_g+m-n-2 \prod_{l=1} S_l'^2}. \tag{A.3}
\end{aligned}$$

The next case is contracting the momenta from a three gluon vertex with a metric tensor from the term and is given by

$$\begin{aligned}
& C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j^{\mu_j} \prod_{k=1}^n g^{\nu_k \sigma_k}}{N_g+m-n-2} \frac{i}{k^2+i0} \frac{-i}{\sqrt{2}} g^{\nu'_1 \sigma'_1} (k_1 - k_2)_{\sigma_1} \\
& \quad \prod_{l=1} S_l^2 \\
&= \frac{C}{\sqrt{2}} \frac{\prod_{i=1}^m P_i \cdot Q_i \left(\prod_{j=1}^{N_g-2n} R_j^{\mu_j} \right) (k_1 - k_2)^{\nu_1} \left(\prod_{k=2}^n g^{\nu_k \sigma_k} \right) g^{\nu'_1 \sigma'_1}}{k^2 \prod_{l=1}^{N_g+m-n-2} S_l^2} \\
&= C' \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N'_g-2n} R_j^{\mu'_j} \prod_{k=1}^n g^{\nu'_k \sigma'_k}}{N'_g+m-n-2} \prod_{l=1} S_l'^2 . \tag{A.4}
\end{aligned}$$

The last case for adding a three gluon vertex is contracting the momenta from the vertex with a momenta in the term which results in

$$\begin{aligned}
& C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j^{\mu_j} \prod_{k=1}^n g^{\nu_k \sigma_k}}{N_g+m-n-2} \frac{i}{k^2+i0} \frac{-i}{\sqrt{2}} g^{\nu'_{n+1} \sigma'_{n+1}} (k_1 - k_2)_{\mu_1} \\
& \quad \prod_{l=1} S_l^2 \\
&= \frac{C}{\sqrt{2}} \frac{\left(\prod_{i=1}^m P_i \cdot Q_i \right) R_1 \cdot (k_1 - k_2) \prod_{j=2}^{N_g-2n} R_j^{\mu_j} \left(\prod_{k=1}^n g^{\nu_k \sigma_k} \right) g^{\nu'_{n+1} \sigma'_{n+1}}}{k^2 \prod_{l=1}^{N_g+m-n-2} S_l^2} \\
&= C' \frac{\prod_{i=1}^{m'} P'_i \cdot Q'_i \prod_{j=1}^{N'_g-2n'} R'_j{}^{\mu'_j} \prod_{k=1}^{n'} g^{\nu'_k \sigma'_k}}{N'_g+m'-n'-2} \prod_{l=1} S_l'^2 , \tag{A.5}
\end{aligned}$$

where in the last line as well as replacing N_g with $N'_g = N_g + 1$, n has been replaced with $n' = n + 1$ and m has been replaced with $m' = m + 1$ which all agree with the expected limits on the different factors.

Lastly for the pure gluon amplitude the contraction of a four gluon vertex with a term must be shown to produce correctly formatted terms. Again each case will be

handled separately. Firstly, if the vertex is contracted with a metric tensor from the term then the combination is given by

$$\begin{aligned}
& C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j^{\mu_j} \prod_{k=1}^n g^{\nu_k \sigma_k}}{N_g+m-n-2} \frac{i}{k^2 + i0} c g_{\sigma_1}^{\sigma'_1} g^{\nu'_{n+1} \sigma'_{n+1}} \\
& \quad \prod_{l=1} S_l^2 \\
& = icC \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j^{\mu_j} \left(\prod_{k=2}^n g^{\nu_k \sigma_k} \right) g^{\nu_1 \sigma'_1} g^{\nu'_{n+1} \sigma'_{n+1}}}{N_g+m-n-2} \\
& \quad \frac{k^2}{\prod_{l=1} S_l^2} \\
& = C' \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N'_g-2n'} R_j^{\mu_j} \prod_{k=1}^{n'} g^{\nu'_k \sigma'_k}}{N'_g+m-n'-2} \frac{i}{k^2 + i0} c g_{\mu_1}^{\mu'_1} g^{\nu'_{n+1} \sigma'_{n+1}} \\
& \quad \prod_{l=1} S_l'^2, \tag{A.6}
\end{aligned}$$

where now N_g is replaced by $N'_g = N_g + 2$ and n is replaced by $n' = n + 1$. The last case for pure gluon amplitudes is contracting the four gluon vertex with a momenta in the term which gives

$$\begin{aligned}
& C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j^{\mu_j} \prod_{k=1}^n g^{\nu_k \sigma_k}}{N_g+m-n-2} \frac{i}{k^2 + i0} c g_{\mu_1}^{\mu'_1} g^{\nu'_{n+1} \sigma'_{n+1}} \\
& \quad \prod_{l=1} S_l^2 \\
& = icC \frac{\prod_{i=1}^m P_i \cdot Q_i \left(\prod_{j=2}^{N_g-2n} R_j^{\mu_j} \right) R_1^{\mu'_1} \left(\prod_{k=1}^n g^{\nu_k \sigma_k} \right) g^{\nu'_{n+1} \sigma'_{n+1}}}{N_g+m-n-2} \\
& \quad \frac{k^2}{\prod_{l=1} S_l^2} \\
& = C' \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N'_g-2n'} R_j^{\mu_j} \prod_{k=1}^{n'} g^{\nu'_k \sigma'_k}}{N'_g+m-n'-2} \frac{i}{k^2 + i0} c g_{\mu_1}^{\mu'_1} g^{\nu'_{n+1} \sigma'_{n+1}} \\
& \quad \prod_{l=1} S_l'^2. \tag{A.7}
\end{aligned}$$

The changes in number of factors of each type that can be introduced by adding each type of vertex are summarised in [Table A.1](#).

Now the form for amplitudes with a single quark line will be derived which is given

Terms of the form	Formula	Three Gluon Vertex		Four Gluon Vertex
$P \cdot Q$	m		+1	
R^μ	$N_g - 2n$	+1	-1	
$g^{\mu\nu}$	n		+1	+1
$\frac{1}{S^2}$	$N_g + m - n - 2$	+1	+1	+1
Number of gluons	N_g	+1	+1	+2

Table A.1: The different forms of factors that can appear in pure gluon amplitudes along with how the number of each factor present is changed by adding extra propagator and vertex combinations of each type.

by

$$C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g - N_{qg} - 2n} R_j \cdot \epsilon_j \prod_{k=1}^n \epsilon_k \cdot \epsilon'_k}{\prod_{l=1}^{N_g + m - n - N_{qR} - 1} S_l^2} \bar{u} \left(\prod_a^{N_{qR} + N_{qg} - 1} \psi_a \not{T}_a \right) \psi_{(N_{qR} + N_{qg})} u, \quad (\text{A.8})$$

where N_{qg} of the U_a s are gluon polarisation vectors and the rest (N_{qR} of them) along with T_a are momentum vectors containing some combination of neighbouring momenta. To derive this formula, the form is built by tracing along the quark line from the spinor towards the conjugate spinor and then at the end adding the conjugate spinor for the other quark to the end of the quark line. The form used to build this up is therefore the equation given above with the conjugate spinor, \bar{u} , removed which is given by

$$C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g - N_{qg} - 2n} R_j \cdot \epsilon_j \prod_{k=1}^n \epsilon_k \cdot \epsilon'_k}{\prod_{l=1}^{N_g + m - n - N_{qR} - 1} S_l^2} \left(\prod_a^{N_{qR} + N_{qg} - 1} \psi_a \not{T}_a \right) \psi_{(N_{qR} + N_{qg})} u. \quad (\text{A.9})$$

The base cases for the proof by induction are therefore the quark combined with a single quark vertex which is itself combined with a gluon propagator and then a term for a combination of gluons, which will be of the form found above for the pure gluon amplitude. There are two different cases depending on whether the gamma matrix introduced by the quark vertex is contracted with a momentum from the gluon term or a metric tensor and therefore eventually with a gluon. If the gamma matrix is

contracted with a momentum vector then the terms will have the form of

$$\begin{aligned}
& C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{(N_g+1)-2n-1} R_j^{\mu_j} \prod_{k=1}^n g^{\nu_k \sigma_k}}{(N_g+1)+m-n-2} \frac{i}{k^2 + i0} \mathcal{R}_{((N_g+1)-2n)u} \\
& \quad \prod_{l=1} S_l^2 \\
& = C' \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j^{\mu_j} \prod_{k=1}^n g^{\nu_k \sigma_k}}{N_g+m-n-1} \mathcal{R}_{(N_g-2n+1)u} \\
& \quad k^2 \prod_{l=1} S_l^2 \\
& = C' \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j^{\mu_j} \prod_{k=1}^n g^{\nu_k \sigma_k}}{N_g+m-n} \mathcal{V}_1 u, \tag{A.10} \\
& \quad \prod_{l=1} S_l^2
\end{aligned}$$

where $N_g + 1$ is used where N_g would have been used in the pure gluon amplitude as one gluon from the factor is replaced with the quark line. This matches the expected form for when $N_{qR} = 1$ and $N_{qg} = 0$.

$$\begin{aligned}
& C \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{(N_g+1)-2n} R_j^{\mu_j} \prod_{k=1}^{n-1} g^{\nu_k \sigma_k}}{(N_g+1)+m-n-2} \frac{i}{k^2 + i0} \gamma^{\sigma_n} u \\
& \quad \prod_{l=1} S_l^2 \\
& = C' \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n} R_j^{\mu_j} \prod_{k=1}^{n-1} g^{\nu_k \sigma_k}}{N_g+m-n-1} \gamma^{\sigma_n} u \\
& \quad k^2 \prod_{l=1} S_l^2 \\
& = C' \frac{\prod_{i=1}^m P_i \cdot Q_i \prod_{j=1}^{N_g-2n'-2} R_j^{\mu_j} \prod_{k=1}^{n'} g^{\nu_k \sigma_k}}{N_g+m-n'-1} \gamma^{\sigma_n} u, \tag{A.11} \\
& \quad \prod_{l=1} S_l'^2
\end{aligned}$$

where in the last line n has been replaced with $n' = n - 1$. Again this matches with the expected form, but now for when $N_{qR} = 0$ and $N_{qg} = 1$. The recursion relations are derived by adding an extra quark propagator and quark vertex with the quark vertex contracted with a gluon propagator and a pure gluon term. The formulas are too complex to show here but the changes induced by each combination are shown in [Table A.2](#).

Terms of the Form	Formula	Type of Vector that is Contracted into the Quark Line	
		Polarisation Vector	Momenta
$P \cdot Q$	m	$+m'$	$+m'$
R^μ	$N_g - 2n - N_{qg}$	$+N'_g - 2n' - 1$	$+N'_g - 2n'$
$g^{\mu\nu}$	n	$+n'$	$+n'$
$\frac{1}{S^2}$	$N_g + m - n + N_{qR} - 1$	$+N'_g + m' - n'$	$+N'_g + m' - n' + 1$
Number of gluons	N_g	$+N'_g$	$+N'_g$
Number of Polarisation Vectors in the Quark Line	N_{qg}	$+1$	
Number of Momenta in the Quark Line	$2N_{qR} + N_{qg}$	$+1$	$+2$

Table A.2: The different forms of factors that can appear in amplitudes with one quark line along with how the number of each factor present is changed by adding extra propagator and vertex combinations of each type.

Appendix B

Source Code for 4

Dimensional Calculations and Comparing to BlackHat

The code created as part of this project can also be downloaded from <http://bit.ly/2oXSBSu>.

B.1 Mathematica Implementation of Tree and One Loop Cut Part Calculations

Listing B.1: Common.txt

```

1 (* Functions to handle a list with wrapping. Handles the modular
   arithmetic with the range of indices being 1 to length rather than
   the more common 0 to length - 1 *)
2 LLMod[l_List, i_Integer] := 1 + Mod[i - 1, Length[l]]
3 LL[l_List, i_Integer] := 1[[LLMod[l, i]]]
4 LL[l_List, i_List] := (LL[l, #1] & ) /@ i
5 LLRange[l_List, i_Integer, j_Integer] :=
6 (LLMod[l, #1] & ) /@ Range[i, i + Mod[j - i, Length[l]]]
7
8 (* Load S@M *)
9 << "Spinors"
10
11 (* Silence the many messages that S@M prints *)

```

```

12 Spinors `Private `PRINT[x_..] := Null
13
14 (* Add an undeclare method *)
15 Spinors `UndeclareSpinorMomentum[s_] := (Spinors `Private `NumLa[s] =.;
16   Spinors `Private `NumCLa[s] =.; Spinors `Private `NumLat[s] =.;
17   Spinors `Private `NumCLat[s] =.; Unprotect[Num4V]; Num4V[s] =.;
18   Protect[Num4V];
19   Spinors `Private `NumSpinorList :=
20     Evaluate[DeleteCases[Spinors `Private `NumSpinorList, s]];
21   Spinors `Private `NumDefs[])
22 Spinors `FullyUndeclareSpinor[s_] := (UndeclareSpinorMomentum[s];
23   UndeclareSpinor[s]);
24
25 (* Declare an improved version of the spinor definition formula that
    handles Indeterminates and Complex Infinities *)
26 Spinors `Private `Sol2Dim1[k_List] :=
27   Module[{sqk1p = Sqrt[k[[1]] + k[[4]]], ptplus = k[[2]] + I k[[3]]},
28   If[MemberQ[k, Indeterminate | ComplexInfinity], {Indeterminate,
29     Indeterminate},
30   If[sqk1p == 0,
31     If[k[[1]] - k[[4]] ==
32       0, {(k[[2]] - I k[[3]])/Sqrt[2 k[[2]]], (k[[2]] + I k[[3]])/
33         Sqrt[2 k[[2]]]}, {(k[[2]] - I k[[3]])/Sqrt[k[[1]] - k[[4]]],
34         Sqrt[k[[1]] - k[[4]]]}], {sqk1p, ptplus/sqk1p}]]]
35 Spinors `Private `Sol2Dim2[k_List] :=
36   Module[{factor = 1, sqk1p = Sqrt[k[[1]] + k[[4]]],
37   ptminus = k[[2]] - I k[[3]]},
38   If[MemberQ[k, Indeterminate | ComplexInfinity], {Indeterminate,
39     Indeterminate},
40   If[sqk1p == 0,
41     If[k[[1]] - k[[4]] ==
42       0, {(k[[2]] + I k[[3]])/Sqrt[2 k[[2]]], (k[[2]] - I k[[3]])/
43         Sqrt[2 k[[2]]]}, {(k[[2]] + I k[[3]])/Sqrt[k[[1]] - k[[4]]],
44         Sqrt[k[[1]] - k[[4]]]}], {sqk1p, ptminus/sqk1p}]]]
45
46 DeclareSpinor[Epsilon[...]]
47
48 (* Use a different sign convention. Can be set to true after this is
    loaded and before calculation is performed *)
49 $UseBHSigns=False
50
51 (* Set the precision if it is not already set *)

```

```

52 If[Hold[Evaluate[$Precision]] == Hold[$Precision], $Precision =
53   MachinePrecision]
54
55 (* Allows increased precision in internal parts of the calculation *)
56 NN[e_] := N[e, $Precision]
57 NN[e_, f : (_Integer | _Real)] := N[e, f $Precision]
58 NN[e_, Max] := N[e, 10 MachinePrecision]
59
60 (* Set the symbol s so that it evaluates to the numeric value of the
61    expression v at f times the current precision but only actually
62    evaluates once per precision at which it is used *)
61 NNCache[s_Symbol, v_, f_ : 1] :=
62   s := (Module[{vv = Evaluate[NN[v, f]], p = $Precision},
63     s := vv /; p == $Precision]; s)
64
65
66 (* Only declare the spinor momentum if the elements of the momenta or
67    spinor are numeric *)
67 MyDeclareSpinorMomentum[a_,
68   xx : {_?NumericQ, _?NumericQ, _?NumericQ, _?NumericQ}] :=
69   DeclareSpinorMomentum[a, xx]
70 MyDeclareSpinorMomentum[
71   a_, {la : {_?NumericQ}, {_?NumericQ}},
72   lat : {{_?NumericQ, _?NumericQ}}] :=
73   DeclareSpinorMomentum[a, la, lat]
74
75 (* Evaluate an expression with the specified sets of definitions for
76    momenta. Will remain unevaluated if the spinor elements are not
77    numeric *)
76 WithSpinors[expr_,
77   spinors : {_, ({_?NumericQ, _?NumericQ, _?NumericQ, _?
78     NumericQ} | {{_?NumericQ}, {_?NumericQ}}, {{_?NumericQ, _?
79     NumericQ}})} ...] :=
80   Module[{tmp}, MyDeclareSpinorMomentum @@@ {spinors}; tmp = expr;
81     FullyUndeclareSpinor[#[[1]]] & /@ {spinors}; tmp];
82 SetAttributes[WithSpinors, HoldFirst]
83 SyntaxInformation[WithSpinors] = {"ArgumentsPattern" -> {_, ---},
84   "LocalVariables" -> {"Table", {2, \[Infinity]}}]
85
86 (* Declare various properties for the different types of particle *)
87 ParticleName[Gluon[h_]] := "Gluon helicity:" <> ToString[h]
88 ParticleName[Phi[1]] := "Phi"

```

```

89 ParticleName [Phi[-1]]:= "Phi dagger"
90 ParticleName [Higgs []]:= "Higgs boson"
91 ParticleName [Quark [1, h_, f_]]:= "Quark of flavour:" <> ToString[h] <> " and
    helicity:" <> ToString[h]
92 ParticleName [Quark [-1, h_, f_]]:= "Anti Quark of flavour:" <> ToString[h] <> "
    and helicity:" <> ToString[h]
93
94 ReverseParticle [phi_Phi]:= phi
95 ReverseParticle [higgs_Higgs]:= higgs
96 ReverseParticle [Gluon [h_]]:= Gluon[-h]
97 ReverseParticle [Quark [p_, h_, f_]]:= Quark[-p, -h, f]
98
99 IsOrderedQ [_Phi | _Higgs]:= False
100 IsOrderedQ [_Gluon | _Quark]:= True
101 IsUnorderedQ [p_]:= !IsOrderedQ [p]
102
103 IsMassiveQ [_Phi | _Higgs]:= True
104 IsMassiveQ [_Gluon | _Quark]:= False
105
106 (* Skip evaluating if the condition is false *)
107 SetAttributes [SkipIf, HoldFirst]
108 SkipIf [val_, False] := val
109 SkipIf [-, True] := Sequence []
110
111 (* Provide a sort key that uniquely sorts particle types into a format
    that gives convenient orderings *)
112 SortKey [Gluon [h_]]:= {0, h, 0, 0, 0}
113 SortKey [Quark [p_, h_, f_]]:= {-1, 0, 0, 0, 0}
114 SortKey [Phi [p_]]:= {100, -p, 0, 0, 0}
115 SortKey [Higgs []]:= {100, 0, 0, 0, 0}

```

Listing B.2: HelAmplN.txt

```

1 (* Split the particles into separate groups for the colour ordered and
    non-colour ordered particles *)
2 HelAmplN [p : {---}, types : {Except[_List] ...}] :=
3 HelAmplN[{Extract [p, Position [types, _?IsUnorderedQ]],
4 Extract [p, Position [types, _?IsOrderedQ]]}, {Cases [
5 types, _?IsUnorderedQ], Cases [types, _?IsOrderedQ]}]
6
7 (* Put the particles into a canonical order *)
8 HelAmplN[{pNonOrd_List, p_List}, {typesNonOrd_List, types_List}] :=
9 Module[{bestI =

```

```

10   Sort[Table[{RotateLeft[SortKey /@ types, i], i}, {i, 0,
11     Length[types] - 1}][[1, 2]]],
12   HelAmplN[{pNonOrd, RotateLeft[p, bestI]}, {typesNonOrd,
13     RotateLeft[types, bestI]}] /; bestI != 0]
14
15 (* Allow the momenta for the particles to be passed in directly rather
    than requiring previously declared momenta names *)
16 HelAmplN[{pVNonOrd : {_List ...},
17   pV : {_List ...}}, {typesNonOrd_List, types_List}] :=
18 Module[{isMassive = IsMassiveQ /@ Join[typesNonOrd, types], n, p, P,
19   allp}, n = Plus @@ isMassive /. {False -> {0, 1}, True -> {1, 0}};
20   p = Range[n[[2]]];
21   P = Table[Symbol["PP" <> ToString[i]], {i, n[[1]]}];
22   allp = isMassive[[_; _]];
23   allp[[_Position[isMassive, True] // Flatten]] = P;
24   allp[[_Position[isMassive, False] // Flatten]] = p;
25   MapThread[
26     If[#1, DeclareSpinorMomentum,
27       DeclareLVectorMomentum][#2, #3] &, {isMassive, allp,
28       Join[pVNonOrd, pV]}, 1];
29   HelAmplN[{allp[[_; Length[pVNonOrd]]],
30     allp[[_Length[pVNonOrd] + 1 _;]]}, {typesNonOrd, types}]
31
32 (* Declare an error message format for use in amplitudes *)
33 HelAmplN::badamp =
34   "Bad amplitude evaluated for particles of types ‘‘ called ‘‘ has \
35   bad value ‘‘";
36
37 (* Declare the vanishing, MHV and anti-MHV pure gluon amplitudes *)
38 HelAmplN[{{},
39   p_List}, {{}, {Gluon[-1] | PatternSequence[], Gluon[1], Gluon[1],
40     Gluon[1] ..}}] := 0
41 HelAmplN[{{},
42   p_List}, {{}, {Gluon[-1], Gluon[-1], Gluon[-1] ..,
43     Gluon[1] | PatternSequence[]}}] := 0
44
45 HelAmplN[{{}, p_List}, {{},
46   types : {Gluon[-1], Gluon[-1], Gluon[1]}}] :=
47 I*Module[{num = Spaa[p[[1]], p[[2]]] // NN,
48   denom = Product[Spaa[LL[p, i], LL[p, i + 1]], {i, Length[p]}] //
49   NN}, If[num == 0 && denom == 0,
50   Message[HelAmplN::badamp, types, p,

```



```

51   HoldForm[Num^4/Denom] /. {Num -> num, Denom -> denom}; 0,
52   num^4/denom]]
53 HelAmplN[{{}, p-List}, {{},
54   types : {Gluon[-1], Gluon[1], Gluon[1]}] :=
55   I*(-1)^Length[p]*
56   Module{num = Spbb[p[[2]], p[[-1]]] // NN,
57   denom = Product[Spbb[LL[p, i], LL[p, i + 1]], {i, Length[p]}] //
58   NN}, If[num == 0 && denom == 0,
59   Message[HelAmplN::badamp, types, p,
60   HoldForm[Num^4/Denom] /. {Num -> num, Denom -> denom}; 0,
61   num^4/denom]]
62
63 HelAmplN[{{}, p-List}, {{}, {Gluon[-1], gp1:(Gluon[1]...), Gluon[-1], Gluon
64   [1]...}}] :=
65   I*Spaa[p[[1]], p[[Length{gp1}+2]]]^4/
66   Product[Spaa[LL[p, i], LL[p, i + 1]], {i, Length[p]}]//NN
67 HelAmplN[{{}, p-List}, {{}, {gm1:Gluon[-1]..., Gluon[1], gm2:Gluon[-1]...,
68   Gluon[1]}]}] :=
69   I*(-1)^Length[p]*Spbb[p[[Length{gm1}+1]], p[[-1]]]^4/
70   Product[Spbb[LL[p, i], LL[p, i + 1]], {i, Length[p]}]//NN
71
72 (* Declare the vanishing, MHV and anti-MHV amplitudes with one quark
73   pair *)
74 HelAmpl[{{}, -List}, {{}, {Quark[-, -, -], Quark[-, -, -],
75   Gluon[hg-], Gluon[hg- ..]}}, - : Null] := 0
76 HelAmpl[{-List, -List}, {-List, {Quark[pq1-, hq1-, f1-],
77   Quark[pq2-, hq2-, f2-], Gluon[-] ...}} - : Null] :=
78   0 /; pq1 == pq2 || hq1 == hq2 || f1 != f2
79
80 HelAmplN[{-, -}, {-List, {-Quark, -Quark}}] := 0
81
82 HelAmplN[{{}, p : {pq1-, pq2-, pg-}},
83   types : {{}, {Quark[Pq1-, hq1-, f-], Quark[Pq2-, hq2-, f-],
84   Gluon[hg-]}}, epsilon- : Epsilon[]] :=
85   If[$UseBHSigns, 1, -Pq1 hq1] If[hq1 == 1, -(-1)^Length[p], 1] I*
86   Module{num =
87   If[hg == hq1, If[hg == -1, Spaa[pg, pq1], Spbb[pg, pq1]],
88   If[hg == -1, Spaa[pg, pq2], Spbb[pg, pq2]]] // NN,
89   denom = If[hg == -1, Spaa[pq2, pq1], Spbb[pq2, pq1]] // NN},
90   If[num == 0 && denom == 0,
91   Message[HelAmplN::badamp, types, p,
92   HoldForm[Num^2/Denom] /. {Num -> num, Denom -> denom}; 0,

```

```

90     num^2/denom]] /; Pq1 == -Pq2 && hq1 == -hq2
91
92 HelAmplN2QuarksH1H3[Sptype_, h1_, h3_, {p1_, p2_, p3_}] :=
93 NN[Sptype[p1, p3]]^2 /; h1 == h3
94 HelAmplN2QuarksH1H3[Sptype_, h1_, h3_, {p1_, p2_, p3_}] :=
95 NN[Sptype[p2, p3]]^2 /; h1 == -h3
96
97 HelAmplN2Quarks[Sptype_, h1_, h3_, {p1_, p2_, p3_}, p-List] :=
98 I*HelAmplN2QuarksH1H3[Sptype, h1, h3, {p1, p2, p3}]*
99 NN[Sptype[p1, p3]*
100   Sptype[p2, p3]/
101   Product[Sptype[LL[p, i], LL[p, i + 1]], {i, Length[p]}]]
102
103 HelAmplN[{{}},
104   p : {pq1_, pq2_, pgs_}, {{}, {Quark[Pq1_, hq1_, f_],
105     Quark[Pq2_, hq2_, f_], go1 : Gluon[hog_] ..., Gluon[hg_],
106     Gluon[hog_] ...}}, epsilon_ : Epsilon[], hg_ | PatternSequence[]] :=
107 If[$UseBHSigs, 1, -Pq1*hq1]*If[hg == 1, -(-1)^Length[p], 1]*
108 HelAmplN2Quarks[If[hg == -1, Spaa, Spbb], hq1,
109   hg, {pq1, pq2, {pgs}[Length[{go1}] + 1]}, p] /; Pq1 == -Pq2 && hq1 ==
110   hq2 && hog == -hg
111
112 (* Declare the vanishing, MHV and anti-MHV amplitudes with a phi *)
113 HelAmplN[{{-}, p-List], {{Phi[1]}, {Gluon[-], Gluon[1]...}}] := 0
114 HelAmplN[{{-}, p-List], {{Phi[1]}, {Gluon[-1], gp1 : (Gluon[1]...), Gluon[-1],
115   Gluon[1]...}}] :=
116 I*Spaa[p[[1]], p[[Length[{gp1}] + 2]]]^4/
117 Product[Spaa[LL[p, i], LL[p, i + 1]], {i, Length[p]}]//NN
118 HelAmplN[{{-}, p-List], {{Phi[1]}, {Gluon[-1]..}}] :=
119 I*(-1)^Length[p]*s @@ p^2/Product[Spbb[LL[p, i], LL[p, i + 1]],
120   {i, Length[p]}]//NN
121
122 (* Declare the 3 negative, Next-to-MHV amplitude. This uses an arbitrary
123   reference vector which is generated at random *)
124 HelAmplPhiRatioSum[p_, l1_, l2_, qs_, qe_, r_] :=
125 HelAmplPhiRatioSum[p, LLMOD[p, l1], LLMOD[p, l2], LLMOD[p, qs],
126   LLMOD[p, qe], r] /; l1 > Length[p] || l2 > Length[p] ||
127   qs > Length[p] || qe > Length[p]
128 HelAmplPhiRatioSum[p_, e_, e_, e_, e_, r_] := 1
129 HelAmplPhiRatioSum[p_, l1_, l2_, qs_, qe_, r_] :=
130 Sum[Spab[LL[p, l1], LL[p, qi], r], {qi, LLRange[p, qs, qe]}/
131 Sum[Spab[LL[p, l2], LL[p, qi], r], {qi, LLRange[p, qs, qe]}]

```

```

129 HelAmplPhi3NegTerm[p_, {m1_, m2_, m3_}, i_, j_, qs_, qe_, epsilon_] :=
130 (Spaa[LL[p, m2], LL[p, m3]]^4*Spaa[LL[p, i], LL[p, i + 1]]*
131   Spaa[LL[p, j], LL[p, j + 1]]*HelAmplPhiRatioSum[p, m1, i, qs, qe,
132     epsilon]*HelAmplPhiRatioSum[p, m1, j, qs, qe, epsilon]*
133   HelAmplPhiRatioSum[p, m1, i + 1, qs, qe, epsilon]*
134   HelAmplPhiRatioSum[p, m1, j + 1, qs, qe, epsilon])/
135   s @@ LL[p, LLRange[p, qs, qe]]
136 HelAmplPhi3Neg[p_, {m1_, m2_, m3_}, epsilon_] :=
137 Sum[HelAmplPhi3NegTerm[p, {m1, m2, m3}, i, j, i + 1, j, epsilon] +
138   HelAmplPhi3NegTerm[p, {m1, m2, m3}, i, j, j + 1, i, epsilon],
139   {i, LLRange[p, m1, m2 - 1]}, {j, LLRange[p, m3, m1 - 1]}]
140
141 RandomNSphere[n_] :=
142 Module{tmp = RandomVariate[MultinormalDistribution[
143   ConstantArray[0, n], IdentityMatrix[n]]]},
144   tmp/Sqrt[Plus @@ (tmp*tmp)]]
145 RandomNBall[n_] := Random[Real]^(1/n)*RandomNSphere[n]
146 Random4Vector[Real, m_:0, scale_:1] :=
147 Module{tmp = scale*RandomNBall[3]},
148   Prepend[tmp, (2*Random[Integer] - 1)*Sqrt[m^2 + Plus @@ (tmp*tmp)]]]
149
150 NWithEpsilons[expr_] :=
151 Module{es = DeleteDuplicates[Cases[expr, Epsilon[---],
152   100]]}, PRINT["Generating Momenta for", es];
153   (DeclareSpinorMomentum[#1, Random4Vector[Real]] & ) /@ es; N[expr]]
154
155 HelAmplN[{{_}, p_List}, {{Phi[1]}, {Gluon[-1], gp1:Gluon[1]... , Gluon[-1],
156   gp2:Gluon[1]... , Gluon[-1], Gluon[1]... }}] :=
157 NWithEpsilons[I*Sum[HelAmplPhi3Neg[p, RotateLeft[{1, Length[{gp1}]+2,
158   Length[{gp1, gp2}]+3}, i], Epsilon[[]], {i, 3}]/
159   Product[Spaa[LL[p, i], LL[p, i + 1]], {i, Length[p]}]]]
160
161 (* Will return all ways of splitting a list of unique elements into two
162   parts *)
163 AllSplits[l_List] := {#, Complement[l, #]} & /@ Subsets[l]
164
165 (* Will return all possible propagators for a BCFW cut *)
166 PropsForCut[{{Gluon[-] ...}, {a_, b_}] := {Gluon[1], Gluon[-1]}
167 PropsForCut[types_List, {a_, b_}] :=
168   PropsForCut[RotateLeft[types, b - 1][[; LLMod[types, a - b]]]
169 PropsForCut[{{left_... , Gluon[-] .. , right_...}] :=
170   PropsForCut[{{left, right}}]

```

```

168 PropsForCut[{left_---, Quark[p1_, h1_, f_], Quark[p2_, h2_, f_],
169   right_---}] := PropsForCut[{left, right}] /; p1 == -p2 && h1 == -h2
170 PropsForCut[{}] := {Gluon[1], Gluon[-1]}
171 PropsForCut[{q : Quark[-, -, -]}] := {q}
172 PropsForCut[{-, -}] := {}
173
174 (* Check if a pair of particles are valid for the BCFW shift particles i
      and j *)
175 InvalidShiftsQ[_List, {Gluon[1] | Quark[-, 1, -],
176   Gluon[-1] | Quark[-, -1, -]}, {_Integer, _Integer}] := True
177 InvalidShiftsQ[
178   types_List, {Quark[pa_, -, f_], Quark[pb_, -, f_]}, {a_Integer,
179   b_Integer}] :=
180   True /; pa == -pb && (LLMod[types, a + 1] == b ||
181     LLMod[types, b + 1] == a)
182 InvalidShiftsQ[types_List, {Quark[-, -1, -], Gluon[-1]} | {Gluon[1],
183   Quark[-, 1, -]}, {a_Integer, b_Integer}] := True /;
184   (LLMod[types, a + 1] == b || LLMod[types, b + 1] == a)
185 InvalidShiftsQ[_List, {-, -}, {_Integer, _Integer}] := False
186
187 (* Check for the factor introduced in the spinors by reversing the
      momenta *)
188 ReversingFactor[mom_List] := (DeclareSpinorMomentum[spinor, mom];
189   DeclareSpinorMomentum[mspinor, -mom];
190   NN[La[spinor]][[1, 1]]/NN[La[mspinor]][[1, 1]])
191 ReversingFactor[sm : {{-, -}, {-, -}}] :=
192   ReversingFactor[PfromSm2[sm]]
193
194 (* The extra factor needed to correct for how the momenta either side of
      the cut are defined *)
195 PropogatorFactor[Gluon[-, -]]:=1
196 PropogatorFactor[Quark[p_, -, -], mom_]:=ReversingFactor[-p*mom]
197
198
199 (* Calculate the momenta or particle types to use for the amplitude on
      one side of a cut *)
200 HelAmplBCFCombineSide[{NonOrd_List, ord_List}, {a_, b_},
201   splitNonOrd_List, prop_] := {Part}[NonOrd, splitNonOrd],
202   Append[LL[ord, LLRange[ord, a, b - 1]], prop]}
203
204 (* Evaluate a BCFW term *)
205 HelAmplNBCFTermImpl[{pNonOrd_List, p_List}, {i_, j_}, {a_,

```

```

206 b_}, {leftNonOrd_, rightNonOrd_}, prop_,
207 lefttypes : {_List, _List}, righttypes : {_List, _List}] :=
208 Module[{z, ii, jj, shiftedp, q, mq, qp, A},
209 z = z /.
210 Flatten[ExpandSToSpinors[
211 SpOpen[Solve[
212 0 == ShiftBA[LL[p, i], LL[p, j], z][
213 s @@ Join[Part[pNonOrd, leftNonOrd],
214 LL[p, LLRange[p, a, b - 1]]], z]]] // NN;
215 WithSpinors[shiftedp = ReplacePart[p, {i -> ii, j -> jj}];
216 qp = PfromSm2[(Sum[
217 Sm2[LL[shiftedp, qs]], {qs, LLRange[p, a, b - 1]} +
218 Sum[Sm2[pp], {pp, Part[pNonOrd, leftNonOrd]}])] // NN];
219 WithSpinors[
220 A = HelAmplN[
221 HelAmplBCFCCombineSide[{pNonOrd, shiftedp}, {a, b}, leftNonOrd,
222 mq], lefttypes]*
223 HelAmplN[
224 HelAmplBCFCCombineSide[{pNonOrd, shiftedp}, {b, a}, rightNonOrd,
225 q], righttypes];
226 A*I/NN[
227 s @@ Join[Part[pNonOrd, leftNonOrd],
228 LL[p, LLRange[p, a, b - 1]]]*PropogatorFactor[prop, qp], {q,
229 qp}, {mq, -qp}], {ii, {La[p[[i]]] // NN,
230 Lat[p[[i]]] - z*Lat[p[[j]]] //
231 NN}}, {jj, {La[p[[j]]] + z*La[p[[i]]] // NN, Lat[p[[j]]] // NN}}]]
232
233 (* Declare some terms that are known to vanish based on the particle
types on each side *)
234 HelAmplNBCFTerm[{pNonOrd_List, p_List}, {i_, j_}, {a_,
235 b_}, {leftNonOrd_, rightNonOrd_}, prop_,
236 lefttypes : {_List, _List},
237 righttypes : {{}, {Gluon[h_] | Quark[-, h_, -],
238 Gluon[h_] | Quark[-, h_, -],
239 Gluon[h_] | Quark[-, h_, -]} | {Gluon[1] | Quark[-, 1, -],
240 Gluon[1] | Quark[-, 1, -],
241 Gluon[-1] | Quark[-, -1, -]} | {Gluon[1] | Quark[-, 1, -],
242 Gluon[-1] | Quark[-, -1, -],
243 Gluon[1] | Quark[-, 1, -]} | {Gluon[-1] | Quark[-, -1, -],
244 Gluon[1] | Quark[-, 1, -], Gluon[1] | Quark[-, 1, -]}] := 0
245 HelAmplNBCFTerm[{pNonOrd_List, p_List}, {i_, j_}, {a_,
246 b_}, {leftNonOrd_, rightNonOrd_}, prop_,

```

```

247 lefttypes : {{}, {Gluon[h-] | Quark[-, h-, -],
248   Gluon[h-] | Quark[-, h-, -],
249   Gluon[h-] | Quark[-, h-, -]} | {Gluon[1] | Quark[-, 1, -],
250   Gluon[-1] | Quark[-, -1, -],
251   Gluon[-1] | Quark[-, -1, -]} | {Gluon[-1] | Quark[-, -1, -],
252   Gluon[1] | Quark[-, 1, -],
253   Gluon[-1] | Quark[-, -1, -]} | {Gluon[-1] | Quark[-, -1, -],
254   Gluon[-1] | Quark[-, -1, -], Gluon[1] | Quark[-, 1, -]}},
255 righttypes : {_List, _List}] := 0
256
257 (* Will swap the sides to put the term in a canonical order. The rule is
   if there is a single non-colour ordered particle it goes in the
   left half, otherwise the left half will be the smaller half in terms
   of number of particles. If neither of these rules choose an order
   the current order is used *)
258 HelAmplNBCFTerm[{pNonOrd_List, p_List}, {i-, j-}, {a-, b-},
259 nonordsplit : {{}, {}} | {{--}, {--}}, prop-,
260 lefttypes : {_List, _List}, righttypes : {_List, _List}] :=
261 HelAmplNBCFTermImpl[{pNonOrd, p}, {i, j}, {b, a},
262   Reverse[nonordsplit], ReverseParticle[prop], righttypes,
263   lefttypes] /; Mod[b - a, Length[p]] > Length[p]/2
264 HelAmplNBCFTerm[{pNonOrd_List, p_List}, {i-, j-}, {a-,
265   b-}, {{leftnonordsplit--}, {}}, prop-, lefttypes : {_List, _List},
266   righttypes : {_List, _List}] :=
267 HelAmplNBCFTermImpl[{pNonOrd, p}, {i, j}, {b,
268   a}, {{}, {leftnonordsplit}}, ReverseParticle[prop], righttypes,
269   lefttypes]
270 HelAmplNBCFTerm[{pNonOrd_List, p_List}, {i-, j-}, {a-,
271   b-}, {leftNonOrd-, rightNonOrd-}, prop-,
272   lefttypes : {_List, _List}, righttypes : {_List, _List}] :=
273 HelAmplNBCFTermImpl[{pNonOrd, p}, {i, j}, {a, b}, {leftNonOrd,
274   rightNonOrd}, prop, lefttypes, righttypes]
275
276 (* Declare some BCFW terms that are known to vanish *)
277 HelAmplNBCFTerm[{{}, p_List}, {{}, _List}, {-, -}, {a-,
278   b-}, {{}, {}}, -] :=
279 0 /; a == LMod[p, b + 1] || LMod[p, a + 1] == b
280 HelAmplNBCFTerm[[_List, p_List], [_List, _List], {-, -}, {a-,
281   b-}, {{}, {--}}, -] := 0 /; LMod[p, a + 1] == b
282 HelAmplNBCFTerm[[_List, p_List], [_List, _List], {-, -}, {a-,
283   b-}, {{--}, {}}, -] := 0 /; a == LMod[p, b + 1]
284

```

```

285 (* Calculate the types of particle on each side of the cut *)
286 HelAmplNBCFTerm[p : {-List, -List},
287   types : {-List, -List}, {i-, j-}, {a-, b-}, {leftNonOrd-,
288     rightNonOrd-}, prop-] :=
289   HelAmplNBCFTerm[p, {i, j}, {a, b}, {leftNonOrd, rightNonOrd}, prop,
290     HelAmplBCFCombineSide[types, {a, b}, leftNonOrd, prop],
291     HelAmplBCFCombineSide[types, {b, a}, rightNonOrd,
292       ReverseParticle[prop]]]
293
294
295 (* Evaluate an amplitude for a given choice of shift particles *)
296 HelAmplNBCF[{pNonOrd_List, p_List}, {typesNonOrd_List,
297   types_List}, {i-, j-}] := Sum[
298   HelAmplNBCFTerm[{pNonOrd, p}, {typesNonOrd, types}, {i, j}, {a, b},
299     splitNonOrd, prop], {a, LLRange[p, i + 1, j]}, {b,
300     LLRange[p, j + 1, i]}, {splitNonOrd,
301     AllSplits[Range[Length[pNonOrd]]]}, {prop,
302     PropsForCut[types, {a, b}]}]
303
304 (* Try different shifts until one evaluates successfully *)
305 HelAmplNBCFSearch[{pNonOrd_List, p_List}, {typesNonOrd_List,
306   types_List}] :=
307   Catch[Do[If[!
308     InvalidShiftsQ[types, {types[[i]], types[[j]]}, {i, j}],
309     Module[{val =
310       HelAmplNBCF[{pNonOrd, p}, {typesNonOrd, types}, {i, j}],
311       If[! MatchQ[
312         val, (- : 1)*Infinity | ComplexInfinity | Indeterminate],
313         Throw[val]]], {i, Length[p]}, {j,
314       LLRange[p, i + 1, i - 1]}; Indeterminate]
315
316 (* Evaluate the amplitude using BCFW. Will only be used if none of the
317   rules for specific types of amplitude match *)
317 HelAmplN[{pNonOrd_List, p_List}, {typesNonOrd_List, types_List}] :=
318   HelAmplNBCFSearch[{pNonOrd, p}, {typesNonOrd, types}] /; Length[pNonOrd
319     ]+Length[p] > 3

```

Listing B.3: Loop-Cuts.txt

```

1 (* A safe list part which only evaluates when the value to index is a
2   list *)
2 SPart[i_-][l_List] := Part[l, i]
3 \[GothicCapitalP] = SPart;

```

```

4
5 (* Declare momenta sets and processes, including which quarks are
   travelling left *)
6 Process[-] := {{}, {}}
7 LeftQuarks[-] := {}
8
9 DeclareProcess[p : {-List, -List}, leftquarks_ : {}] :=
10 Module[{process}, Process[process] = p; LeftQuarks[process] =
   leftquarks;
11 process]
12
13 MomConf[-] := {{}, {}}
14
15 DeclareMomConf[p : {-List, -List}] :=
16 Module[{momconf}, MomConf[momconf] = p; momconf]
17
18 (* Convert names to objects that can be added *)
19 ToSp[l_List] := ToSp /@ l
20 ToSp[i_Integer] := Sp[i]
21 ToSp[Sp[i_]] := Sp[i]
22 ToSp[s_Symbol /; LVectorQ[s]] := s
23
24 (* Declare the reverse spinors for an already defined momenta *)
25 DeclareReverseSpinor[spinors_List, reverse_List] :=
26 MapThread[DeclareReverseSpinor, {spinors, reverse}]
27 DeclareReverseSpinor[sp : (-Integer | -?SpinorQ),
28 r : (-Integer | -Symbol)] := (DeclareSpinorMomentum[r, I*La[sp] // NN,
29 I*Lat[sp] // NN]; {sp, r})
30
31 (* The factor introduced by reversing a particle *)
32 LoopReversingFactor[Quark[p_, ---]] := p I
33 LoopReversingFactor[_Gluon] := -1
34
35 (* Calculate the amplitude for a corner given the cut particles and the
   loop momenta defined as pointing around the loop *)
36 CalculateLoopCornerAmplitude[{pNonOrd_List,
37 p_List}, {typesNonOrd_List, {g1 : (Gluon[-] ...), Quark[q1_, h1_, f_
   ],
38 mid_..., Quark[q2_, h2_, f_], g2 : (Gluon[-] ...)}}, {Quark[q1_, h1_,
   f_],
39 Quark[q1_, h1_, f_]}, {l1_, l2_}] :=
40 CalculateLoopCornerAmplitude[{pNonOrd,

```



```

41 p}, {typesNonOrd, {g1, Quark[q1, h1, f], mid, Quark[q2, h2, -f - 1],
42   g2}}, {Quark[q1, h1, f], Quark[q1, h1, -f - 1]}, {l1, l2}]
43 CalculateLoopCornerAmplitude[{pNonOrd_List, p_List}, {typesNonOrd_List,
44   types_List}, {t1_, t2_}, {l1_, l2_}] :=
45 WithSpinors[
46   Module[{tmp}, DeclareReverseSpinor[l11, l11r];
47     tmp = HelAmplN[{pNonOrd, Join[p, {l12, l11r}]}, {typesNonOrd,
48       Join[types, {t2, ReverseParticle[t1]}]}]*LoopReversingFactor[t1];
49     FullyUndeclareSpinor[l11r]; tmp], {l11, l1}, {l12, l2}]
50 CalculateLoopCornerAmplitude[{pNonOrd_List, p_List}, {typesNonOrd_List,
51   types_List}, {t1_,
52   t2_}, {{{Indeterminate | {(Indeterminate | {Indeterminate ...}) ...}
53     ...}},
54   l2_}] := Indeterminate
55 CalculateLoopCornerAmplitude[{pNonOrd_List, p_List}, {typesNonOrd_List,
56   types_List}, {t1_,
57   t2_}, {l1_, {(Indeterminate | {(Indeterminate | {Indeterminate ...})
58     ...}) \
59   ...}}] := Indeterminate
60 (* Split the external particles into the different corners *)
61 SplitCorners[{Unord_List, Ord_List}, {splits_List, splitUnord_List}] :=
62   Transpose[{Extract[Unord, Position[splitUnord, #]] & /@
63     Range[Length[splits]],
64     RotateLeft[Ord, #1 - 1][[;; #2 - #1]] & @@@
65     Partition[splits, 2, 1, 1, Length[Ord] + splits[[1]]]}]
66
67 (* Combine the external particles in each corner with the loop
68   propagators to give the particles in each corner *)
69 CombineCornerPropsTypes[corners : {{_List, _List} ...}, props_List] :=
70   MapThread[{#1[[1]],
71     Join[#1[[2]], #2]} &, {corners, {#1[[2]], ReverseParticle[#1[[1]]]} &
72     /@
73     Partition[props, 2, 1, 1]}]
74
75 (* Calculate at what index a new split must be inserted to maintain the
76   order of the cuts *)
77 InsertionIndex[split_, corner_, newsplit_] :=
78   0 /; corner == Length[split] &&
79   newsplit <= split[[1]] && ! newsplit == split[[corner]]
80 InsertionIndex[split_, corner_, newsplit_] := corner

```

78

```

79 (* Calculate the name for a new combined split *)
80 SplitName[{{}, nonord_List], {corner_, newsplit_Integer,
81   newNonord_List}] := {{newsplit},
82   PadLeft[{}, Length[newNonord], 1]}
83 SplitName[{splits_List, nonord_List}, {corner_Integer,
84   newsplit_Integer, newNonord_List}] :=
85 Module[{insert = InsertionIndex[splits, corner, newsplit]}, {Insert[
86   splits, newsplit, insert + 1],
87   MapThread[
88     Mod[#1 + If[(#1 == corner) && #2 > 0, 1, 0] +
89       If[#1 > insert, 1, 0] - 1, Length[splits] + 1] + 1 &, {nonord,
90       newNonord}]}]}
91
92 (* Replace new splits that are completely invalid with Sequence[] *)
93 RemoveInvalidSplits[
94   process_, {split_, splitNonord_}, {c_, i_, nonordshift_}] :=
95   Sequence[] /; split[[c]] == i && Count[nonordshift, 0] == 0
96 RemoveInvalidSplits[
97   process_, {split_, splitNonord_}, {c_, i_, nonordshift_}] :=
98   Sequence[] /; (Length[split] !=
99     c || (i <= split[[1]] && ! split[[c]] == i)) &&
100   split[[Mod[c, Length[split]] + 1]] == i &&
101   Count[nonordshift, 1] == 0
102 RemoveInvalidSplits[process_, {split_, splitNonord_},
103   newsplit_] := newsplit
104
105 (* Calculate the list of possible new splits given an existing split *)
106 SplitOptions[process_, {{}, splitNonord_List}] :=
107   Table[{None, i, Table[0, {Length[Process[process][[1]]]}]}, {i,
108     Length[Process[process][[2]]]}]
109 SplitOptions[process_, {splits_List, splitNonord_List}] :=
110   RemoveInvalidSplits[process, {splits, splitNonord}, #] & /@
111   Flatten[Table[
112     Module[{nonordindex = Position[splitNonord, c]},
113       Table[{c,
114         Mod[i + splits[[c]] - 1, Length[Process[process][[2]]]] + 1,
115         ReplacePart[PadRight[{}, Length[splitNonord], None],
116           MapThread[Rule, {nonordindex, non}]}], {i, 0,
117         If[c == Length[splits],
118           splits[[1]] - splits[[c]] + Length[Process[process][[2]]],
119           splits[[c + 1]] - splits[[c]]}], {non,

```

```

120     Tuples[[0, 1], Length[nonordindex]]], {c, 1,
121     Length[splits]], 2]
122
123 (* Calculate all possible splits into n corners *)
124 SplitOptions[process_, 0] := {{{}, {}}
125 SplitOptions[process_, n_Integer] :=
126 Union @@
127 Function[name,
128     SplitName[name, #] & /@ SplitOptions[process, name]] /@
129     SplitOptions[process, n - 1]
130
131 MatchQuarks[Quark[p_, -, f_], process_] :=
132 MemberQ[LeftQuarks[process], Quark[p, f]]
133
134 SplitPropOptions[process_, {splits : {}, _List},
135     loopprops : {}, {corner_, newsplit_Integer, _List}] :=
136 Module[{quarks = Cases[LeftQuarks[process], Quark[p_, -1] :> p],
137     cornertypes =
138     RotateLeft[Process[process][[2]], newsplit - 1] /.,
139     Gluon[_] -> Sequence[] /. {o1_... ,
140     Quark[p1_, h1_, f_]?(MatchQuarks[#, process] &),
141     Quark[p2_, h2_, f_], o2_...} :> {o1, o2} /;
142     p1 == -p2 && h1 == -h2},
143 If[Length[cornertypes] == 0,
144 If[Length[quarks] ==
145     1, {{Quark[quarks[[1]], 1, -1]}, {Quark[
146     quarks[[1]], -1, -1]}}, {{Gluon[1]}, {Gluon[-1]}},
147 If[And[Length[quarks] == 0,
148     MatchQ[cornertypes, {Quark[p1_, h1_, f_],
149     Quark[p2_, h2_, f_]?(MatchQuarks[#, process] &)} /;
150     p1 == -p2 && h1 == -h2]], {{cornertypes[[1]]}, {}]]]
151
152 (* Calculate the possible loop propagators given a split and its
153     propagators and the new cut being added *)
154 SplitPropOptions[process_, {splits : {_Integer}, nonord_List},
155     loopprops : {_List ...}, {corner_Integer, newsplit_Integer,
156     newNonord_List}] :=
157 Flatten[SplitPropOptions[
158     process, {splits, nonord}, #, {corner, newsplit, newNonord}] & /@
159     loopprops, 1]
160 SplitPropOptions[process_, {splits : {_Integer}, nonord_List},
161     loopprops_List, {corner_Integer, newsplit_Integer,

```

```

161   newNonord_List}} :=
162   Module[{index = InsetIndex[splits, corner, newsplit],
163     name = SplitName[{splits, nonord}, {corner, newsplit, newNonord}],
164     cornerparts, options},
165     cornerparts =
166     CombineCornerPropsTypes[SplitCorners[Process[process], name],
167       Insert[looppops, None, index + 1][[index + 1, 2]];
168     cornerparts =
169     cornerparts /.
170     Gluon[_] | ReverseParticle[None] -> Sequence[] //. {o1_---,
171       Quark[p1_, h1_, f_]?(MatchQuarks[#, process] &),
172       Quark[p2_, h2_, f_], o2_---} :> {o1, o2} /;
173     p1 == -p2 && h1 == -h2;
174     options =
175     If[Length[cornerparts] == 0, {Gluon[1], Gluon[-1]},
176     If[Length[cornerparts] == 1, cornerparts, {}]];
177     Insert[looppops, #, index + 1] & /@ options]
178
179 (* Calculate all possible propagators for a given split *)
180 SplitPropOptions[process_, {splits_List, nonord_List}] :=
181   Module[{name = {{}, {}}, options = {}},
182     Do[options =
183       SplitPropOptions[process, name,
184         options, {i - 1, splits[[i]], nonord - i + 1}];
185     name = SplitName[name, {i - 1, splits[[i]], nonord - i + 1}], {i,
186       Length[splits]}; options]
187
188 (* Calculate the helicity of a corner if it has one, or return None *)
189 CornerHelicity[nonels_List, els_List] :=
190   CornerHelicity[Sort[nonels],
191     Sort[els]] /; ! (OrderedQ[els] && OrderedQ[nonels])
192
193 CornerHelicity[{}, {Gluon[h_], Quark[-, -, -], Quark[-, -, -]}] := h
194 CornerHelicity[{}, {Gluon[-1], Gluon[-1], Gluon[1]}] := -1
195 CornerHelicity[{}, {Gluon[-1], Gluon[1], Gluon[1]}] := 1
196 CornerHelicity[_List, _List] := None
197
198 (* Remove from the list of propagators all combinations that are known
199   to vanish for any reason *)
200 RemoveIgnorableOptions[process_, {splits : {_Integer}, nonord_List},
201   looppops : {_List ...}] :=
202   RemoveIgnorableOptions[process, {splits, nonord}, #] & /@ looppops

```

```

202 RemoveIgnorableOptions[process_, {splits : {_-}, nonord_List],
203   loopprops_List] := Sequence[]
204 RemoveIgnorableOptions[process_, {splits : {i-, j-}, {2 ...}},
205   loopprops_List] :=
206   Sequence[] /; Mod[i, Length[Process[process][[2]]]] + 1 == j
207 RemoveIgnorableOptions[process_, {splits : {i-, j-}, {1 ...}},
208   loopprops_List] :=
209   Sequence[] /; i == Mod[j, Length[Process[process][[2]]]] + 1
210 RemoveIgnorableOptions[process_, {splits : {_Integer}, nonord_List},
211   loopprops_List] :=
212   RemoveIgnorableOptions[process, {splits, nonord}, loopprops,
213     Map[Sort,
214       CombineCornerPropsTypes[
215         SplitCorners[Process[process], {splits, nonord}], loopprops], 2]]
216 RemoveIgnorableOptions[process_, {splits : {_Integer}, nonord_List},
217   loopprops_List, {---, {{}, {Gluon[h_] ...}}, ---}] := Sequence[]
218 RemoveIgnorableOptions[process_, {splits : {_Integer}, nonord_List},
219   loopprops_List, {---, {{}, {Gluon[h1-], Gluon[h2-], Gluon[h2-],
220     Gluon[h2-] ..} | {Gluon[h1-], Gluon[h1-], Gluon[h1-] ..,
221     Gluon[h2-]}}, ---}] := Sequence[]
222 RemoveIgnorableOptions[process_, {splits : {_Integer}, nonord_List},
223   loopprops_List, {---, {{}, {Gluon[h-],
224     Gluon[h-] .., -Quark, -Quark}}, ---}] := Sequence[]
225 RemoveIgnorableOptions[process_, {splits : {_Integer}, nonord_List},
226   loopprops_List, {---, {{Phi[p-]}, {Gluon[p-],
227     Gluon[h-]} | {Gluon[h-], Gluon[p-]}}, ---}] := Sequence[]
228 RemoveIgnorableOptions[process_, {splits : {_Integer}, nonord_List},
229   loopprops_List, {---, {{Phi[
230     p-]}, {Gluon[p-] ..., -Quark, -Quark}}, ---}] := Sequence[]
231 RemoveIgnorableOptions[process_, {splits : {_Integer}, nonord_List},
232   loopprops_List, {---, {{}, {Gluon[h1-], Gluon[h : (h1- | h2-)],
233     Gluon[h2-]} | {Gluon[h-], -Quark, -Quark}}, {{}, {Gluon[h1-],
234     Gluon[h : (h1- | h2-)],
235     Gluon[h2-]} | {Gluon[h-], -Quark, -Quark}}, ---}] := Sequence[]
236 RemoveIgnorableOptions[process_, {splits : {_Integer}, nonord_List},
237   loopprops_List, {{{}, {Gluon[h1-], Gluon[h : (h1- | h2-)],
238     Gluon[h2-]} | {Gluon[h-], -Quark, -Quark}}, ---, {{}, {Gluon[
239     h1-], Gluon[h : (h1- | h2-)],
240     Gluon[h2-]} | {Gluon[h-], -Quark, -Quark}}}] := Sequence[]
241 RemoveIgnorableOptions[process_, {splits : {_Integer}, nonord_List},
242   loopprops_List, {pre---, {{}, {Gluon[h1-], Gluon[h1-],
243     Gluon[h2-]}}, -, {{}, {Gluon[h1-], Gluon[h2-], Gluon[h2-]}},

```

```

244   post_...}] :=
245   Sequence [] /; Length{pre, post} == 1 (* total length==4 *)
246   RemoveIgnorableOptions[process_, {splits : {--Integer}, nonord_List},
247   loopprops_List, {pre_..., {{}}, {Gluon[h2_], Gluon[h1_],
248     Gluon[h1_]}}, -, {{}}, {Gluon[h2_], Gluon[h2_], Gluon[h1_]}},
249   post_...}] :=
250   Sequence [] /; Length{pre, post} == 1 (* total length==4 *)
251
252   RemoveIgnorableOptions[_, -, loopprops_List, _] := loopprops
253
254   SplitValidPropOptions[process_, split_] := RemoveIgnorableOptions[
255     process, split, SplitPropOptions[process, split]]
256
257   (* Calculate the momentum solution for a box *)
258   CalcBoxLoopMom[{{{}}, {k1_}}, K2 : {_List, _List}, K3 : {_List, _List},
259     K4 : {_List, _List}}, 1] :=
260   Module{k2 = Plus @@ ToSp[K2 // Flatten],
261     k3 = Plus @@ ToSp[K3 // Flatten],
262     k4 = Plus @@ ToSp[K4 // Flatten], {{La[k1],
263     CLa[k1].CSm2[k2].Sm2[k3].CSm2[k4]/Spaa[k1, k2, k4, k1]}}, {La[k1],
264     CLa[k1].CSm2[k4].Sm2[k3].CSm2[k2]/
265     Spaa[k1, k2, k4, k1]}}, {CSm2[k3].Sm2[k4].La[k1]/
266     Spaa[k1, k2, k4, k1],
267     CLa[k1].CSm2[k2]}}, {CSm2[k3].Sm2[k2].La[k1]/Spaa[k1, k2, k4, k1],
268     CLa[k1].CSm2[k4]} // NN]
269   CalcBoxLoopMom[{{{}}, {k1_}}, K2 : {_List, _List}, K3 : {_List, _List},
270     K4 : {_List, _List}}, -1] :=
271   Module{k2 = Plus @@ ToSp[K2 // Flatten],
272     k3 = Plus @@ ToSp[K3 // Flatten],
273     k4 = Plus @@
274     ToSp[K4 // Flatten], {{CSm2[k4].Sm2[k3].CSm2[k2].CLat[k1]/
275     Spbb[k1, k4, k2, k1],
276     Lat[k1]}}, {CSm2[k2].Sm2[k3].CSm2[k4].CLat[k1]/
277     Spbb[k1, k4, k2, k1], Lat[k1]}}, {CSm2[k2].CLat[k1],
278     Lat[k1].Sm2[k4].CSm2[k3]/
279     Spbb[k1, k4, k2, k1]}}, {CSm2[k4].CLat[k1],
280     Lat[k1].Sm2[k2].CSm2[k3]/Spbb[k1, k4, k2, k1]} // NN]
281
282   CalcBoxLoopMom[momconf_, {split_List, nonord_List},
283     masslessindex_Integer, h_] :=
284   CalcBoxLoopMom[momconf, {split, nonord}, masslessindex, h] =
285   RotateRight[

```

```

286 CalcBoxLoopMom[
287   RotateLeft[SplitCorners[MomConf[momconf], {split, nonord}],
288     masslessindex - 1], h], masslessindex - 1]
289
290 (* Determine if a box has a massless corner and if so where it is *)
291 MasslessBoxCorner[
292   process_, {split : {pre---, i-, j-, ---}, nonord_List}] :=
293   Length[{pre}] + 1 /; j == i + 1 && FreeQ[nonord, Length[{pre}] + 1]
294 MasslessBoxCorner[process_, {split : {1, ---, i-}, nonord_List}] :=
295   4 /; Length[Process[process][[2]]] == i && FreeQ[nonord, 4]
296 MasslessBoxCorner[
297   process_, {split : {---Integer}, nonord : {---Integer}}] := None
298
299 (* Determine which momenta solution is valid for a given box *)
300 ValidBoxLoopSolution[{nonord_List, ord_List}] :=
301   ValidBoxLoopSolution[Sort[nonord], Sort[ord]]
302 ValidBoxLoopSolution[{}, {Gluon[h-], Quark[-, -, -],
303   Quark[-, -, -]}] := h
304 ValidBoxLoopSolution[{}, {Gluon[-1], Gluon[-1], Gluon[1]}] := -1
305 ValidBoxLoopSolution[{}, {Gluon[-1], Gluon[1], Gluon[1]}] := 1
306 ValidBoxLoopSolution[{}, {-, -, -}] := None
307
308 (* Calculate the contribution from a single box *)
309 (* masslessindex is an integer so there is a massless corner. If there
    are no massless corners, a version where masslessindex is None must
    be implemented *)
310 CalculateBoxContribution[process_,
311   momconf_, {split_List, nonord_List}, masslessindex_Integer, props_,
312   solutionhel_] :=
313   Module[{solution,
314     cornertypes = SplitCorners[Process[process], {split, nonord}],
315     cornernames = SplitCorners[MomConf[momconf], {split, nonord}],
316     If[solutionhel == None, 0,
317     solution =
318       CalcBoxLoopMom[momconf, {split, nonord}, masslessindex,
319       solutionhel];
320     I/2*Product[
321       CalculateLoopCornerAmplitude[cornernames[[i]],
322       cornertypes[[i]], {props[[i]],
323       props[[Mod[i, 4] + 1]]}, {solution[[i]],
324       solution[[Mod[i, 4] + 1]]}], {i, Length[split]}] /;
325   ValidBoxLoopSolution@(CombineCornerPropsTypes[

```

```

326     SplitCorners[Process[process], {split, nonord}], props][[
327     masslessindex]]) == solutionhel
328
329 CalculateBoxContribution[process_,
330   momconf_, {split_List, nonord_List}, masslessindex_Integer, props_,
331   solutionhel_] :=
332 0 /; ValidBoxLoopSolution@(CombineCornerPropsTypes[
333   SplitCorners[Process[process], {split, nonord}], props][[
334   masslessindex]]) != solutionhel
335
336 CalculateBoxContribution[process_,
337   momconf_, {split_List, nonord_List}, props_,
338   solutionhel_Integer] := (CalculateBoxContribution[process,
339   momconf, {split, nonord}, props, solutionhel] =
340   CalculateBoxContribution[process, momconf, {split, nonord},
341   MasslessBoxCorner[process, {split, nonord}], props, solutionhel])
342 CalculateBoxContribution[process_,
343   momconf_, {split_List, nonord_List}, props_,
344   None | PatternSequence[]] :=
345 CalculateBoxContribution[process, momconf, {split, nonord},
346   props, -1] +
347 CalculateBoxContribution[process, momconf, {split, nonord}, props,
348   1]
349
350 (* Calculate the equation for the triangle loop momentum as a function
351    of t, evaluating as much as possible, as early as possible *)
351 CalcTriLoopMomEqn[{kk1_List, _, kk3_List}, None, h_] :=
352 Module[{k1 = Plus @@ ToSp[kk1 // Flatten],
353   k3 = Plus @@ ToSp[kk3 // Flatten], K1, K3, S1, S3,
354   k1k3, \[CapitalDelta], \[Gamma], K3t, K1t}, K1 = Num4V[k1];
355 K3 = Num4V[k3]; S1 = MP2[k1] // NN; S3 = MP2[k3] // NN;
356 k1k3 = MP[k1, k3] // NN; \[CapitalDelta] = k1k3^2 - S1 S3; \[Gamma] =
357 k1k3 + h*Sqrt[\[CapitalDelta]]; K3t = K3 - S3/\[Gamma] K1;
358 K1t = K1 - S1/\[Gamma] K3;
359 Function[t,
360 Evaluate[
361   WithSpinors[{{t La[k1t] +
362     S1*(S3 + \[Gamma])/(4*\[CapitalDelta]) La[
363     k3t], -S3*(S1 + \[Gamma])/(4*\[CapitalDelta]*t) Lat[k1t] +
364     Lat[k3t]}, {t La[k1t] +
365     S1*S3*(S1 + \[Gamma])/(4*\[Gamma]*\[CapitalDelta]) La[
366     k3t], -\[Gamma]*(S3 + \[Gamma])/(4*\[CapitalDelta]*t) Lat[

```



```

367         k1t] + Lat[
368         k3t}], {t La[
369         k1t] + \[Gamma]*(S1 + \[Gamma])/(4*\[CapitalDelta]) La[
370         k3t], -S1*
371         S3*(S3 + \[Gamma])/(4*\[Gamma]*\[CapitalDelta]*t) Lat[k1t] +
372         Lat[k3t]} // NN, {k1t, K1t}, {k3t, K3t}]]]]
373
374 CalcTriLoopMomEqn[corners_List, masslessindex_Integer, 1] :=
375 Module[{k1 = Plus @@ ToSp[corners[[masslessindex]] // Flatten],
376         k3 = Plus @@
377         ToSp[corners[[Mod[masslessindex + 1, 3] + 1]] // Flatten], c,
378         K3t}, c = MP2[k3]/(2 MP[k1, k3]) // NN;
379 K3t = Num4V[k3] - c Num4V[k1];
380 Function[t,
381 Evaluate[
382 RotateRight[
383 WithSpinors[{{t La[k1], -c/t Lat[k1] +
384             Lat[k3t]}, {t La[k1], -(c + 1)/t Lat[k1] +
385             Lat[k3t]}, {t La[k1] + La[k3t], Lat[k3t]} // NN, {k3t,
386             K3t}], masslessindex - 1]]]]]
387 CalcTriLoopMomEqn[corners_List, masslessindex_Integer, -1] :=
388 Module[{k1 = Plus @@ ToSp[corners[[masslessindex]] // Flatten],
389         k3 = Plus @@
390         ToSp[corners[[Mod[masslessindex + 1, 3] + 1]] // Flatten], c,
391         K3t}, c = MP2[k3]/(2 MP[k1, k3]) // NN;
392 K3t = Num4V[k3] - c Num4V[k1];
393 Function[t,
394 Evaluate[
395 RotateRight[
396 WithSpinors[{{-c /t La[k1] + La[k3t],
397             t Lat[k1]}, {-(c + 1)/ t La[k1] + La[k3t],
398             t Lat[k1]}, {La[k3t], t*Lat[k1] + Lat[k3t]} // NN, {k3t,
399             K3t}], masslessindex - 1]]]]]
400
401 CalcTriLoopMomEqn[momconf_, {split_List, nonord_List}, masslessindex_,
402 h_] := Module[{corners =
403 SplitCorners[MomConf[momconf], {split, nonord}], params},
404 CalcTriLoopMomEqn[momconf, {split, nonord}, masslessindex, h] =
405 CalcTriLoopMomEqn[corners, masslessindex, h]]
406
407 CalcTriLoopMom[momconf_, {split_List, nonord_List}, masslessindex_,
408 h_, t_] :=

```

```

409 CalcTriLoopMomEqn[momconf, {split, nonord}, masslessindex, h][t]
410
411 (* Calculate the subtraction of a box from a triangle as a function of t
      , evaluating as much as possible, as early as possible *)
412 CalculateTriSubtractionEqnImpl[process_, momconf_, split_,
413   masslessindex_Integer, solution : (1 | -1), props_] :=
414   Function[t,
415     Evaluate[Module[{cornernames =
416       SplitCorners[MomConf[momconf], split], k1, K3, C},
417       k1 = Plus @@ ToSp[cornernames[[masslessindex]] // Flatten];
418       K3 = Plus @@
419       ToSp[cornernames[[Mod[masslessindex + 1, 3] + 1]] // Flatten];
420       C = MP2[K3]/2/MP[k1, K3] // NN;
421       WithSpinors[
422         Sum[Module[{boxname = SplitName[split, boxsplit],
423           insindex =
424             InsertionIndex[split[[1]], boxsplit[[1]], boxsplit[[2]]], k},
425           k =
426             Plus @@ ToSp[
427               Flatten[If[insindex >= masslessindex,
428                 SplitCorners[MomConf[momconf], boxname][[
429                   masslessindex ;; insindex]],
430                 RotateLeft[SplitCorners[MomConf[momconf], boxname],
431                   masslessindex][[
432                     1 ;; insindex - masslessindex + 4]]]]]; (1/
433               If[solution == 1,
434                 Spab[k1, k,
435                   k3T]*(t - (MP[k, k] + 2 C MP[k, k1])/Spab[k1, k, k3T]),
436                 Spab[k3T, k,
437                   k1]*(t - (MP[k, k] + 2 C MP[k, k1])/Spab[k3T, k, k1])] //
438               NN)*Sum[CalculateBoxContribution[process, momconf, boxname,
439                 boxprops,
440                 solution*(-1)^(masslessindex +
441                   If[insindex < masslessindex, 1, 0] -
442                   MasslessBoxCorner[process, boxname])], {boxprops,
443                 RemoveIgnorableOptions[process, boxname,
444                 SplitPropOptions[process, split, props,
445                 boxsplit]]}], {boxsplit,
446                 SplitOptions[process, split]}, {k3T,
447                 Num4V[K3] - C*Num4V[k1]}]]]]]
448
449 CalculateTriSubtractionEqnImpl[process_, momconf_, split_,

```

```

450 masslessindex : None, h_, props_] :=
451 Module[{cornernames = SplitCorners[MomConf[momconf], split], k1, k3,
452   K1, K3, S1, S3, k1k3, \[CapitalDelta]},
453 k1 = Plus @@ ToSp[cornernames[[1]] // Flatten];
454 k3 = Plus @@ ToSp[cornernames[[3]] // Flatten]; K1 = Num4V[k1];
455 K3 = Num4V[k3]; S1 = MP2[k1] // NN; S3 = MP2[k3] // NN;
456 k1k3 = MP[k1, k3] // NN; \[CapitalDelta] = k1k3^2 - S1 S3;
457 Function[t,
458 Evaluate[
459   Module[{\[Gamma] = k1k3 + h*Sqrt[\[CapitalDelta]]},
460     WithSpinors[
461       Sum[Module[{boxname = SplitName[split, boxesplit],
462         insindex =
463           InsertionIndex[split[[1]], boxesplit[[1]], boxesplit[[2]]], k,
464           f, b, c, j, boxmasslessindex},
465         k = Plus @@
466           ToSp[Flatten[
467             If[insindex >= 1,
468               SplitCorners[MomConf[momconf], boxname][[1 ;; insindex]],
469               RotateLeft[SplitCorners[MomConf[momconf], boxname], 1][[
470                 1 ;; insindex + 3]]]]; f = Spab[k1T, k, k3T] // NN;
471             b = (MP[k,
472               k] + (S3*(S1 + \[Gamma])*MP[k, k1T] -
473               S1*(S3 + \[Gamma])*
474               MP[k, k3T])/(2 \[CapitalDelta]))/(2 f) // NN;
475             c = -S1*S3*(S1 + \[Gamma])*
476               Spab[k3T, k, k1T]/((4 \[CapitalDelta])^2*f) // NN;
477             j = Sqrt[b^2 - c];
478             boxmasslessindex = MasslessBoxCorner[process, boxname];
479             Sum[CalculateBoxContribution[process, momconf, boxname,
480               boxprops, boxh]*
481               If[WithSpinors[(MP[ltm, lb] - MP[ltp, lb])/MP[ltp, ltm] //
482                 NN // Re, {lb,
483                 CalcBoxLoopMom[momconf, boxname, boxmasslessindex,
484                   boxh][[1]]}, {ltp, \[GothicCapitalP][1]@
485                   CalcTriLoopMom[momconf, split, masslessindex, h,
486                   b + j]}, {ltm, \[GothicCapitalP][1]@
487                   CalcTriLoopMom[momconf, split, masslessindex, h,
488                   b - j]}] <
489                 0, -(b - j)/(j*(t - (b - j))), (b +
490                   j)/(j*(t - (b + j))), {boxh, -1, 1}, {boxprops},
491               RemoveIgnorableOptions[process, boxname,

```

```

492         SplitPropOptions[process, split, props, boxsplit]]]/f/
493     2 ], {boxsplit, SplitOptions[process, split]}, {k1T,
494     K1 - S1/\[Gamma] K3}, {k3T, K3 - S3/\[Gamma] K1}}]]]]
495
496 CalculateTriSubtractionEqn[process_,
497     momconf_, {split_List, nonord_List}, masslessindex_, soln_,
498     props_] :=
499 CalculateTriSubtractionEqn[process, momconf, {split, nonord},
500     masslessindex, soln, props] =
501 CalculateTriSubtractionEqnImpl[process, momconf, {split, nonord},
502     masslessindex, soln, props]
503
504 CalculateTriSubtraction[args_, t_] :=
505 CalculateTriSubtractionEqn[args][
506     t] (* args is always process_, momconf_, split_, masslessindex_, soln_,
507         props_ *)
508
509 (* Check which momentum solutions are valid for a given triangle *)
510 ValidTriLoopSolution[{nonord_List, ord_List}] :=
511 ValidTriLoopSolution[Sort[nonord], Sort[ord]]
512 ValidTriLoopSolution[{}], {Gluon[h_], Quark[-, -, -],
513     Quark[-, -, -]}] := h
514 ValidTriLoopSolution[{}], {Gluon[-1], Gluon[-1], Gluon[1]}] := -1
515 ValidTriLoopSolution[{}], {Gluon[-1], Gluon[1], Gluon[1]}] := 1
516 ValidTriLoopSolution[{}], {-, -, -}] := None
517
518 (* Try and find a massless corner in a triangle *)
519 MasslessTriCorner[
520     process_, {split : {pre_---, i_, j_, ---}, nonord_List},
521     props_List] :=
522     Length[{pre}] + 1 /;
523     j == i + 1 && FreeQ[nonord, Length[{pre}] + 1] &&
524     1 == ValidTriLoopSolution@(CombineCornerPropsTypes[
525         SplitCorners[Process[process], {split, nonord}], props][[
526             Length[{pre}] + 1]])
527 MasslessTriCorner[process_, {split : {1, ---, i_}, nonord_List},
528     props_List] :=
529     3 /; Length[Process[process][[2]]] == i && FreeQ[nonord, 3] &&
530     1 == ValidTriLoopSolution@(CombineCornerPropsTypes[
531         SplitCorners[Process[process], {split, nonord}], props][[3]])
532 MasslessTriCorner[
533     process_, {split : {pre_---, i_, j_, ---}, nonord_List},

```

```

533 props_List] :=
534 Length[{pre}] + 1 /; j == i + 1 && FreeQ[nonord, Length[{pre}] + 1]
535 MasslessTriCorner[process_, {split : {1, ---, i_}, nonord_List},
536 props_List] :=
537 3 /; Length[Process[process][[2]]] == i && FreeQ[nonord, 3]
538 MasslessTriCorner[
539 process_, {split : {---Integer}, nonord : {---Integer}},
540 props_List] := None
541
542 (* Calculate a raw triangle contribution as a function of t *)
543 CalculateRawTriContributionImpl[cornernames_, cornertypes_, props_,
544 solution_] :=
545 1/2 Product[
546 CalculateLoopCornerAmplitude[\[GothicCapitalP][i]@
547 cornernames, \[GothicCapitalP][i]@
548 cornertypes, {\[GothicCapitalP][i]@
549 props, \[GothicCapitalP][Mod[i, 3] + 1]@
550 props}, {\[GothicCapitalP][i]@
551 solution, \[GothicCapitalP][Mod[i, 3] + 1]@solution}], {i, 3}]
552
553 CalculateRawTriContribution[process_,
554 momconf_, {split_List, nonord_List}, masslessindex_Integer, props_,
555 solutionh_, t_] :=
556 Module[{cornertypes =
557 SplitCorners[Process[process], {split, nonord}],
558 cornernames = SplitCorners[MomConf[momconf], {split, nonord}]},
559 CalculateRawTriContributionImpl[cornernames, cornertypes, props,
560 CalcTriLoopMom[momconf, {split, nonord}, masslessindex,
561 solutionh, t]] -
562 CalculateTriSubtraction[process, momconf, {split, nonord},
563 masslessindex, solutionh, props, t] /;
564 solutionh ==
565 ValidTriLoopSolution@(CombineCornerPropsTypes[cornertypes,
566 props][[masslessindex]])]
567
568 CalculateRawTriContribution[process_,
569 momconf_, {split_List, nonord_List}, masslessindex_Integer, props_,
570 solutionh_, t_] :=
571 Module[{cornertypes =
572 SplitCorners[Process[process], {split, nonord}]},
573 0 /; solutionh !=
574 ValidTriLoopSolution@(CombineCornerPropsTypes[cornertypes,

```

```

575         props][[masslessindex]])]
576
577 CalculateRawTriContribution[process_,
578   momconf_, {split_List, nonord_List}, masslessindex : None, props_,
579   solutionh_, t_] :=
580 Module{{cornertypes =
581   SplitCorners[Process[process], {split, nonord}],
582   cornernames = SplitCorners[MomConf[momconf], {split, nonord}]},
583   CalculateRawTriContributionImpl[cornernames, cornertypes, props,
584   CalcTriLoopMom[momconf, {split, nonord}, masslessindex, solutionh,
585   t]] - CalculateTriSubtraction[process, momconf, {split, nonord},
586   masslessindex, solutionh, props, t]}
587
588 CalculateRawTriContribution[process_,
589   momconf_, {split_List, nonord_List}, props_, h_, t_] :=
590 CalculateRawTriContribution[process, momconf, {split, nonord},
591   MasslessTriCorner[process, {split, nonord}, props], props, h, t]
592
593 (* Set up the constants used for the extraction of the different
594 coefficients *)
594 NNCache[t0, 3 Sqrt[2] - Pi, 2]
595 pt0[---] := t0
596 p = 9; (* Number of points to use *)
597 pp[---] := p
598 mp = 4; (* Maximum power - high enough for a Higgs boson to work *)
599 pmp[---] := mp
600
601 (* Calculate the equation for the coefficient of any power in a triangle
602 *)
602 CalculateTriContributionEquation[process_,
603   momconf_, {split_List, nonord_List}, props_, h_] :=
604 CalculateTriContributionEquation[process, momconf, {split, nonord},
605   props, h] =
606 Module{{p = pp[process], t0 = pt0[process]},
607   Function[n,
608     Evaluate[
609       Sum[t0-n*
610         CalculateRawTriContribution[process,
611           momconf, {split, nonord}, props, h,
612           t0*Exp[2 Pi I j/(2 p + 1)]]*
613         Exp[-2 Pi I j n/(2 p + 1)], {j, -p, p}]/(2 p + 1) // Factor]]]
614

```

```

615 CalculateTriContribution[process_,
616   momconf_, {split_List, nonord_List}, props_, h_Integer, n_: 0] :=
617 CalculateTriContributionEquation[process, momconf, {split, nonord},
618   props, h][n]
619 CalculateTriContribution[process_,
620   momconf_, {split_List, nonord_List}, props_, h : None, n_: 0] :=
621 CalculateTriContribution[process, momconf, {split, nonord}, props, 1,
622   n] + CalculateTriContribution[process, momconf, {split, nonord},
623   props, -1, n]
624 CalculateTriContribution[process_,
625   momconf_, {split_List, nonord_List}, props_] :=
626 CalculateTriContribution[process, momconf, {split, nonord}, props,
627   None, 0]
628
629 (* The arbitrary vector used in the bubble momentum parametrisation *)
630 DeclareSpinorMomentum[\[Chi], {1 + 2 I, 1 - 2 I, 1 + 4 I,
631   NN[Sqrt[15], 100]}]
632
633 (* Calculate the bubble loop momentum as a function of t and y,
634   evaluating as much as possible, as early as possible *)
635 CalcBubbleLoopMomEqn[{kk1_List, _}] :=
636 Module[{k1 = Plus @@ ToSp[kk1 // Flatten], K1t, S1o\[Chi]},
637   S1o\[Chi] = MP2[k1]/2/MP[k1, \[Chi]] // NN;
638   K1t = Num4V[k1] - S1o\[Chi] Num4V[\[Chi]];
639   Function[{t, y},
640     Evaluate[
641       WithSpinors[{{t La[k1t] + (1 - y) S1o\[Chi] La[\[Chi]],
642         y/t Lat[k1t] + Lat[\[Chi]]}, {-t La[k1t] +
643         y S1o\[Chi] La[\[Chi]], (1 - y)/t Lat[k1t] - Lat[\[Chi]]}] //
644       NN, {k1t, K1t}]]]]
645 CalcBubbleLoopMomEqn[momconf_, {split_List, nonord_List}] :=
646 Module[{corners = SplitCorners[MomConf[momconf], {split, nonord}],
647   params},
648   CalcBubbleLoopMomEqn[momconf, {split, nonord}] =
649   CalcBubbleLoopMomEqn[corners]]
650 CalcBubbleLoopMom[momconf_, {split_List, nonord_List}, t_, y_] :=
651 CalcBubbleLoopMomEqn[momconf, {split, nonord}][t, y // NN]
652
653 (* Calculate the raw and unsubtracted bubble contribution, evaluating as
654   much as possible, as early as possible *)
655 CalculateRawBubbleContributionImpl[cornernames_, cornertypes_, props_,
656   solution_] := -I Product[

```

```

655 CalculateLoopCornerAmplitude[cornernames[[i]],
656   cornertypes[[i]], {props[[i]],
657   props[[Mod[i, 2] + 1]]}, {solution[[i]],
658   solution[[Mod[i, 2] + 1]]}], {i, 2}
659
660 (* Calculate the t to use in a triangle from the bubble's t and y *)
661 CalculateTriTFromBubble[momconf_, k1t_, S1o\[Chi]_, triname_,
662   tricorner_Integer, h_Integer, -1, t_, y_] := 0
663 CalculateTriTFromBubble[momconf_, k1t_, S1o\[Chi]_, triname_,
664   tricorner_Integer, 1, power_, t_, y_] :=
665 Module[{corners = SplitCorners[MomConf[momconf], triname], tk1, tk3},
666   tk1 = Plus @@ ToSp[Flatten[corners[[tricorner]]]];
667   tk3 = Plus @@ ToSp[Flatten[corners[[Mod[tricorner + 1, 3] + 1]]]];
668   WithSpinors[(t Spaa[k1t, tK3t] -
669     S1o\[Chi] (1 - y) Spaa[tK3t, \[Chi]]) (t Spbb\[Chi], tk1) -
670     y Spbb[tk1, k1t])/(t s[tk1, tK3t]) // NN, {tK3t,
671     Num4V[tk3] - NN[MP2[tk3]/2 /MP[tk1, tk3]] Num4V[tk1]}]]
672 CalculateTriTFromBubble[momconf_, k1t_, S1o\[Chi]_, triname_,
673   tricorner_Integer, -1, power_, t_, y_] :=
674 Module[{corners = SplitCorners[MomConf[momconf], triname], tk1, tk3},
675   tk1 = Plus @@ ToSp[Flatten[corners[[tricorner]]]];
676   tk3 = Plus @@ ToSp[Flatten[corners[[Mod[tricorner + 1, 3] + 1]]]];
677   WithSpinors[(t Spaa[k1t, tk1] -
678     S1o\[Chi] (1 - y) Spaa[tk1, \[Chi]]) (t Spbb\[Chi], tK3t) -
679     y*Spbb[tK3t, k1t])/(t s[tk1, tK3t]) // NN, {tK3t,
680     Num4V[tk3] - NN[MP2[tk3]/2 /MP[tk1, tk3]] Num4V[tk1]}]]
681 CalculateTriTFromBubble[momconf_, k1t_, S1o\[Chi]_, triname_, None,
682   h_, 1, t_, y_] :=
683 Module[{corners = SplitCorners[MomConf[momconf], triname], tk1, tk3,
684   tK1, tK3, S1, S3, \[Gamma]},
685   tk1 = Plus @@ ToSp[Flatten[corners[[1]]]];
686   tk3 = Plus @@ ToSp[Flatten[corners[[3]]]]; tK1 = Num4V[tk1];
687   tK3 = Num4V[tk3]; S1 = NN[MP2[tk1]];
688   S3 = NN[MP2[tk3]]; \[Gamma] =
689   NN[MP[tk1, tk3]] + h Sqrt[NN[MP[tk1, tk3]]^2 - S1 S3];
690   WithSpinors[(t Spaa[k1t, tK3t] -
691     S1o\[Chi] (1 - y) Spaa[tK3t, \[Chi]]) (t Spbb\[Chi], tK1t) -
692     y Spbb[tK1t, k1t])/(t s[tK1t, tK3t]) // NN, {tK1t,
693     tK1 - S1 tK3/\[Gamma]}, {tK3t, tK3 - S3 tK1/\[Gamma]}]]
694 CalculateTriTFromBubble[momconf_, k1t_, S1o\[Chi]_, triname_, None,
695   h_, -1, t_, y_] :=
696 Module[{corners = SplitCorners[MomConf[momconf], triname], tk1, tk3,

```



```

697   tK1, tK3, S1, S3, \[CapitalDelta], \[Gamma]},
698   tk1 = Plus @@ ToSp[Flatten[corners[[1]]]];
699   tk3 = Plus @@ ToSp[Flatten[corners[[3]]]]; tK1 = Num4V[tk1];
700   tK3 = Num4V[tk3]; S1 = NN[MP2[tk1]];
701   S3 = NN[MP2[tk3]]; \[CapitalDelta] =
702   MP[tK1, tK3]^2 - S1 S3 // NN; \[Gamma] =
703   NN[MP[tk1, tk3]] + h Sqrt[\[CapitalDelta]];
704   WithSpinors[-16 \[CapitalDelta]^2 (t Spaa[k1t, tK1t] -
705     S1o\[Chi] (1 - y) Spaa[tK1t, \[Chi]]) (t Spbb[\[Chi], tK3t] -
706     y Spbb[tK3t, k1t])/(t S1 S3 (S1 + \[Gamma]) (S3 + \[Gamma]) s[
707     tK1t, tK3t]) // NN, {tK1t, tK1 - S1 tK3/\[Gamma]}, {tK3t,
708     tK3 - S3 tK1/\[Gamma]}]]
709
710 (* Calculate the bubble subtraction contributions, evaluating as much as
711    possible, as early as possible *)
712 CalculateBubbleSubtractionEqnImpl[process_, momconf_, split_,
713   props_] :=
714 Module[{cornernames = SplitCorners[MomConf[momconf], split], K1,
715   S1o\[Chi], ypoles, res, tts, ttsy, tripropss, triconts},
716 K1 = Plus @@ ToSp[cornernames[[1]] // Flatten];
717 S1o\[Chi] = MP2[K1]/2/MP[K1, \[Chi]] // NN;
718 WithSpinors[
719 Do[Module[{triname = SplitName[split, tri],
720   insindex = InsertionIndex[split[[1]], tri[[1]], tri[[2]]], K2,
721   S2, \[Chi]K2k1t, a, at, bt, c, sqrtterm},
722 K2 =
723 Plus @@ ToSp[
724 Flatten[If[insindex == 0,
725   RotateLeft[SplitCorners[MomConf[momconf], triname], 1][[
726   1 ;; insindex + 3 - 1]],
727   SplitCorners[MomConf[momconf], triname][[1 ;; insindex]]]];
728 S2 = MP2[K2] // NN;
729 c = 1/(S1o\[Chi]*Spab[\[Chi], K2, bk1t]) // NN;
730 bt = c*MP[K1, K2] - (2*MP[K2, \[Chi]])/Spab[\[Chi], K2, bk1t] //
731 NN; a = c*MP[K2, K2] - 2 MP[K2, \[Chi]]/Spab[\[Chi], K2, bk1t] //
732 NN; at = -c*Spab[bk1t, K2, \[Chi]] // NN;
733 ypoles[tri, t_, h_] :=
734 Evaluate[(1/2 + t bt) + h*Sqrt[(1/2 + t bt)^2 - t (a + t at)]];
735 res[tri, t_] :=
736 Evaluate[c t/Sqrt[(1/2 + t bt)^2 - t (a + t at)]];
737 ttsy[tri, corner_, h_, trih_,
738   power_] := (ttsy[tri, corner, h, trih, power] =

```

```

738     Function[{t, y},
739     Evaluate[
740         CalculateTriTFromBubble[momconf, bk1t, S1o\[Chi], triname,
741             corner, trih, power, t, y]]];
742     tts[tri, corner_, h_, trih_,
743         power_] := (tts[tri, corner, h, trih, power] =
744     Function[{t},
745     Evaluate[
746         ttsy[tri, corner, h, trih, power][t, ypoles[tri, t, h]]]);
747     triconts[tri, tricorner_, h_, trih_, triprops_, ttp_, ttm_] :=
748
749         CalculateTriContribution[process, momconf, triname,
750             triprops, trih, 0] +
751     Sum[ttp^i CalculateTriContribution[process, momconf, triname,
752             triprops, trih, i] +
753         ttm^i CalculateTriContribution[process, momconf, triname,
754             triprops, trih, -i], {i, 1, pmp[process] - 1}];
755     tripropss[tri] =
756     RemoveIgnorableOptions[process, triname,
757         SplitPropOptions[process, split, props, tri]], {tri,
758         SplitOptions[process, split]}], {bk1t,
759         Num4V[K1] - S1o\[Chi] Num4V[\[Chi]]};
760     Function[{t,
761         y}, -WithSpinors[
762     Sum[Module[{triname = SplitName[split, tri],
763         insindex = InsertionIndex[split[[1]], tri[[1]], tri[[2]]}],
764         res[tri, t] Sum[
765     Module[{tricorner =
766         MasslessTriCorner[process, triname, triprops], trihs},
767         trihs[h_] :=
768     Evaluate[
769         If[MatchQ[tricorner, _Integer],
770         If[(Abs[tts[tri, tricorner, -1, -1, 1][t]] -
771             Abs[tts[tri, tricorner, -1, 1, 1][t]])/(Abs[
772             tts[tri, tricorner, -1, -1, 1][t]] +
773             Abs[tts[tri, tricorner, -1, 1, 1][t]]) <
774             0, {-h}, {h}], {1, -1}]];
775     Sum[Sum[CalculateTriSubtraction[process, momconf, triname,
776         tricorner, trih, triprops,
777         tts[tri, tricorner, h, trih, 1][
778             t]] h/(y - ypoles[tri, t, h]) +
779         triconts[tri, tricorner, h, trih, triprops,

```

```

780      ttsy[tri, tricorner, h, trih, 1][t, y],
781      ttsy[tri, tricorner, h, trih, -1][t,
782      y]] (h/(y - ypoles[tri, t, h]) -
783      h/(y - ypoles[tri, t, -h])) +
784      triconts[tri, tricorner, h, trih, triprops,
785
786      ttsy[tri, tricorner, h, trih, 1][t,
787      ypoles[tri, t, -h]],
788      ttsy[tri, tricorner, h, trih, -1][t,
789      ypoles[tri, t, -h]]] h/(y -
790      ypoles[tri, t, -h]), {trih, trihs[h]}/
791      Length[trihs[h]], {h, {-1, 1}}], {triprops,
792      tripropss[tri]}], {tri,
793      SplitOptions[process, split]}, {bk1t,
794      Num4V[K1] - S1o\[Chi] Num4V\[Chi]}}]]
795
796 CalculateBubbleSubtractionEqn[process_,
797   momconf_, {split_List, nonord_List}, props_] :=
798 CalculateBubbleSubtractionEqn[process, momconf, {split, nonord},
799   props] =
800 CalculateBubbleSubtractionEqnImpl[process, momconf, {split, nonord},
801   props]
802
803 CalculateBubbleSubtraction[process_,
804   momconf_, {split_List, nonord_List}, props_List, t_, y_] :=
805 CalculateBubbleSubtractionEqn[process, momconf, {split, nonord},
806   props][t, y]
807
808 (* Calculate the raw but subtracted bubble contribution, evaluating as
      much as possible, as early as possible *)
809 CalculateRawBubbleContribution[process_,
810   momconf_, {split_List, nonord_List}, props_, t_, y_] :=
811 Module[{cornertypes =
812   SplitCorners[Process[process], {split, nonord}],
813   cornernames = SplitCorners[MomConf[momconf], {split, nonord}]},
814 CalculateRawBubbleContributionImpl[cornernames, cornertypes, props,
815   CalcBubbleLoopMom[momconf, {split, nonord}, t, y]] -
816 CalculateBubbleSubtraction[process, momconf, {split, nonord},
817   props, t, y]]
818
819 (* The coefficients and values for y needed to extract the correct
      combinations for different maximum powers of y *)

```

```

820 GetyExtraction[1 | 2] := {{1, 1/2}}
821 GetyExtraction[
822   3 | 4] := {{1/2, (3 - Sqrt[3])/6}, {1/2, (3 + Sqrt[3])/6}}
823 GetyExtraction[
824   5 | 6] := {{4/9,
825     1/2}, {5/18, (5 - Sqrt[15])/10}, {5/18, (5 + Sqrt[15])/10}}
826 GetyExtraction[
827   7 | 8] := {{(18 + Sqrt[30])/72, (35 - Sqrt[525 - 70 Sqrt[30]])/
828     70}, {(18 + Sqrt[30])/72, (35 + Sqrt[525 - 70 Sqrt[30]])/
829     70}, {(18 - Sqrt[30])/72, (35 - Sqrt[525 + 70 Sqrt[30]])/
830     70}, {(18 - Sqrt[30])/72, (35 + Sqrt[525 + 70 Sqrt[30]])/70}}
831 GetyExtraction[
832   9 | 10] := {{64/225,
833     1/2}, {(322 + 13 Sqrt[70])/1800, (21 - Sqrt[245 - 14 Sqrt[70]])/
834     42}, {(322 + 13 Sqrt[70])/1800, (21 + Sqrt[245 - 14 Sqrt[70]])/
835     42}, {(322 - 13 Sqrt[70])/1800, (21 - Sqrt[245 + 14 Sqrt[70]])/
836     42}, {(322 - 13 Sqrt[70])/1800, (21 + Sqrt[245 + 14 Sqrt[70]])/42}}
837
838 (* Calculate the needed bubble contribution *)
839 CalculateBubbleContribution[process_,
840   momconf_, {split_List, nonord_List}, props_] :=
841   CalculateBubbleContribution[process, momconf, {split, nonord},
842     props] =
843   Module[{p = pp[process], t0 = pt0[process], yex},
844     yex = GetyExtraction[p];
845     Sum[yex[[k, 1]] CalculateRawBubbleContribution[process,
846       momconf, {split, nonord}, props, t0*Exp[2 Pi I j/(2 p + 1)],
847       yex[[k, 2]]], {k, Length[yex]}, {j, -p, p}]/(2 p + 1) //
848     Factor]

```

B.2 Comparing Mathematica Implementation and BlackHat

Listing B.4: Test.txt

```

1 SetAttributes[PrintTiming, HoldFirst]
2 PrintTiming[expr_] := Module[{tmp = AbsoluteTiming[expr]},
3   Print[tmp[[1]], " to evaluate ", HoldForm[expr]]; tmp[[2]]]
4
5 FailedTests = {};
6 TestIndex = 0;
7

```

```

8 (* Run a test, show its result and if it failed record for later *)
9 SetAttributes[TestEqual, HoldFirst]
10 TestEqual[expr_, value_] := Module[{i, result},
11   i = TestIndex = TestIndex + 1; Print["Starting Test ", i,
12     ": " *Defer[expr], "=", value]; result = PrintTiming[expr];
13   If[TrueQ[result == value], Print["Test succeeded"],
14     Print["Test ", i, " Failed: ", result, "!=" , value];
15     FailedTests = Append[FailedTests, i]]; ]
16 TestEqual[expr_, value_, error_] := Module[{i, result},
17   i = TestIndex = TestIndex + 1; Print["Starting Test ", i,
18     ": " *Defer[expr], "=", value]; result = PrintTiming[expr];
19   If[TrueQ[If[TrueQ[value == 0], Abs[result], (Abs[result - value]*2)/
20     (Abs[result] + Abs[value])] < error], Print["Test succeeded"],
21     Print["Test ", i, " Failed: ", result, "!=" , value];
22     FailedTests = Append[FailedTests, i]]; ]
23
24 (* Show any tests that have failed *)
25 ShowFailedTests [] := If[Length[FailedTests] == 0,
26   Print["No Tests Failed"], Print["Test(s) ",
27     Sequence @@ Riffle[FailedTests, ", ", " ", " Failed"]]]

```

Listing B.5: Rules.txt

```

1 epsilon = 10-7;
2
3 (* Compare two sets of rules recursively. Rules match if the values for
   the same keys match within tolerance except for keys that are only
   in one set of rules whose values must all be zero within tolerance
   *)
4 CompareRules[keys_, math : _Real | _Complex | _Integer ,
5   bh : _Real | _Complex | _Integer] :=
6   If[TrueQ[Abs[math - bh]*2/(0.1 + Abs[math] + Abs[bh]) < epsilon],
7     True, Print["Difference in valid values for ", keys, " ", math,
8       "!=" , bh]; False]
9 CompareRules[keys_, math : {}, bh : _Real | _Complex | _Integer] :=
10  If[TrueQ[Abs[bh] < epsilon], True,
11    Print["bh value not 0 and math missing for ", keys, " ", bh];
12    False]
13 CompareRules[keys_, math : _Real | _Complex | _Integer , bh : {}] :=
14  If[TrueQ[Abs[math] < epsilon], True,
15    Print["math value not 0 and bh missing for ", keys, " ", math];
16    False]
17 CompareRules[keys_, math_List , bh_List] :=

```

```

18 Module[{result = True},
19 If[Not[And @@ (CompareRules[Append[keys, #], # /. math, # /. bh] & /@
20     Intersection[First /@ math, First /@ bh])], result = False;
21 Print["error in common elements for ", keys]];
22 If[Not[And @@ (CompareRules[Append[keys, #], # /. math, {}] & /@
23     Complement[First /@ math, First /@ bh])], result = False;
24 Print["error in elements only in math ", keys]];
25 If[Not[And @@ (CompareRules[Append[keys, #], {}, # /. bh] & /@
26     Complement[First /@ bh, First /@ math])], result = False;
27 Print["error in elements only in bh ", keys]]; result]
28
29 (* Generate a set of nested rules with split as the outer key and
    propagators as the inner key that map to the value of the
    corresponding coefficient *)
30 GenerateMathRules[process_, momconf_, n_, genfunc_] :=
31 Module[{tree = HelAmplN[MomConf[momconf], Process[process]]},
32 Function[split,
33     split -> ((# :>
34         genfunc[process, momconf, split, #, None]/tree) & /@
35         RemoveIgnorableOptions[process, split,
36             SplitPropOptions[process, split]])] /@
37     SplitOptions[process, n]]
38 GenerateMathRules[process_, momconf_, n : 4] :=
39 GenerateMathRules[process, momconf, n, CalculateBoxContribution]
40 GenerateMathRules[process_, momconf_, n : 3] :=
41 GenerateMathRules[process, momconf, n, CalculateTriContribution]
42 GenerateMathRules[process_, momconf_, n : 2] :=
43 GenerateMathRules[process, momconf, n, CalculateBubbleContribution]
44
45 $UseBHSigns = False;

```

Listing B.6: rambo.py by Daniel Maitre

```

1 """ Generate random sets of on-shell momenta that conserve momenta """
2
3 import random
4 import math
5 from numpy import array
6 random.seed(1234)
7
8 def doublePi():
9     return math.pi
10

```

```

11 math.get_pi=doublePi
12
13 def dot(a,b):
14     return sum([a[i]*b[i] for i in range(len(a))],type(a[0])(0) )
15
16 def getRandomQ(RandomGenerator=random.random,Type=float,mathLib=math):
17     c=Type(2)*RandomGenerator()-Type(1)
18     phi=Type(2)*mathLib.get_pi()*RandomGenerator()
19     q0=-mathLib.log(RandomGenerator()*RandomGenerator())
20     qx=q0*mathLib.sqrt(Type(1)-c*c)*mathLib.cos(phi)
21     qy=q0*mathLib.sqrt(Type(1)-c*c)*mathLib.sin(phi)
22     qz=q0*c
23     return (q0,qx,qy,qz)
24
25 def boost(q,x,gamma,b):
26     Type=type(x)
27     p0=x*(gamma*q[0]+dot(b,q[1:]))
28     p= array(q[1:])
29     p+= b*q[0]
30     f=Type(1)/(Type(1)+gamma)
31     f*=dot(b,q[1:])
32     p+=b*f
33     p*=x
34     return (p0,)+tuple(p)
35
36 def finalStatePS(w,n,Type=float,mathLib=math,**kargs):
37     qs = [ getRandomQ(Type=Type,mathLib=mathLib,**kargs) for i in range(
n) ]
38     Q = array([ sum([q[j] for q in qs ],Type(0)) for j in range(4) ])
39     M = mathLib.sqrt(Q[0]*Q[0]-dot(Q[1:],Q[1:]))
40     b = array( [-q/M for q in Q[1:] ])
41     x = Type(w)/M
42     gamma = Q[0]/M
43     ps=[ boost(q,x,gamma,b) for q in qs ]
44     return ps
45
46 def PS(n,Type=float,mathLib=math,RandomGenerator=random.random,**kargs):
47     ctheta=Type(2)*RandomGenerator()-Type(1)
48     stheta=mathLib.sqrt(Type(1)-ctheta*ctheta)
49     phi=Type(2)*RandomGenerator()*mathLib.get_pi()
50     sphi=mathLib.sin(phi)
51     cphi=mathLib.cos(phi)

```

```

52     w=Type(n)
53     E=w/Type(2)
54     c=mathLib
55     p1=(-E,
56         -E*stheta*sphi,
57         -E*stheta*cphi,
58         -E*ctheta
59     )
60     p2 = ( -E , -p1[1] , -p1[2] , -p1[3])
61     ps=finalStatePS(n,n-2,RandomGenerator=RandomGenerator,Type=Type,
62     mathLib=mathLib,**kargs)
63     return [ p1,p2] + ps

```

Listing B.7: BHtools.py based on code by Daniel Maitre

```

1  """ Load BlackHat and declare a few functions to map between strings and
2      BlackHat types """
3  ## uncomment this and insert the correct path to find the
4  ## blackhat python library if it is not already in the
5  ## module import path
6  ##import sys
7  ##sys.path.append('/path/to/blackhat/python/library')
8  import BH
9
10 import re
11
12 import rambo
13
14 def getRandomMC(n):
15     ps =rambo.PS(n)
16     cms = [ BH.Cmomd(*p) for p in ps ]
17     return BH.mcd(*cms)
18
19 stringToParticleMap={
20     'm' : BH.cvar.m ,
21     'p' : BH.cvar.p ,
22     'qm' : BH.cvar.qm ,
23     'qp' : BH.cvar.qp ,
24     'Qm' : BH.cvar.q2m ,
25     'Qp' : BH.cvar.q2p ,
26     'qbm' : BH.cvar.qbm ,
27     'qbp' : BH.cvar.qbp ,

```



```

28     'Qbm' : BH.cvar.qb2m ,
29     'Qbp' : BH.cvar.qb2p ,
30     'ym'  : BH.cvar.ym ,
31     'yp'  : BH.cvar.yp ,
32     'lp'  : BH.cvar.lp ,
33     'lm'  : BH.cvar.lm ,
34     'lbp' : BH.cvar.lbp ,
35     'lbm' : BH.cvar.lbm ,
36     'Um'  : BH.cvar.Qm ,
37     'Up'  : BH.cvar.Qp ,
38     'Ubm' : BH.cvar.Qbm ,
39     'Ubp' : BH.cvar.Qbp ,
40     'ph'  : BH.cvar.ph ,
41     'phd' : BH.cvar.phd ,
42     'H'   : BH.cvar.H ,
43     'q0m' : BH.cvar.qm ,
44     'q0p' : BH.cvar.qp ,
45     'qb0m' : BH.cvar.qbm ,
46     'qb0p' : BH.cvar.qbp ,
47     'q1m' : BH.cvar.q2m ,
48     'q1p' : BH.cvar.q2p ,
49     'qb1m' : BH.cvar.qb2m ,
50     'qb1p' : BH.cvar.qb2p ,
51     'q2m' : BH.cvar.q3m ,
52     'q2p' : BH.cvar.q3p ,
53     'qb2m' : BH.cvar.qb3m ,
54     'qb2p' : BH.cvar.qb3p ,
55     'q3m' : BH.cvar.q4m ,
56     'q3p' : BH.cvar.q4p ,
57     'qb3m' : BH.cvar.qb4m ,
58     'qb3p' : BH.cvar.qb4p
59 }
60
61 def stringToParticles(st):
62     ps=st.split(' ')
63     return [ stringToParticleMap[p] for p in ps ]
64
65 def stringToProcess(st):
66     return BH.process(BH.vectorpID(stringToParticles(st)))

```

Listing B.8: BMathLink.py

```

1 """ Functions to convert between BlackHat and Mathematica implementation
   representations of various objects """
2 import BHtools as BHT
3 import BH
4 from math import log10, ceil
5 import numpy as np
6 from collections import defaultdict
7
8 def genmc(pro):
9     """ Generate a set of momenta for the given process taking account of
   massive particles """
10    nparticles=len(pro)
11    imassive=[i for i in range(nparticles) if pro[i].mass()!=0]
12    nmassive=len(imassive)
13    imassless=[i for i in range(nparticles) if pro[i].mass()==0]
14    nmassless=len(imassless)
15    n=nparticles+nmassive #nmassless+2*nmassive
16    mc=BHT.getRandomMC(n)
17    inds=np.zeros((nparticles),dtype=int)
18    inds[imassless]=range(1,nmassless+1)
19    inds[imassive]=[mc.Sum(i,i+1) for i in range(nparticles-nmassive+1,n
   +1,2)]
20    return mc,[int(i) for i in inds]
21
22 def dasmath(d):
23     """ Write a double in a form that can be read as Mathematica input """
24     if d==0:
25         return "0"
26     exp=int(log10(abs(d)))
27     return str(d/10**exp)+"*"+str(exp)
28 def casmath(c):
29     """ Write a complex number in a form that can be read as Mathematica
   input """
30     return dasmath(c.real)+"+I*"+dasmath(c.imag)
31
32 def makequarkmap(cut,n):
33     """
34     Calculate a map to correct quark flavours to Mathematica versions.
35
36     In corners of a cut, BlackHat relabels quarks to ensure that if the
   same quark line traverses a corner twice, the correct pairs of
   quarks will connect up. The Mathematica implementation does this at

```

```

    the last possible step before evaluating and therefore expects
    quarks to be labelled using their external flavours
37 """
38 quarkmap={}
39 cutquarks=defaultdict(set)
40 for i in range(n):
41     for j in range(cut.corner_size(i+1)):
42         ind=cut.corner_ind(i+1,j+1)
43         externpart=cut.extern_process().p(ind)
44         internpart=cut.get_process(i+1).p(j+2)
45         if externpart.is_a(BH.cvar.quark):
46             # if the particle is a quark then correct from the
47             # type used in the corner to the type used in the
48             # external process
49             quarkmap[internpart.flavor()]=externpart.flavor()
50         cutpart=cut.get_process(i+1).front()
51         # map cut quarks from the type in one corner to the
52         # type used in the previous corner
53         if cutpart.is_a(BH.cvar.quark):
54             otherpart=cut.get_process((i-1)%n+1).back()
55             cutquarks[cutpart.flavor()].add(otherpart.flavor())
56             cutquarks[otherpart.flavor()].add(cutpart.flavor())
57         while cutquarks:
58             # if the quark is not found then there must be a closed quark loop
59             value=-1
60
61             # start from a cut quark and repeatedly search for quarks
62             # that are equivalent and also check if any match an
63             # external quark
64             quarkset=[cutquarks.keys()[0]]
65             newquarks=quarkset[:]
66             while newquarks:
67                 quark=newquarks.pop()
68                 if quark in quarkmap:
69                     value=quarkmap[quark]
70                 for otherquark in cutquarks.pop(quark):
71                     if otherquark not in quarkset:
72                         quarkset.append(otherquark)
73                         newquarks.append(otherquark)
74             for quark in quarkset:
75                 quarkmap[quark]=value
76         return quarkmap

```

```

77
78 _partToMathMap={
79     BH.cvar.m:"Gluon[-1]" ,
80     BH.cvar.p:"Gluon[1]" ,
81     BH.cvar.ph:"Phi[1]" ,
82     BH.cvar.phd:"Phi[-1]" ,
83     BH.cvar.H:"Higgs[]"
84 }
85
86 def partToMathMap(p,quarkmap={}):
87     """ Convert a BlackHat particle to its Mathematica type name """
88     if p.is_a(BH.cvar.quark):
89         return "Quark["+( "-1" if p.is_anti_particle() else "1" )+"," +str(p.
90             helicity())+" ,"+str(quarkmap.get(p.flavor(),p.flavor()))+"]"
91     else:
92         return _partToMathMap[p]
93
94 def partname(p,i):
95     """ Generate a name for a particle compatible with S@M """
96     if p.mass()==0:
97         return str(i)
98     else:
99         return "P"+str(i)
100
101
102 def mcmathprint(pro,mc,inds):
103     """ Convert the momentum configuration to the code to declare the
104         corresponding S@M momenta """
105     s=""
106     for p,i in zip(pro.particle_IDs(),inds):
107         if p.mass()==0:
108             s=s+"DeclareSpinorMomentum["+str(i)+",{ "+casmath(mc.p(i).E())+" ,"+
109                 casmath(mc.p(i).X())+" ,"+casmath(mc.p(i).Y())+" ,"+casmath(mc.p(i).Z
110                 ())+"}]\n"
111         else:
112             s=s+"DeclareLVectorMomentum["+str(i)+",{ "+casmath(mc.p(i).E())+" ,
113                 "+casmath(mc.p(i).X())+" ,"+casmath(mc.p(i).Y())+" ,"+casmath(mc.p(i).
114                 Z())+""}]\n"
115     return s
116
117 def generate_label_maps(pro):

```

```

113 """ Generate a pair of maps containing information needed to map cuts
      from BlackHat to their Mathematica names """
114 ordered_label_map={}
115 unordered_label_map={}
116 i=0
117 j=0
118 for k in range(len(pro)):
119     if pro[k].get_ordered() == 0:
120         j=j+1
121         ordered_label_map[k+1]=j
122     else:
123         i=i+1
124         unordered_label_map[k+1]=i
125 return ordered_label_map , unordered_label_map
126
127 def get_label_indexes((ordered_map , unordered_map)):
128 """ Generate lists containing the indexes of the ordered and unordered
      particles in the BlackHat process """
129 ordered_indexes=[-1]*len(ordered_map)
130 unordered_indexes=[-1]*len(unordered_map)
131 for k,v in ordered_map.items():
132     ordered_indexes[v-1]=k
133 for k,v in unordered_map.items():
134     unordered_indexes[v-1]=k
135 return ordered_indexes , unordered_indexes
136
137 def calculate_raw_code(c,(ordered_map , unordered_map)):
138 """ Calculate the split code from a BlackHat cut, in a form that
      matches the Mathematica implementation """
139 cut_labels=[]
140 unordered_labels=[-1]*len(unordered_map)
141 skipped = 0
142 for i in range(c.nbr_props()):
143     corner = [c.corner_ind(i+1,j+1) for j in range(c.corner_size(i+1))]
144     for ind in [ind for ind in corner if ind in unordered_map]:
145         unordered_labels[unordered_map[ind]-1]=i+1
146         corner.remove(ind)
147     if len(corner) == 0:
148         skipped = skipped+1
149     else:
150         cut_labels.extend([ordered_map[corner[0]]*(skipped+1)])
151         skipped=0

```

```

152 cut_labels.extend([cut_labels[0]*skipped)
153 shifted = 0;
154 target = min(cut_labels)
155 while cut_labels[0] != target or cut_labels[-1] == target:
156     shifted = shifted+1
157     cut_labels.append(cut_labels.pop(0))
158 if shifted != 0:
159     unordered_labels = [((i+len(cut_labels)-shifted-1) % len(cut_labels)
160     ) + 1 for i in unordered_labels]
161 return cut_labels, unordered_labels, shifted
162
163 def calculate_mycode(c,maps):
164     """ Calculate the split code as a string from a BlackHat cut, in a
165     form that matches the Mathematica implementation """
166     cut_labels, unordered_labels, shifted=calculate_raw_code(c,maps)
167     cuts = ".join(str(el) for el in cut_labels)
168     if len(unordered_labels) != 0:
169         cuts += "-".join(str(el) for el in unordered_labels)
170     return cuts, shifted
171
172 def calculate_mathsplit(cut,maps):
173     """ Calculate the Mathematica code for the label for a split """
174     cut_labels, unordered_labels, shifted=calculate_raw_code(cut,maps)
175     return "{{"+"",".join(str(s) for s in cut_labels)+"},{"+"",".join(str(s)
176     for s in unordered_labels)+"}}",shifted
177
178 def sortedcuts(cutfunction, cutcount, maps):
179     """
180     Sort the cuts into a consistent order.
181     Returns a list of tuples containing the BlackHat index of the cut and
182     the cut itself
183     """
184     cuts=[(i+1,cutfunction(i+1)) for i in range(cutcount)]
185     def key((i, cut)):
186         return (calculate_mycode(cut,maps),)+tuple(cut.l(i+1).conjugate()
187         for i in range(cut.size()))
188     cuts.sort(key=key)
189     return cuts
190
191 def getprops(cut, n, shift):

```

```

188 """ Return the propagators for a cut in the order needed for the
      Mathematica implementation """
189 return tuple(p.p(p.n()) for p in (cut.get_process((i+shift-1)%n+1) for
      i in range(n)))
190
191 # Dictionaries to map colour structures to strings and back
192 colourstructToBH={
193     "glue":BH.glue ,
194     "nf":BH.nf ,
195     "LT":BH.LT,
196     "RT":BH.RT,
197     "LLT":BH.LLT,
198     "LRT":BH.LRT,
199     "RLT":BH.RLT,
200     "RRT":BH.RRT}
201
202 colourstructureFromBH={v:k for k,v in colourstructToBH.iteritems()}
203
204 # Dictionary to map colour structures to the Mathematica code for them
205 colourstructureBHtoMath={
206     BH.glue:"{}",
207     BH.nf:"{Quark[1,-1]}",
208     BH.LT:"{Quark[-1,1]}",
209     BH.RT:"{Quark[1,1]}",
210     BH.LLT:"{Quark[-1,1],Quark[-1,2]}",
211     BH.LRT:"{Quark[-1,1],Quark[1,2]}",
212     BH.RLT:"{Quark[1,1],Quark[-1,2]}",
213     BH.RRT:"{Quark[1,1],Quark[1,2]}"}

```

Listing B.9: TreeTests.py

```

1 #!/usr/bin/python
2
3 import BHtools as BHT
4 import BH
5 from BHMathlink import *
6 import sys
7 import os
8 import random
9 random.seed()
10 from itertools import combinations
11 import numpy as np
12

```

```

13 def calctest(pro,mc,inds):
14     ep=BH.ep(mc,inds)
15     A=BH.TreeHelAmpl(pro)
16     return A.eval(ep)
17 def mathcode(pro,mc,inds, val):
18     s=mcmathprint(pro,mc,inds)
19     s=s+"\n"
20     s=s+"TestEqual[HelAmplN[{ "+", ".join(partname(p,i) for p,i in zip(pro.
        particle_IDs(),inds))+"}, {" "+", ".join(partToMathMap(p) for p in pro.
        particle_IDs())+"}], "+casmath(val)+" ,1*^-9]\n"
21     return s
22 def runtest(pro):
23     mc,inds=genmc(pro)
24     val=calctest(pro,mc,inds)
25     return mathcode(pro,mc,inds, val)
26
27 def allcombs(l):
28     for i in range(0,len(l)+1):
29         for c in combinations(l,i):
30             yield c
31
32 def shuffle(els,counts):
33     if sum(counts)==0:
34         yield ()
35     else:
36         for i in range(len(els)):
37             counts[i]-=1
38             for s in shuffle(els,counts):
39                 yield (el[i],)+s
40             counts[i]+=1
41
42 def noshuffle(els,counts):
43     return [tuple(e for e,c in zip(els,counts) for i in range(c))]
44
45 shuffle=noshuffle
46
47 def rungluetests():
48     s=""
49     for n in range(4,10):
50         for nneg in range(n+1):
51             for shuf in shuffle(("p", "m"),(n-nneg,nneg)):
52                 s=s+runtest(BHT.stringToProcess(" ".join(shuf))+"\n\n")

```



```

53  return s
54
55  def runphigluetests():
56  s=""
57  for n in range(4,10):
58  for nneg in range(n+1):
59  for shuf in shuffle(("p","m"),(n-nneg,nneg)):
60  s=s+runtest(BHT.stringToProcess("ph "+ " ".join(shuf))+"\n\n")
61  return s
62
63  def quarkoptions(nquarks,i=0):
64  if nquarks==0:
65  yield []
66  else:
67  qo=[("qp","qbm"),("qm","qbp"),("qbp","qm"),("qbm","qp")]
68  for option in quarkoptions(nquarks-1,i+1):
69  for o in qo:
70  yield [tuple(q[:-1]+str(i)+q[-1] for q in o)]+option
71
72  def insertquarks(others,quarks):
73  others=tuple(others)
74  if len(quarks)==0:
75  yield others
76  else:
77  for i in range(len(others)+1):
78  for rest in insertquarks(others[i:],quarks[1:]):
79  yield others[:i]+quarks[0]+rest
80
81  def pad(it,el):
82  for e in it:
83  yield e
84  while True:
85  yield el
86
87  def grouper(l,n):
88  for i in range(0,len(l),n):
89  yield l[i:i+n]
90
91  def insertquarksone(others,quarks):
92  others=tuple(others)
93  per=max(len(others)/len(quarks),1)
94  l=[e for o,q in zip(pad(grouper(others,per),()),quarks) for e in q+o]

```

```

95  return [tuple(1)+others[len(1):]]
96
97  insertquarks=insertquarksone
98
99  def runquarkgluetests():
100     s=""
101     for N in range(4,10):
102         for nneg in range(N-2,-1,-1):
103             for nquarks in range(1,min((N-nneg)/2,3)+1):
104                 n=N-nquarks*2
105                 for quarks in quarkoptions(nquarks):
106                     for shuf in shuffle(("p","m"),(n-nneg,nneg)):
107                         for withq in insertquarks(shuf,quarks):
108                             print withq;
109                             s=s+runtest(BHT.stringToProcess(" ".join(withq))+"\n\n")
110     return s
111
112  def runphiquarkgluetests():
113     s=""
114     for N in range(4,10):
115         for nneg in range(N-2,-1,-1):
116             for nquarks in range(1,min((N-nneg)/2,3)+1):
117                 n=N-nquarks*2
118                 for quarks in quarkoptions(nquarks):
119                     for shuf in shuffle(("p","m"),(n-nneg,nneg)):
120                         for withq in insertquarks(shuf,quarks):
121                             print withq;
122                             s=s+runtest(BHT.stringToProcess("ph "+withq).join(withq))+"\n\n")
123     return s
124
125  def mainglue():
126     from subprocess import Popen,PIPE
127
128     p = Popen(["math8","-noprompt"], cwd=os.path.dirname(os.path.realpath(
129         __file__)), stdin=PIPE)
130     pipe = p.stdin
131     pipe.write("SetOptions[#,FormatType->OutputForm]&/@Streams[]\n")
132     pipe.write("<<\Common.txt\n")
133     pipe.write("<<\HelAmplN.txt\n")
134     pipe.write("<<\Test.txt\n")

```

```
135
136 pipe.write(runluetests())
137
138 pipe.write("ShowFailedTests []\n")
139 pipe.write("Exit []")
140 pipe.close()
141 p.wait()
142
143 def mainphiglu():
144     from subprocess import Popen, PIPE
145
146     p = Popen(["math8", "-noprompt"], cwd=os.path.dirname(os.path.realpath(
147         __file__)), stdin=PIPE)
148     pipe = p.stdin
149     pipe.write("SetOptions[#,FormatType->OutputForm]&/@Streams []\n")
150
151     pipe.write("<<\nCommon.txt\n\n")
152     pipe.write("<<\nHelAmplN.txt\n\n")
153     pipe.write("<<\nTest.txt\n\n")
154
155     pipe.write(runphigluetests())
156
157     pipe.write("ShowFailedTests []\n")
158     pipe.write("Exit []")
159     pipe.close()
160     p.wait()
161
162 def mainquarkglue():
163     from subprocess import Popen, PIPE
164
165     p = Popen(["math8", "-noprompt"], cwd=os.path.dirname(os.path.realpath(
166         __file__)), stdin=PIPE)
167     pipe = p.stdin
168     pipe.write("SetOptions[#,FormatType->OutputForm]&/@Streams []\n")
169
170     pipe.write("<<\nCommon.txt\n\n")
171     pipe.write("<<\nHelAmplN.txt\n\n")
172     pipe.write("<<\nTest.txt\n\n")
173
174     pipe.write(runquarkgluetests())
175
176     pipe.write("ShowFailedTests []\n")
```

```

175 pipe.write("Exit []")
176 pipe.close()
177 p.wait()
178
179 def mainphiquarkglue():
180     from subprocess import Popen,PIPE
181
182     p = Popen(["math8","-noprompt"], cwd=os.path.dirname(os.path.realpath(
183         __file__)), stdin=PIPE)
184     pipe = p.stdin
185     pipe.write("SetOptions[#,FormatType->OutputForm]&/@Streams []\n")
186
187     pipe.write("<<<\nCommon.txt\n\n")
188     pipe.write("<<<\nHelAmplN.txt\n\n")
189     pipe.write("<<<\nTest.txt\n\n")
190
191     pipe.write(runphiquarkgluetests())
192
193     pipe.write("ShowFailedTests []\n")
194     pipe.write("Exit []")
195     pipe.close()
196     p.wait()
197
198 if __name__=="__main__":
199     import sys
200     name=sys.argv[1] if len(sys.argv)>=2 else "glue"
201     globals()["main"+name]()

```

Listing B.10: OneLoopTests.py

```

1 #!/usr/bin/python
2 print "(*)"
3 import sys
4 import os.path
5 import BHtools as BHT
6 import BH
7 from BHTMathlink import *
8 from collections import defaultdict
9 from subprocess import Popen,PIPE
10
11 BH.use_setting("USEKNOWNFORMULAE no")
12 print "(*)"

```

```

13
14 def getprocesses (cut ,n) :
15     return [cut .get_process (i+1) for i in range(n)]
16
17 def collectcuts (cuts ,ncuts ,n,maps) :
18     splits=defaultdict (dict)
19     for i in range(ncuts) :
20         cut=cuts (i+1)
21         split ,shifted = calculate_mathsplit (cut ,maps)
22         splits [split ][getprops (cut ,n ,shifted)]=(makequarkmap (cut ,n) ,cut)
23     return splits
24
25 def run (PRO,mode=BH .glue ,doprint=True ,mathtestcode=True ,plotgraph=False ,
26         outfile=sys .stdout) :
27
28     mc,inds=genmc (PRO)
29
30     ep=BH .ep (mc,inds)
31
32     print "(*)#make the next few lines output be treated by mathematica as
33         a comment
34
35     A=BH .TreeHelAmpl (PRO)
36
37     tree=A .eval (ep)
38
39     if doprint :
40         print PRO
41         print tree
42
43     AA=BH .One_Loop_Helicity_Amplitude (PRO,mode)
44
45     cc=AA .cut_part ()
46     cp=cc .makeDarrenCutPart ()
47     if cp==None :
48         cp=cc .makeHiggsCutPart ()
49     if cp==None :
50         print "can't convert cut part to known type"
51         return
52     print cp
53
54     print "Generating Maps: "+str (PRO)

```

```

53  print len(PRO)
54
55  maps = generate_label_maps(PRO)
56
57  print maps
58
59  if doprint:
60      for (i,c) in sortedcuts(cp.bubble, cp.nbr_bubbles(), maps):
61          res=c.eval(ep)
62          print "Bubble {i:{pad}} code {code}/{mycode} [ {processes[0]} |
        {processes[1]} ]: {res.real:< 16.10g}{res.imag:<+16.10g}i | {norm.
        real:< 16.10g}{norm.imag:<+16.10g}i".format(i=i, pad=len(str(cp.
        nbr_bubbles())), code=c.get_code(), mycode=calculate_mycode(c, maps)
        [0], processes=getprocesses(c,2), res=res, norm=res/tree)
63
64
65  for (i,c) in sortedcuts(cp.triangle, cp.nbr_triangles(), maps):
66      res=c.eval(ep)
67      print "Triangle {i:{pad}} code {code}/{mycode} [ {processes[0]}
        | {processes[1]} | {processes[2]} ]: {res.real:< 16.10g}{res.imag
        :<+16.10g}i | {norm.real:< 16.10g}{norm.imag:<+16.10g}i".format(i=i,
        pad=len(str(cp.nbr_triangles())), code=c.get_code(), mycode=
        calculate_mycode(c, maps)[0], processes=getprocesses(c,3), res=res, norm
        =res/tree)
68
69  for (i,c) in sortedcuts(cp.box, cp.nbr_boxes(), maps):
70      res=c.eval(ep)
71      print "Box {i:{pad}} code {code}/{mycode} [ {processes[0]} | {
        processes[1]} | {processes[2]} | {processes[3]} ]: {res.real:< 16.10
        g}{res.imag:<+16.10g}i | {norm.real:< 16.10g}{norm.imag:<+16.10g}i".
        format(i=i, pad=len(str(cp.nbr_boxes())), code=c.get_code(), mycode=
        calculate_mycode(c, maps)[0], processes=getprocesses(c,4), res=res, norm
        =res/tree)
72
73
74  print "*"
75
76  if mathtestcode:
77      try:
78          if mode==BH.nf:
79              sign=-1
80          else:

```

```

81     sign=1
82
83     s=mcmathprint(PRO,mc,inds)+"\n"
84
85     ordered_indexes , unordered_indexes=get_label_indexes (maps)
86
87     s+=" process=DeclareProcess [{"+" , ".join (partToMathMap(PRO[i-1])
for i in unordered_indexes)+"}, {"+" , ".join (partToMathMap(PRO[i-1])
for i in ordered_indexes)+"}], "+colourstructureBHtoMath [mode]+" "\n"
88     s+=" momconf=DeclareMomConf [{"+" , ".join (partname(PRO[i-1], inds [i
-1]) for i in unordered_indexes)+"}, {"+" , ".join (partname(PRO[i-1],
inds [i-1]) for i in ordered_indexes)+"} ] "\n"
89     s+="\n"
90     s+=" CompareRules [ \ "TREE\ " , Abs [ HelAmplN [ MomConf [ momconf ] , Process [
process ] ] ] , "+dasmath ( abs ( tree ) ) + " "\n"
91     s+=" HelAmplN [ MomConf [ momconf ] , Process [ process ] ] / ( "+casmath ( tree ) +
) "\n\n"
92
93     if not doprint :
94         # force the cuts to evaluate themselves (only first copy is
apparently needed)
95         for ( i , c ) in sortedcuts ( cp . bubble , cp . nbr_bubbles () , maps ) :
96             res=c. eval ( ep )
97         for ( i , c ) in sortedcuts ( cp . triangle , cp . nbr_triangles () , maps ) :
98             res=c. eval ( ep )
99
100     splits=collectcuts ( cp . box , cp . nbr_boxes () , 4 , maps )
101     s+=" bh={"+" , ".join ( split+"->{"+" , ".join ( "{"+" , ".join (
partToMathMap(p, quarkmap) for p in prop)+"}->"+casmath ( sign*cut. eval
(ep)/tree) for prop, (quarkmap, cut) in els.iteritems ())+"} "\n" for
split , els in splits.iteritems ())+" } ; "\n"
102     s+=" CompareRules [ {} , GenerateMathRules [ process , momconf , 4 ] , bh ] "\n"
103
104     splits=collectcuts ( cp . triangle , cp . nbr_triangles () , 3 , maps )
105     s+=" bh={"+" , ".join ( split+"->{"+" , ".join ( "{"+" , ".join (
partToMathMap(p, quarkmap) for p in prop)+"}->"+casmath ( sign*cut. eval
(ep)/tree) for prop, (quarkmap, cut) in els.iteritems ())+"} "\n" for
split , els in splits.iteritems ())+" } ; "\n"
106     s+=" CompareRules [ {} , GenerateMathRules [ process , momconf , 3 ] , bh ] "\n"
107
108     splits=collectcuts ( cp . bubble , cp . nbr_bubbles () , 2 , maps )

```

```

109     s+="bh={"+", ".join(split+"->{"+"", ".join("{"+"", " ".join(
partToMathMap(p,quarkmap) for p in prop)+"->" +casmath(sign*cut.eval
(ep)/tree) for prop,(quarkmap,cut) in els.iteritems()+"}\n" for
split,els in splits.iteritems()+"};\n"
110     s+="CompareRules[{}],GenerateMathRules[process,momconf,2],bh]\n"
111
112
113     except Exception as ex:
114         print "(* ERROR *)"
115         print s
116         raise
117
118     print "\n\n(*****MATH CODE*****)\n\n"
119     if not (outfile==sys.stdout or outfile==sys.stderr):
120         print s;
121         print >>outfile,s,"\n\n"
122         print "(*****END MATH CODE*****)\n"
123
124     if plotgraph:
125         path='tree_structure-'+str(PRO)+'-'+colourstructureFromBH[mode]
126
127         import os
128         os.system('mkdir -p \''+s+'\'' % path)
129         BH.print_cut_part_graph(cp,path)
130         os.system('cd \''+s+'\'' ; make all > /dev/null' % path)
131
132     return A,AA,cp
133
134     class MathStream(object):
135     def __init__(self):
136         self.p=None
137     def connect(self):
138         if self.p==None:
139             self.p = Popen(["math8","-noprompt"], cwd=os.path.dirname(os.path.
realpath(__file__)), stdin=PIPE)
140             self.p.stdin.write("SetOptions[#,FormatType->OutputForm]&/@Streams
[]\n")
141             self.p.stdin.write("<<<Common.txt\n\n")
142             self.p.stdin.write("<<<HelAmplN.txt\n\n")
143             self.p.stdin.write("<<<Loop-Cuts.txt\n\n")
144             self.p.stdin.write("<<<Rules.txt\n\n")
145     def write(self,st):

```



```

146     if self.p==None:
147         self.connect()
148         self.p.stdin.write(st)
149 def close(self):
150     if self.p!=None:
151         self.p.stdin.close()
152         self.p.wait()
153         self.p=None
154
155 def parseargs(args,mathstream):
156     if len(args) == 0:
157         print "Script to run tests of the BlackHat and Mathematica
158             implementations of one loop cut amplitudes."
159         print
160         print "Arguments are: ([OPTIONS] PROCESS|'!' COLOUR.STRUCTURE|'*')
161             ..."
162         print "Valid options are '--[no]print' to turn on/off printing
163             output to the console"
164         print "
165             '--[no]mathtestcode' to turn on/off
166             generating Mathematica code"
167         print "
168             '--[no]plotgraph' to turn on/off generating
169             the graph of cuts contributing to the amplitude."
170         print "
171             '--out' FILE to specify where the
172             Mathematica code should be written to."
173         print "
174             FILE can be one of the special values '
175             stdout' or 'stderr'. It can also be '/math/' to request that the
176             Mathematica code be sent directly to a command line instance of
177             Mathematica"
178
179     files={"stdout":sys.stdout,"stderr":sys.stderr,"/math/":mathstream}
180     currentoptions={"doprint":True,"mathtestcode":True,"plotgraph":False,"
181         outfile":sys.stdout}
182     longoptions={"print":{"doprint":True},
183         "noprint":{"doprint":False},
184         "mathtestcode":{"mathtestcode":True},
185         "nomathtestcode":{"mathtestcode":False},
186         "plotgraph":{"plotgraph":True},
187         "noplotgraph":{"plotgraph":False}}
188     shortoptions={"t":{"mathtestcode":True},
189         "n":{"mathtestcode":False},
190         "T":{"doprint":False,"mathtestcode":True,"plotgraph":
191         False},

```

```

177         "p":{"doprint":True},
178         "q":{"doprint":False},
179         "g":{"plotgraph":True}}
180 changesincelast=False
181 torun=[]
182 args=list(args)
183 while args:
184     arg=args.pop(0)
185     if arg[0]=="-":
186         if arg[1]=="-":
187             if arg[2]=="out":
188                 fname=args.pop(0);
189                 if fname not in files:
190                     try:
191                         files[fname]=open(fname,"w")
192                     except IOError as ex:
193                         print ex
194                 if fname in files:
195                     currentoptions["outfile"]=files[fname]
196             else:
197                 currentoptions.update(longoptions[arg[2:]])
198         else:
199             for a in arg[1:]:
200                 currentoptions.update(shortoptions[a])
201     changesincelast=True
202 else:
203     if arg!="!":
204         PRO=BHT.stringToProcess(arg)
205     arg=args.pop(0)
206     if arg!="*":
207         mode=colourstructToBH[arg]
208     torun.append(((PRO,mode),dict(currentoptions)))
209     changesincelast=False
210 if changesincelast:
211     print>(*WARNING: Trailing options what will be ignored *)
212 if len(torun)==0:
213     print>(*ERROR: Nothing to run *)
214 return torun
215
216 def main(args):
217     mathstream=MathStream()
218     for a,kwa in parseargs(args,mathstream):

```

```

219     run(*a,**kwa)
220     mathstream.close()
221
222     import sys
223     main(sys.argv[1:])

```

Listing B.11: DrawCuts.txt

```

1 LaTeXValToSign[1] := "+"
2 LaTeXValToSign[-1] := "-"
3 MakeLaTeXName[Gluon[h_], name_] :=
4 "g_{ } <> ToString[name] <> }^" <> LaTeXValToSign[h]
5 MakeLaTeXName[Phi[1], name_] := "\\phi_{ } <> ToString[name] <> }"
6 MakeLaTeXName[Phi[-1], name_] :=
7 "\\bar\\phi_{ } <> ToString[name] <> }"
8 MakeLaTeXName[Quark[1, h_, f_], name_] :=
9 "q_{ } <> ToString[f] <> ,," <> ToString[name] <> }^" <>
10 LaTeXValToSign[h]
11 MakeLaTeXName[Quark[-1, h_, f_], name_] :=
12 "\\bar q_{ } <> ToString[f] <> ,," <> ToString[name] <> }^" <>
13 LaTeXValToSign[h]
14 MakeLaTeXName[-, name_] := "?_{ } <> ToString[name] <> }"
15 LaTeXLineType[_Phi] := "dashes"
16 LaTeXLineType[Quark[-1, -, -]] := "plain"
17 LaTeXLineType[Quark[1, -, -]] := "fermion"
18 LaTeXLineType[_Gluon] := "gluon"
19 LaTeXLineType[-] := "dots"
20 ReverseLine[Quark[-1, -, -]] := True;
21 ReverseLine[-] := False;
22
23 MakeFeynMFSide[0] := ""
24 MakeFeynMFSide[x.?Positive] := ",left=" <> ToString[x];
25 MakeFeynMFSide[x.?Negative] := ",left=" <> ToString[x];
26
27 MakeFeynMFLine[v1_, v2_, prop_, options_: "", side_: 0] :=
28 MakeFeynMFLine[v2, v1, ReverseParticle[prop], options, -side] /;
29 ReverseLine[prop]
30 MakeFeynMFLine[v1_, v2_, prop_, options_: "", side_: 0] :=
31 StringJoin["\\fmf{", LaTeXLineType[prop], options,
32 MakeFeynMFSide[side], "}" , v1, ",", v2, "]"]
33 MakeFeynMFLabels[i_, v1_, v2_, prop_] :=
34 MakeFeynMFLabels[i, v2, v1, ReverseParticle[prop], "right"] /;
35 ReverseLine[prop]

```

```

36 MakeFeynMFLabels[i_, v1_, v2_, prop_, side_: "left"] :=
37 Module[{path = StringJoin["vpath", ToString[i], "--", v1, "--", v2, ""]
    }],
38 Sow[StringJoin["\\fmfi{phantom,label=${",
39   MakeLaTeXName[ReverseParticle[prop], "1" <> ToString[i]],
40   "}${label.side=", side, "{subpath(0.6 length(", path,
41   ")", 0.8 length(", path, ")} of ", path, "}"]];
42 Sow[StringJoin["\\fmfi{phantom,label=${",
43   MakeLaTeXName[prop, "1" <> ToString[i]], "}${label.side=", side,
44   "}${subpath(0.2 length(", path, ")", 0.4 length(", path, ")} of ",
45   path, "}"]]]]
46
47 MakeFeynMF[process_, momconf_, split_, props_] :=
48 StringJoin[
49   Riffle[Reap[
50     Module[{vname, cornernames, cornertypes},
51       cornernames = SplitCorners[MomConf[momconf], split];
52       cornertypes = SplitCorners[Process[process], split];
53       Sow[StringJoin["\\fmfsurround{",
54         Riffle[ToString /@ Range[Length[Flatten[cornernames]]] //
55         Reverse, ",", "}], "}],];
56       Table[vname[Flatten[cornernames][[i]]] = ToString[i];
57       Sow[StringJoin["\\fmfv{label=${",
58         MakeLaTeXName[Flatten[cornertypes][[i]],
59         Flatten[cornernames][[i]], "}${", ToString[i], "}], {i,
60         Length[Flatten[cornernames]]};
61       Table[Table[
62         Sow[MakeFeynMFLine["v" <> ToString[i],
63           vname[Flatten[cornernames][[i]][[j]],
64           Flatten[cornertypes][[i]][[j]], "", 0], {j,
65           Length[Flatten[cornertypes][[i]]}];
66         Sow[MakeFeynMFLine[
67           "v" <> ToString[Mod[i - 2, Length[split[[1]]] + 1],
68           "v" <> ToString[i], props[[i]], " ,tension=0.4,tag=" <> ToString[i
69         ],
69         If[Length[split[[1]]] == 2, 0.6, 0]], {i,
70         Length[split[[1]]}]; Sow["\\fmffreeze"];
71       Table[
72         MakeFeynMFLabels[i,
73         "v" <> ToString[Mod[i - 2, Length[split[[1]]] + 1],
74         "v" <> ToString[i], props[[i]], {i,
75         Length[split[[1]]}]]][[2, 1], "\n"]

```

```

76
77 MyRun[cmd_] :=
78   Module[{val = Run[cmd]},
79     If[TrueQ[val == 0], 0, Throw[{cmd, val}, Run]]]
80
81 MakeFeyn[process_, momconf_, split_, props_] := Module[{code = "
82   \nonstopmode
83   \pdfminorversion=3
84   \documentclass{letter}
85   \usepackage[usenames]{color} %used for font color
86   \usepackage{amssymb} %maths
87   \usepackage{amsmath} %maths
88   \usepackage[utf8]{inputenc} %useful to type directly diacritic \
89 characters
90   \usepackage{graphicx}
91   \usepackage[outdir=./]{epstopdf}
92   \DeclareGraphicsRule{*}{mps}{*}{}
93   \usepackage{feynmp}
94   \begin{document}
95   \thispagestyle{empty}
96   \begin{fmffile}{mathfig}
97   \begin{fmfgraph*}(120,120)
98     <> MakeFeynMF[process, momconf, split, props] <> "
99   \end{fmfgraph*}
100  \end{fmffile}
101  \end{document}", dir = CreateDirectory[], img},
102 SetDirectory[dir]; Export[dir <> "/math.tex", code, "Text"];
103 Catch[MyRun["pdflatex math"]; MyRun["mpost mathfig.mp"];
104 MyRun["pdflatex math"]; MyRun["pdflatex math"];
105 MyRun["pdfcrop --margin 10 math math-crop.pdf"];
106 img = Import[dir <> "/math-crop.pdf", ImageSize -> 1000][[1]];
107 ResetDirectory[]; DeleteDirectory[dir, DeleteContents -> True];
108 Show[img, ImageSize -> 300],
109 Run, (ResetDirectory[]);
110 Print[Row[{"Error: ", #1, " for ", dir}]] &]]
111
112 GetFeyn[process_, momconf_, split_, props_] :=
113   Module[{val = MakeFeyn[process, momconf, split, props]},
114     If[TrueQ[val[[0]] == Graphics],
115       GetFeyn[process, momconf, split, props] = val, val]]
116
117 GenerateAllDiagrams[proc_, momconf_, n_, f_ : (Grid[#1] -> #3 &)] :=

```

```

118 Function[split ,
119   f[split , #, GetFeyn[proc , momconf, split , #]] & /@
120   RemoveIgnorableOptions[proc , split ,
121   SplitPropOptions[proc , split]]] /@ SplitOptions[proc , n] //
122 Flatten
123
124 ForAllDiagrams[proc_ , momconf_ , n_ , f_ : (#1 -> #2 &)] :=
125 Function[split ,
126   f[split , #] & /@
127   RemoveIgnorableOptions[proc , split ,
128   SplitPropOptions[proc , split]]] /@ SplitOptions[proc , n] //
129 Flatten
130
131 CalculateDiagramDependancies[proc_ , momconf_ , n_] :=
132 ForAllDiagrams[proc , momconf , n ,
133 Function[{split , props} ,
134 Function[
135   newSplit , ({SplitName[split , newSplit] , #} -> {split , props}) & /@
136   RemoveIgnorableOptions[proc , SplitName[split , newSplit] ,
137   SplitPropOptions[proc , split , props , newSplit]]] /@
138   SplitOptions[proc , split]]]
139 ShowDependencyGraph[proc_ , momconf_] :=
140 Module[{FMFVertexShapeFunction} ,
141   FMFVertexShapeFunction[pos_ , {split_ , props_} , size_ , args_...] :=
142   Inset[GetFeyn[proc , momconf, split , props] , pos , {0, 0} , size];
143   Graph[{CalculateDiagramDependancies[proc , momconf, 2] ,
144     CalculateDiagramDependancies[proc , momconf, 3]} // Flatten ,
145     VertexShapeFunction -> FMFVertexShapeFunction , VertexSize -> 1.5 ,
146     DirectedEdges -> False , GraphLayout -> "SpringEmbedding"]
147
148 GenerateDependencyGraphDot[procP_ , momconfP_ , name_ : "graph"] :=
149 Module[{ToN$Value = 0 , ToN$Values = {} , ToN , dir = CreateDirectory [] ,
150   spec , graph} , ToN[1_] := (ToN$Value = ToN$Value + 1;
151   ToN$Values = Union[Append[ToN$Values , 1]]);
152   ToN[1] = ToString[ToN$Value];
153   spec = StringJoin[
154     Riffle[{ "digraph" , "node[label=\"\ , shape=none]" ,
155       ToN[#[[1]]] < "→" < ToN[#[[2]]] < ";" & /@
156       CalculateDiagramDependancies[procP , momconfP , 2] ,
157       ToN[#[[1]]] < "→" < ToN[#[[2]]] < ";" & /@
158       CalculateDiagramDependancies[procP , momconfP , 3] ,
159       StringJoin["{rank=same;" , Riffle[# , ";" , "]" & /@

```

```

160     Map[ToString[#[[1]]] &,
161       GatherBy[
162         SortBy[{ToN[#], Length[#[[1, 1]]]} & /@ ToN$Values, Last],
163         Last], {2}],
164     ToN[#] <> "[image=\"" <>
165       Export[dir <> "/" <> ToN[#] <> ".eps",
166         GetFeyn[procP, momconfP, #[[1]], #[[2]]] <> "\"];"] & /@
167     ToN$Values, "}] // Flatten, "\n"]];
168 graph = Export[dir <> "/graph.txt", spec];
169 Catch[MyRun[
170   "dot -Tps < " <> graph <> " > " <> dir <> "/graph.eps"];
171 MyRun["eps2eps " <> dir <> "/graph.eps " <> dir <> "/graph2.eps"];
172 MyRun["ps2pdf -dEPSCrop " <> dir <> "/graph2.eps " <> name <>
173   ".pdf"]; DeleteDirectory[dir, DeleteContents -> True];,
174 Run, Print[Row[{"Error: ", #1, " for ", dir}]] &]]

```

Appendix C

6 Dimensional Spinor Helicity Implementation

The code created as part of this project can also be downloaded from <http://bit.ly/2oXSBSu>.

Listing C.1: 6DSpinorHelicity.txt

```

1 (* Declare the dimension of vectors to enable extra simplifications *)
2 Dimension[p_] := None;
3 DeclareVectorDimension[p_, d_?EvenQ] := Dimension[p] = d;
4 UndeclareVectorDimension[p_] := Dimension[p] =.;
5
6 (* TraditionalForm and/or StandardForm representations for the various
   objects *)
7 SetAttributes[DoWith, HoldAll]
8 DoWith[s_, v_, e_] := Module[{tmp}, s = v; tmp = e; s =.; tmp]
9
10 MakeBoxes[spinor[p_, hs_], f : TraditionalForm | StandardForm] ^:=
11   MakeBoxes[Subscript[u, hs][p], f]
12 MakeBoxes[spinorbar[p_, hs_], f : TraditionalForm | StandardForm] ^:=
13   MakeBoxes[Subscript[UnderBar[u], hs][p], f]
14 Subscript[u, hs_][p_] := spinor[p, hs]
15 Subscript[UnderBar[u], hs_][p_] := spinorbar[p, hs]
16 AngleBracket = Sp;
17 MakeBoxes[Sp[a_], f : TraditionalForm] ^:=
18   MakeBoxes[AngleBracket[a], f]
19 MpIndexSymbolIndex = 0;
20 MakeBoxes[Mp[(ma : Mom | MomM)[a_], (mb : Mom | MomM)[b_]],

```



```

21  f : TraditionalForm] ^:=
22  RowBox[{SubscriptBox[If[TrueQ[ma == Mom], "p", "P"],
23    MakeBoxes[a, f]], "\[CenterDot]",
24    SubscriptBox[If[TrueQ[mb == Mom], "p", "P"], MakeBoxes[b, f]]}]
25  MakeBoxes[Mp[a_, b_], f : TraditionalForm] ^:=
26  RowBox[{Block[{MpIndexSymbol = MakeBoxes[\[Mu], f],
27    MpIndexUp = True}, MakeBoxes[a, f]], "\[CenterDot]",
28    Block[{MpIndexSymbol = MakeBoxes[\[Mu], f], MpIndexUp = False},
29    MakeBoxes[b, f]]}]
30  MakeBoxes[\[Gamma], f : TraditionalForm] :=
31  SubscriptBox["\[Gamma]", MpIndexSymbol] /; ! MpIndexUp
32  MakeBoxes[\[Gamma], f : TraditionalForm] :=
33  SuperscriptBox["\[Gamma]", MpIndexSymbol] /; MpIndexUp
34  MakeBoxes[\[Sigma][1], f : TraditionalForm] :=
35  SubscriptBox["\[Sigma]", MpIndexSymbol] /; ! MpIndexUp
36  MakeBoxes[\[Sigma][1], f : TraditionalForm] :=
37  SuperscriptBox["\[Sigma]", MpIndexSymbol] /; MpIndexUp
38  MakeBoxes[\[Sigma][-1], f : TraditionalForm] :=
39  SubscriptBox[OverscriptBox["\[Sigma]", "~"], MpIndexSymbol] /; !
40  MpIndexUp
41  MakeBoxes[\[Sigma][-1], f : TraditionalForm] :=
42  SuperscriptBox[OverscriptBox["\[Sigma]", "~"], MpIndexSymbol] /;
43  MpIndexUp
44  MakeBoxes[Mom[i_], f : TraditionalForm] :=
45  SubscriptBox["p", RowBox[{MakeBoxes[i, f], MpIndexSymbol}]] /; !
46  MpIndexUp
47  MakeBoxes[Mom[i_], f : TraditionalForm] :=
48  SubsuperscriptBox["p", MakeBoxes[i, f], MpIndexSymbol] /; MpIndexUp
49  MakeBoxes[MomM[i_], f : TraditionalForm] :=
50  SubscriptBox["P", RowBox[{MakeBoxes[i, f], MpIndexSymbol}]] /; !
51  MpIndexUp
52  MakeBoxes[MomM[i_], f : TraditionalForm] :=
53  SubsuperscriptBox["P", MakeBoxes[i, f], MpIndexSymbol] /; MpIndexUp
54  MakeBoxes[Mp[a_, b_, i_], f : TraditionalForm] ^:=
55  RowBox[{DoWith[MpIndexUp$$[i], True, MakeBoxes[a, f]],
56    DoWith[MpIndexUp$$[i], False, MakeBoxes[b, f]]}]
57  MakeBoxes[\[Gamma][i_], f : TraditionalForm] :=
58  SubscriptBox["\[Gamma]", MakeBoxes[i, f]] /; ! MpIndexUp$$[i]
59  MakeBoxes[\[Gamma][i_], f : TraditionalForm] :=
60  SuperscriptBox["\[Gamma]", MakeBoxes[i, f]]
61  MakeBoxes[\[Sigma][1][i_], f : TraditionalForm] :=
62  SubscriptBox["\[Sigma]", MakeBoxes[i, f]] /; ! MpIndexUp$$[i]

```

```

63 MakeBoxes[\[Sigma][1][i_], f : TraditionalForm] :=
64 SuperscriptBox["\[Sigma]", MakeBoxes[i, f]]
65 MakeBoxes[\[Sigma][-1][i_], f : TraditionalForm] :=
66 SubscriptBox[OverscriptBox["\[Sigma]", "~"], MakeBoxes[i, f]] /; !
67 MpIndexUp$$[i]
68 MakeBoxes[\[Sigma][-1][i_], f : TraditionalForm] :=
69 SuperscriptBox[OverscriptBox["\[Sigma]", "~"], MakeBoxes[i, f]]
70 MakeBoxes[Mom[p_][i_], f : TraditionalForm] :=
71 SubscriptBox["p", RowBox[{MakeBoxes[p, f], MakeBoxes[i, f]}]] /; !
72 MpIndexUp$$[i]
73 MakeBoxes[Mom[p_][i_], f : TraditionalForm] :=
74 SubsuperscriptBox["p", MakeBoxes[p, f], MakeBoxes[i, f]]
75 MakeBoxes[Metric[i_, j_], f : TraditionalForm] :=
76 SuperscriptBox["g", RowBox[{MakeBoxes[i, f], MakeBoxes[j, f]}]]
77 MakeBoxes[Metric[i_, j_], f : TraditionalForm] :=
78 SubscriptBox["g",
79 RowBox[{MakeBoxes[i, f], MakeBoxes[j, f]}]] /; !
80 MpIndexUp$$[i] && ! MpIndexUp$$[j]
81 MakeBoxes[Metric[i_, j_], f : TraditionalForm] :=
82 SubscriptBox[SuperscriptBox["g", MakeBoxes[i, f]],
83 MakeBoxes[j, f]] /; ! MpIndexUp$$[j]
84 MakeBoxes[Metric[i_, j_], f : TraditionalForm] :=
85 SuperscriptBox[SubscriptBox["g", MakeBoxes[i, f]],
86 MakeBoxes[j, f]] /; ! MpIndexUp$$[i]
87 MakeBoxes[CalculatedP[mom_], f : TraditionalForm] :=
88 RowBox[{SubscriptBox["p", "calc"], " ", MakeBoxes[mom, f], " "}]
89 MakeBoxes[Mom[i_], f : TraditionalForm] :=
90 SubscriptBox["p", MakeBoxes[i, f]]
91 MakeBoxes[Mom[i_], f : TraditionalForm] :=
92 SubscriptBox["p", MakeBoxes[i, f]]
93
94 MakeBoxes[shift[{p_, q_}, z_, h_List, l_List][p_],
95 f : TraditionalForm | StandardForm] :=
96 MakeBoxes[Subsuperscript[OverHat[p] -> q, Row[h], Row[1]][z], f]
97 MakeBoxes[shift[{p_, q_}, z_, h_List, l_List][q_],
98 f : TraditionalForm | StandardForm] :=
99 MakeBoxes[Subsuperscript[p -> OverHat[q], Row[h], Row[1]][z], f]
100 MakeBoxes[Mom[shift[{p_, q_}, z_, h_List, l_List][p_]],
101 f : TraditionalForm] :=
102 SubscriptBox[
103 MakeBoxes[
104 Subsuperscript[OverHat[Subscript["p", p]] -> Subscript["p", q],

```

```

105      Row[h], Row[l]][z], f], MpIndexSymbol] /; ! MpIndexUp
106 MakeBoxes[Mom[shift[{p-, q-}, z-, h_List, l_List][q-]],
107   f : TraditionalForm] :=
108   SubscriptBox[
109     MakeBoxes[
110       Subsuperscript[Subscript["p", p] -> OverHat[Subscript["p", q]],
111         Row[h], Row[l]][z], f], MpIndexSymbol] /; ! MpIndexUp
112 MakeBoxes[Mom[shift[{p-, q-}, z-, h_List, l_List][p-]],
113   f : TraditionalForm] :=
114   SuperscriptBox[
115     MakeBoxes[
116       Subsuperscript[OverHat[Subscript["p", p]] -> Subscript["p", q],
117         Row[h], Row[l]][z], f], MpIndexSymbol] /; MpIndexUp
118 MakeBoxes[Mom[shift[{p-, q-}, z-, h_List, l_List][q-]],
119   f : TraditionalForm] :=
120   SuperscriptBox[
121     MakeBoxes[
122       Subsuperscript[Subscript["p", p] -> OverHat[Subscript["p", q]],
123         Row[h], Row[l]][z], f], MpIndexSymbol] /; MpIndexUp
124 MakeBoxes[Mom[shift[{p-, q-}, z-, h_List, l_List][p-]][i-],
125   f : TraditionalForm] :=
126   SubscriptBox[
127     MakeBoxes[
128       Subsuperscript[OverHat[Subscript["p", p]] -> Subscript["p", q],
129         Row[h], Row[l]][z], f], MakeBoxes[i, f]] /; ! MpIndexUp$$[i]
130 MakeBoxes[Mom[shift[{p-, q-}, z-, h_List, l_List][q-]][i-],
131   f : TraditionalForm] :=
132   SubscriptBox[
133     MakeBoxes[
134       Subsuperscript[Subscript["p", p] -> OverHat[Subscript["p", q]],
135         Row[h], Row[l]][z], f], MakeBoxes[i, f]] /; ! MpIndexUp$$[i]
136 MakeBoxes[Mom[shift[{p-, q-}, z-, h_List, l_List][p-]][i-],
137   f : TraditionalForm] :=
138   SuperscriptBox[
139     MakeBoxes[
140       Subsuperscript[OverHat[Subscript["p", p]] -> Subscript["p", q],
141         Row[h], Row[l]][z], f], MakeBoxes[i, f]]
142 MakeBoxes[Mom[shift[{p-, q-}, z-, h_List, l_List][q-]][i-],
143   f : TraditionalForm] :=
144   SuperscriptBox[
145     MakeBoxes[
146       Subsuperscript[Subscript["p", p] -> OverHat[Subscript["p", q]],

```

```

147      Row[h], Row[l]][z], f], MakeBoxes[i, f]]
148 MakeBoxes[Mom[shift[{p-, q-}, z-, h-List, l-List][p-]],
149   f : TraditionalForm] :=
150 MakeBoxes[
151   Subsuperscript[OverHat[Subscript["p", p]] -> Subscript["p", q],
152     Row[h], Row[l]][z], f]
153 MakeBoxes[Mom[shift[{p-, q-}, z-, h-List, l-List][q-]],
154   f : TraditionalForm] :=
155 MakeBoxes[
156   Subsuperscript[Subscript["p", p] -> OverHat[Subscript["p", q]],
157     Row[h], Row[l]][z], f]
158 Subsuperscript[OverHat[p-] -> q-, Row[h-], Row[l-]][z-] :=
159 shift[{p, q}, z, h, l][p]
160 Subsuperscript[p- -> OverHat[q-], Row[h-], Row[l-]][z-] :=
161 shift[{p, q}, z, h, l][q]
162
163 Subscript["\[PlusMinus]", p-] := HelicitySign[p]
164 MakeBoxes[HelicitySign[p-], f : TraditionalForm | StandardForm] ^=
165 MakeBoxes[Subscript["\[PlusMinus]", p], f]
166
167 (* An object representing a sign whose value may not be known yet but
      which can still be simplified since it is known that it must be
      either 1 or -1 *)
168 HelicitySign[a-, b-] := HelicitySign[a] HelicitySign[b]
169 HelicitySign[a-^n-] := HelicitySign[a]^n
170 HelicitySign[a- b-] := HelicitySign[a, b]
171 HelicitySign[-a-] := -HelicitySign[a]
172 HelicitySign[1] := 1
173 HelicitySign[-1] := -1
174 HelicitySign[a-^n-?OddQ] ^= HelicitySign[a]
175 HelicitySign[a-^n-?EvenQ] ^= 1
176
177 (* Declare spinors, conjugate spinors, slashed matrices, momenta and
      their products along with many basic simplifications that are always
      applied *)
178 Attributes[Sp] = {Flat};
179 Attributes[Metric] = {Orderless};
180 Default[Mp, 3] := Sequence[]
181 Mp[a-, b-, i-] := Mp[b, a, i] /; Order[a, b] == 1
182
183 Sp[pre---, spinorbar[p1-, hs1-], mid---, spinor[p2-, hs2-],
184   post-] := Sp[pre, post] Sp[spinorbar[p1, hs1], mid, spinor[p2, hs2]]

```

```

185 Sp[pre--, spinorbar[p1-, hs1--], mid---, spinor[p2-, hs2--],
186   post---] :=
187 Sp[pre, post] Sp[spinorbar[p1, hs1], mid, spinor[p2, hs2]]
188 Sp[pre---, a_ + b_, post---] := Sp[pre, a, post] + Sp[pre, b, post]
189 Sp[pre---, a_ b_, post---] :=
190 a Sp[pre, b, post] /;
191 FreeQ[a //.
192   Sp[spinorbar[p1-, hs1--], mid---, spinor[p2-, hs2--]] :> 1,
193   spinor | spinorbar | \[Gamma] | Sm | SmM]
194 Mp[a_ + b_, c_, i_] := Mp[a, c, i] + Mp[b, c, i]
195 Mp[a_, b_ + c_, i_] := Mp[a, b, i] + Mp[a, c, i]
196 (* Factors that contain no free objects with a space-time index can be
   moved to outside of the product *)
197 Mp[a_ b_, c_] :=
198 a Mp[b, c] /;
199 FreeQ[a //. Mp[--] :> 1, Mom | MomM | \[Gamma] | \[Sigma] | Metric]
200 Mp[a_, b_ c_] :=
201 b Mp[a, c] /;
202 FreeQ[b //. Mp[--] :> 1, Mom | MomM | \[Gamma] | \[Sigma] | Metric]
203 Mp[a_ b_, c_, i_] :=
204 a Mp[b, c, i] /;
205 FreeQ[a, Mom[-][i] | MomM[-][i] | \[Gamma][i] | \[Sigma][-][i] |
206   Metric[i, -] | Metric[-, i]]
207 Mp[a_, b_ c_, i_] :=
208 b Mp[a, c, i] /;
209 FreeQ[b,
210   Mom[-][i] | MomM[-][i] | \[Gamma][i] | \[Sigma][-][i] |
211   Metric[i, -] | Metric[-, i]]
212 Mp[a_[i-], b_[i-], i_] := Mp[a, b]
213 Mp[a_[i-], Metric[i-, j-], i_] := a[j]
214 Mp[Metric[i-, j-], a_[i-], i_] := a[j]
215 Mp[Metric[i-, j-], Metric[i-, k-], i_] := Metric[j, k]
216 Mp[Sp[pre---, \[Gamma][i-], post---], b_[i-], i_] :=
217 Mp[Sp[pre, \[Gamma], post], b]
218 Mp[a_[i-], Sp[pre---, \[Gamma][i-], post---], i_] :=
219 Mp[a, Sp[pre, \[Gamma], post]]
220 Mp[Sp[pre---, \[Gamma][i-], post---],
221   Sp[pre2---, \[Gamma][i-], post2---], i_] :=
222 Mp[Sp[pre, \[Gamma], post], Sp[pre2, \[Gamma], post2]]
223 Mp[Sp[pre---, \[Gamma][i-], post---], Metric[i-, j-], i_] :=
224 Sp[pre, \[Gamma][j], post]
225 Mp[Metric[i-, j-], Sp[pre---, \[Gamma][i-], post---], i_] :=

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226 Sp[pre, \[Gamma][j], post]
227 Mp[Sp[pre---, \[Sigma][1-][i-], post---], b-[i-], i-] :=
228 Mp[Sp[pre, \[Sigma][1], post], b]
229 Mp[a-[i-], Sp[pre---, \[Sigma][1-][i-], post---], i-] :=
230 Mp[a, Sp[pre, \[Sigma][1], post]]
231 Mp[Sp[pre---, \[Sigma][1-][i-], post---],
232 Sp[pre2---, \[Sigma][12-][i-], post2---], i-] :=
233 Mp[Sp[pre, \[Sigma][1], post], Sp[pre2, \[Sigma][12], post2]]
234 Mp[Sp[pre---, \[Sigma][1-][i-], post---], Metric[i-, j-], i-] :=
235 Sp[pre, \[Sigma][1][j], post]
236 Mp[Metric[i-, j-], Sp[pre---, \[Sigma][1-][i-], post---], i-] :=
237 Sp[pre, \[Sigma][1][j], post]
238 Mp[Mom[a-], Mom[a-]] := 0
239 Mp[0, -, i-.] := 0
240 Mp[-, 0, i-.] := 0
241 Mp2[a-] := Mp[a, a]
242
243 Sp[pre---, s1 : {---List}, s2 : {---List}, post---] :=
244 Sp[pre, s1.s2, post]
245 b_ Sp[{a---}] ^:= Sp[b {a}]
246 Sp /: Sp[{a---}] + Sp[{b---}] := Sp[{a} + {b}]
247 Sp[{a---}] == Sp[{b---}] ^:= {a} == {b}
248 Sp[{{n-}}] := n
249 Mp[p1-List, p2-List] :=
250 p1[[1]] p2[[1]] -
251 Sum[p1[[$i]] p2[[$i]], {$i, 2, Min[Length[p1], Length[p2]]}]
252
253 (* Convert the Minkowski product of momenta with a spinor chain
      containing a gamma matrix to a slashed matrix in the appropriate
      location *)
254 DoMpToSm[expr_] :=
255 expr //. {Mp[Mom[a-], Sp[pre--, \[Gamma], post---]] |
256 Mp[Sp[pre--, \[Gamma], post---], Mom[a-]] :>
257 Sp[pre, Sm[a], post],
258 Mp[Mom[a-], Sp[pre--, \[Sigma][1-], post---]] |
259 Mp[Sp[pre--, \[Sigma][1-], post---], Mom[a-]] :>
260 Sp[pre, Sm[a, 1], post],
261 Mp[MomM[a-], Sp[pre--, \[Gamma], post---]] |
262 Mp[Sp[pre--, \[Gamma], post---], MomM[a-]] :>
263 Sp[pre, SmM[a], post],
264 Mp[MomM[a-], Sp[pre--, \[Sigma][1-], post---]] |
265 Mp[Sp[pre--, \[Sigma][1-], post---], MomM[a-]] :>

```

```

266   Sp[pre, SmM[a, 1], post]}
267
268 (* Convert the Minkowski product of two spinor chains, both containing a
      gamma matrix, to products of spinor chains *)
269 DoMpToSpinorChains[expr_] :=
270   expr //.
271   {Mp[Sp[pre1_---, Sm[p_, mh_], \[Sigma][h_], post1_---],
272     Sp[pre2_---, \[Sigma][h_],
273     post2_---]] := -Mp[Sp[pre1, \[Sigma][mh], Sm[p, h], post1],
274     Sp[pre2, \[Sigma][h], post2]] +
275     2 Sp[pre1, post1] Sp[pre2, Sm[p, h], post2] /; mh == -h,
276   Mp[Sp[pre1_---, \[Sigma][h_], Sm[p_, mh_], post1_---],
277     Sp[pre2_---, \[Sigma][h_],
278     post2_---]] := -Mp[Sp[pre1, Sm[p, h], \[Sigma][mh], post1],
279     Sp[pre2, \[Sigma][h], post2]] +
280     2 Sp[pre1, post1] Sp[pre2, Sm[p, h], post2] /; mh == -h,
281   Mp[Sp[pre2_---, \[Sigma][h_], post2_---],
282     aa : Sp[pre1_---, Sm[p_, mh_], \[Sigma][h_],
283     post1_---]] := -Mp[Sp[pre1, \[Sigma][mh], Sm[p, h], post1],
284     Sp[pre2, \[Sigma][h], post2]] +
285     2 Sp[pre1, post1] Sp[pre2, Sm[p, h], post2] /; mh == -h,
286   Mp[Sp[pre2_---, \[Sigma][h_], post2_---],
287     aa : Sp[pre1_---, \[Sigma][h_], Sm[p_, mh_],
288     post1_---]] := -Mp[Sp[pre1, Sm[p, h], \[Sigma][mh], post1],
289     Sp[pre2, \[Sigma][h], post2]] +
290     2 Sp[pre1, post1] Sp[pre2, Sm[p, h], post2] /;
291   mh == -h} //.
292 {Mp[
293   Sp[spinorbar[p1_, hp1_, lp1_],
294     midpre : (Sm[-, -] ...), \[Sigma][h_], midpost : (Sm[-, -] ...),
295     spinor[p2_, hp2_, lp2_]],
296   Sp[pre_---, \[Sigma][mh_], post_---]] :=
297   2 (Sp[pre, midpost, spinor[p2, hp2, lp2], spinorbar[p1, hp1, lp1],
298     midpre, post] -
299     HelicitySign[lp1, lp2, (-1)^(Length[{midpre, midpost})] Sp[
300     pre, Sequence @@ Reverse[{midpre}], spinor[p1, hp1, lp1],
301     spinorbar[p2, hp2, lp2], Sequence @@ Reverse[{midpost}],
302     post]) /; mh == -h,
303   Mp[Sp[pre_---, \[Sigma][mh_], post_---],
304     aaaa : Sp[spinorbar[p1_, hp1_, lp1_],
305     midpre : (Sm[-, -] ...), \[Sigma][h_],
306     midpost : (Sm[-, -] ...), spinor[p2_, hp2_, lp2_]]] :=

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```

307     2 (Sp[pre, midpost, spinor[p2, hp2, lp2], spinorbar[p1, hp1, lp1],
308         midpre, post] -
309         HelicitySign[lp1, lp2, (-1)^(Length[{midpre, midpost})]) Sp[
310         pre, Sequence @@ Reverse[{midpre}], spinor[p1, hp1, lp1],
311         spinorbar[p2, hp2, lp2], Sequence @@ Reverse[{midpost}],
312         post]) /; mh == -h,
313 Mp[Sp[spinorbar[p1-, hp1-, lp1-],
314     midpre : (Sm[-, -] ...), \[Sigma][h-], post---],
315     Sp[pre---, \[Sigma][mh-], midpost : (Sm[-, -] ...),
316     spinor[p2-, hp2-, lp2-]]] :>
317     2 (Sp[spinorbar[p1, hp1, lp1], midpre, midpost,
318         spinor[p2, hp2, lp2]] Sp[pre, post] -
319         HelicitySign[lp1, lp2, (-1)^(Length[{midpre, midpost})]) Sp[
320         pre, Sequence @@ Reverse[{midpre}], spinor[p1, hp1, lp1],
321         spinorbar[p2, hp2, lp2], Sequence @@ Reverse[{midpost}],
322         post]) /; mh == -h,
323 Mp[Sp[pre---, \[Sigma][mh-], midpost : (Sm[-, -] ...),
324     spinor[p2-, hp2-, lp2-]],
325     aaaa : Sp[spinorbar[p1-, hp1-, lp1-],
326     midpre : (Sm[-, -] ...), \[Sigma][h-], post---]] :>
327     2 (Sp[spinorbar[p1, hp1, lp1], midpre, midpost,
328         spinor[p2, hp2, lp2]] Sp[pre, post] -
329         HelicitySign[lp1, lp2, (-1)^(Length[{midpre, midpost})]) Sp[
330         spinorbar[p2, hp2, lp2], Sequence @@ Reverse[{midpost}],
331         post] Sp[pre, Sequence @@ Reverse[{midpre}],
332         spinor[p1, hp1, lp1]]) /; mh == -h}
333
334 (* Declare more simplifications and rearrangements for spinor products
335    that are always applied *)
336 Sp[---, spinorbar[p-, hs1--], spinor[p-, hs2--], ---] := 0
337 Sp[---, Sm[p-, -], spinor[p-, hs--], ---] := 0
338 Sp[---, spinorbar[p-, hs--], Sm[p-, -], ---] := 0
339 Sp[---, Sm[p-], Sm[p-], ---] := 0
340 Sp[pre---, SmM[p-], SmM[p-], post---] := Sp[pre, post] Mp2[MomM[p]]
341 Sp[---, Sm[p-, -], Sm[p-, -], ---] := 0
342 Sp[pre---, SmM[p-, -], SmM[p-, -], post---] :=
343     Sp[pre, post] Mp2[MomM[p]]
344 Sp[pre---, (Sm | SmM)[p1-, h1-] | \[Sigma][h1-],
345     Sm[p2-, h2-] | \[Sigma][h2-], post---] := 0 /; h1 == h2
346 Sp[pre---, (sm : Sm | SmM)[p1-],
347     o : ((Sm | SmM)[- , h2-] | \[Sigma][h2-]), post---] :=

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348 Sp[pre, sm[p1, -h2], o, post]
349 Sp[pre---,
350   o : ((Sm | SmM)[- , h1-] | \[Sigma][h1-]), (sm : Sm | SmM)[p2-],
351   post---] := Sp[pre, o, sm[p2, -h1], post]
352 Sp[pre---, \[Gamma], o : ((Sm | SmM)[- , h2-] | \[Sigma][h2-]),
353   post---] := Sp[pre, \[Sigma][-h2], o, post]
354 Sp[pre---, o : ((Sm | SmM)[- , h1-] | \[Sigma][h1-]), \[Gamma],
355   post---] := Sp[pre, o, \[Sigma][-h1], post]
356
357 Sp[pre---, (Sm | SmM)[p2-, h2-], spinor[p-, h-, hs---], post---] :=
358   0 /; h2 == -h
359 Sp[pre---, \[Sigma][h2-], spinor[p-, h-, hs---], post---] :=
360   0 /; h2 == -h
361 Sp[pre---, spinorbar[p-, hs--], (Sm | SmM)[p2-, h2-], post---] :=
362   0 /; h2 == -{ConvertHels[hs]}[[1]]
363 Sp[pre---, spinorbar[p-, hs--], \[Sigma][h2-], post---] :=
364   0 /; h2 == -{ConvertHels[hs]}[[1]]
365 Sp[pre---, (sm : Sm | SmM)[p2-], spinor[p-, h-, hs---], post---] :=
366   Sp[pre, sm[p2, h], spinor[p, h, hs], post]
367 Sp[pre---, \[Gamma], spinor[p-, h-, hs---], post---] :=
368   Sp[pre, \[Sigma][h], spinor[p, h, hs], post]
369 Sp[pre---, spinorbar[p-, hs--], (sm : Sm | SmM)[p2-], post---] :=
370   Sp[pre, spinorbar[p, hs], sm[p2, {ConvertHels[hs]}[[1]]], post]
371 Sp[pre---, spinorbar[p-, hs--], \[Gamma], post---] :=
372   Sp[pre, spinorbar[p, hs], \[Sigma][{ConvertHels[hs]}[[1]]], post]
373
374 Sp[pre---, spinorbar[p1-, hs1---], spinor[p2-, h2-, hs2---],
375   post---] := 0 /; h2 == {ConvertHels[hs1]}[[1]]
376 Sp[pre---, spinorbar[p1-, hs1---], spinor[p2-, hs2---], post---] :=
377   0 /; Dimension[p1] != None && Dimension[p2] != None &&
378     MemberQ[Drop[{hs2}, -(Max[Dimension[p1], Dimension[p2]]/2 - 2)] -
379     Drop[{ConvertHels[
380       hs1}], -(Max[Dimension[p1], Dimension[p2]]/2 - 2)], 0]
381 Sp[pre---, spinorbar[p1-, hs1---], sms : Sm[-, -] ...,
382   spinor[p2-, hs2---], post---] :=
383   0 /; Dimension[p1] != None &&
384     Dimension[p2] != None && !
385     MemberQ[Dimension[#[[1]]] & /@ {sms}, None] &&
386     MemberQ[Drop[{hs2}, -(Max[Dimension[p1],
387       Dimension[#[[1]]] & /@ {sms}, Dimension[p2]]/2 -
388       2)] - (-1)^
389     Length[{sms}] Drop[{ConvertHels[

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390         hs1}}, -(Max[Dimension[p1], Dimension[#[[1]]] & /@ {sms},
391         Dimension[p2]]/2 - 2)], 0]
392
393 CompareHels[{}, {}] := False
394 CompareHels[{h1_, hh1_...}, {h2_, hh2_...}] :=
395 Module[{order = Order[h1, h2]},
396   If[order == 0, CompareHels[{hh1}, {hh2}], order > 0]]
397
398 CalculatedP /: Mom[CalculatedP[mom_]] := mom
399 CalculatedP /: MomM[CalculatedP[mom_]] := mom
400 CalculatedP /: Sm[CalculatedP[a_.. Mom[b_] + mom_], h_] :=
401   a Sm[b, h] + Sm[CalculatedP[mom], h]
402 CalculatedP /: Sm[CalculatedP[a_.. Mom[b_]], h_] := a Sm[b, h]
403
404 RevP /: Mom[RevP[mom_]] := -Mom[mom]
405 RevP /: MomM[RevP[mom_]] := -MomM[mom]
406 RevP /: Sm[RevP[a_], h_] := -Sm[a, h]
407 RevP /: spinor[RevP[a_], hs_..] := I spinor[a, hs]
408 RevP /: spinorbar[RevP[a_], hs_..] := I spinorbar[a, hs]
409
410 FlipHelicities[Sp[els_...]] := Sp @@ FlipHelicities[{els}]
411 FlipHelicities[{el_, rest_...}] :=
412   Join[{FlipHelicities[el]}, FlipHelicities[{rest}]]
413 FlipHelicities[{}] := {}
414 FlipHelicities[Sm[p_, i_]] := Sm[p, -i]
415 FlipHelicities[SmM[p_, i_]] := SmM[p, -i]
416 FlipHelicities[\[Sigma][i_]] := \[Sigma][-i]
417
418 Sp[spinorbar[p1_, hs1_..], pre : (Sm | SmM)[_..] ..., \[Sigma][h_],
419   post : (Sm | SmM)[_..] ..., spinor[p2_, hs2_..]] :=
420   SpFlipSign[{hs1}, {hs2}, Length[{pre, post}] + 1] Sp[
421     spinorbar[p2, hs2],
422     ConvertMidHelicities[Reverse[Sp[pre, \[Sigma][h], post]],
423       Length[{hs1}]], spinor[p1, hs1]] /;
424   CompareHels[{-Length[{pre}], hs1, {pre}, p1}, {-Length[{post}], hs2,
425     Reverse[{post}], p2}]
426 Sp[spinorbar[p1_, hs1_..], sms : (Sm | SmM)[_..] ...,
427   spinor[p2_, hs2_..]] :=
428   SpFlipSign[{hs1}, {hs2}, Length[{sms}]] Sp[spinorbar[p2, hs2],
429     ConvertMidHelicities[Reverse[Sp[sms]], Length[{hs1}]],
430     spinor[p1, hs1]] /;
431   CompareHels[{hs1, Sp[sms], p1}, {hs2,

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```

432   ConvertMidHelicities [Reverse[Sp[sms]], Length[{hs1}]], p2}]
433
434   barspinor[p_, hs_] := spinorbar[p, ConvertHels[hs]]
435
436   (* Replace slashed matrices by spinors. d_ is the number of dimensions
      the expression is in. It is also possible for the replacement to be
      in terms of a specified set of helicity labels, in which case the
      dimension is set by the number of labels needed for a spinor *)
437   ConvertSmToSpinors[d_, Sm[p_, h_]] :=
438     ConvertSmToSpinors[Sm[p, h], PadLeft[{}, d/2 - 2, 1]]
439   ConvertSmToSpinors[Sm[p_, h_], hs_List] :=
440     Sum[Sp[spinor[p, -h, Sequence @@ (hs is)],
441         barspinor[p, -h, Sequence @@ (hs is)], {is,
442         Tuples[{1, -1}, {Length[hs]}]}]]
443   ConvertSmToSpinors[d_, p_][expr_] :=
444     expr //. Sp[pre_..., spinorbar[p2_, h2_, hs_], Sm[pp : p, h_],
445         post_...] :>
446         Sp[pre, spinorbar[p2, h2, hs],
447             ConvertSmToSpinors[Sm[pp, h], {hs}], post] //.
448         Sp[pre_..., Sm[pp : p, h_], spinor[p2_, h2_, hs_], post_...] :>
449         Sp[pre, ConvertSmToSpinors[Sm[pp, h], {hs}], spinor[p2, h2, hs],
450             post] //. Sm[pp : p, h_] :> ConvertSmToSpinors[d, Sm[pp, h]]
451   ConvertSmToSpinors[p_, {h_, hs_...}][expr_] :=
452     expr //. Sm[pp : p, hh : h] :> ConvertSmToSpinors[Sm[pp, hh], {hs}]
453
454   (* Commute the given pair of slashed matrices wherever they occur in the
      expression *)
455   CommuteMatrices[m1_\[Sigma], m2p_][expr_] :=
456     expr /. {Sp[pre_..., m1, Sm[m2 : m2p, _], post_...] :>
457         2 Mom[m2] Sp[pre, post] - Sp[pre, Sm[m2], \[Gamma], post],
458         Sp[pre_..., m1, SmM[m2 : m2p, _], post_...] :>
459         2 MomM[m2] Sp[pre, post] - Sp[pre, SmM[m2], \[Gamma], post]}
460   CommuteMatrices[m1p_, m2_\[Sigma]][expr_] :=
461     expr /. {Sp[pre_..., Sm[m1 : m1p, _], m2, post_...] :>
462         2 Mom[m1] Sp[pre, post] - Sp[pre, \[Gamma], Sm[m1], post],
463         Sp[pre_..., SmM[m1 : m1p, _], m2, post_...] :>
464         2 MomM[m1] Sp[pre, post] - Sp[pre, \[Gamma], SmM[m1], post]}
465   CommuteMatrices[m1p_, m2p_][expr_] :=
466     expr /. {Sp[pre_..., Sm[m1 : m1p, a_], Sm[m2 : m2p, _], post_...] :>
467         2 Mp[Mom[m1], Mom[m2]] Sp[pre, post] -
468         Sp[pre, Sm[m2], Sm[m1], post],
469         Sp[pre_..., SmM[m1 : m1p, a_], Sm[m2 : m2p, _], post_...] :>

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470     2 Mp[MomM[m1], Mom[m2]] Sp[pre, post] -
471     Sp[pre, Sm[m2], SmM[m1], post],
472 Sp[pre---, Sm[m1 : m1p, a-], SmM[m2 : m2p, -], post---] :=>
473     2 Mp[Mom[m1], MomM[m2]] Sp[pre, post] -
474     Sp[pre, SmM[m2], Sm[m1], post],
475 Sp[pre---, SmM[m1 : m1p, a-], SmM[m2 : m2p, -], post---] :=>
476     2 Mp[MomM[m1], MomM[m2]] Sp[pre, post] -
477     Sp[pre, SmM[m2], SmM[m1], post]}
478
479 (* Commute matrices if this can cause a term to vanish by placing two
      elements for the same massless momenta next to each other *)
480 CommuteMatricesAway[expr_] :=
481 expr //. {Sp[pre---, Sm[m1-, h-], Sm[m2-, l-], Sm[m1-, h-],
482     post---] :=> 2 Mp[Mom[m2], Mom[m1]] Sp[pre, Sm[m1, h], post],
483 Sp[pre---, Sm[m1-, h-], SmM[m2-, l-], Sm[m1-, h-], post---] :=>
484     2 Mp[MomM[m2], Mom[m1]] Sp[pre, Sm[m1, h], post],
485 Sp[pre---, SmM[m1-, h-], Sm[m2-, l-], SmM[m1-, h-], post---] :=>
486     2 Mp[Mom[m2], Mom[m1]] Sp[pre, Sm[m1, h], post] -
487     Mp2[MomM[m1]] Sp[pre, SmMp[m2, h], post],
488 Sp[pre---, SmM[m1-, h-], SmM[m2-, l-], SmM[m1-, h-], post---] :=>
489     2 Mp[MomM[m2], Mom[m1]] Sp[pre, Sm[m1, h], post] -
490     Mp2[MomM[m1]] Sp[pre, SmMp[m2, h], post],
491 Sp[pre---, Sm[m1-, l-], Sm[m2-, l-], spinor[m1-, h---],
492     post---] :=> 2 Mp[Mom[m2], Mom[m1]] Sp[pre, spinor[m1, h], post],
493 Sp[pre---, Sm[m1-, l-], SmM[m2-, l-], spinor[m1-, h---],
494     post---] :=>
495     2 Mp[MomM[m2], Mom[m1]] Sp[pre, spinor[m1, h], post],
496 Sp[pre---, spinorbar[m1-, h---], Sm[m2-, l-], Sm[m1-, m-],
497     post--] :=>
498     2 Mp[Mom[m2], Mom[m1]] Sp[pre, spinorbar[m1, h], post],
499 Sp[pre---, spinorbar[m1-, h---], SmM[m2-, l-], Sm[m1-, m-],
500     post--] :=>
501     2 Mp[MomN[m2], Mom[m1]] Sp[pre, spinorbar[m1, h], post]} //. {Sp[
502     pre---, Sm[a-, h-], Sm[b-, l-], Sm[c-, m-], Sm[a-, n-],
503     post---] :=>
504     4 Mp[Mom[a], Mom[b]] Mp[Mom[c], Mom[a]] Sp[pre, post] -
505     2 Mp[Mom[a], Mom[b]] Sp[pre, Sm[a], Sm[c], post] -
506     2 Mp[Mom[a], Mom[c]] Sp[pre, Sm[b], Sm[a], post],
507 Sp[pre---, Sm[a-, h-], SmM[b-, l-], Sm[c-, m-], Sm[a-, n-],
508     post---] :=>
509     4 Mp[Mom[a], MomM[b]] Mp[Mom[c], Mom[a]] Sp[pre, post] -
510     2 Mp[Mom[a], MomM[b]] Sp[pre, Sm[a], Sm[c], post] -

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```

511      2 Mp[Mom[a], Mom[c]] Sp[pre, SmM[b], Sm[a], post],
512 Sp[pre---, Sm[a-, h-], Sm[b-, l-], SmM[c-, m-], Sm[a-, n-],
513   post---] :=>
514      4 Mp[Mom[a], Mom[b]] Mp[MomM[c], Mom[a]] Sp[pre, post] -
515      2 Mp[Mom[a], Mom[b]] Sp[pre, Sm[a], SmM[c], post] -
516      2 Mp[Mom[a], MomM[c]] Sp[pre, Sm[b], Sm[a], post],
517 Sp[pre---, Sm[a-, h-], SmM[b-, l-], SmM[c-, m-], Sm[a-, n-],
518   post---] :=>
519      4 Mp[Mom[a], MomM[b]] Mp[MomM[c], Mom[a]] Sp[pre, post] -
520      2 Mp[Mom[a], MomM[b]] Sp[pre, Sm[a], SmM[c], post] -
521      2 Mp[Mom[a], MomM[c]] Sp[pre, SmM[b], Sm[a], post],
522 Sp[pre---, SmM[a-, h-], Sm[b-, l-], Sm[c-, m-], SmM[a-, n-],
523   post---] :=>
524      4 Mp[Mom[a], Mom[b]] Mp[Mom[c], Mom[a]] Sp[pre, post] -
525      2 Mp[Mom[a], Mom[b]] Sp[pre, Sm[a], Sm[c], post] -
526      2 Mp[Mom[a], Mom[c]] Sp[pre, Sm[b], Sm[a], post] +
527      Mp2[MomM[a]] Sp[pre, Sm[b], Sm[c], post],
528 Sp[pre---, SmM[a-, h-], SmM[b-, l-], Sm[c-, m-], SmM[a-, n-],
529   post---] :=>
530      4 Mp[Mom[a], MomM[b]] Mp[Mom[c], Mom[a]] Sp[pre, post] -
531      2 Mp[Mom[a], MomM[b]] Sp[pre, Sm[a], Sm[c], post] -
532      2 Mp[Mom[a], Mom[c]] Sp[pre, SmM[b], Sm[a], post] +
533      Mp2[MomM[a]] Sp[pre, SmM[b], Sm[c], post],
534 Sp[pre---, SmM[a-, h-], Sm[b-, l-], SmM[c-, m-], SmM[a-, n-],
535   post---] :=>
536      4 Mp[Mom[a], Mom[b]] Mp[MomM[c], Mom[a]] Sp[pre, post] -
537      2 Mp[Mom[a], Mom[b]] Sp[pre, Sm[a], SmM[c], post] -
538      2 Mp[Mom[a], MomM[c]] Sp[pre, Sm[b], Sm[a], post] +
539      Mp2[MomM[a]] Sp[pre, Sm[b], SmM[c], post],
540 Sp[pre---, SmM[a-, h-], SmM[b-, l-], SmM[c-, m-], SmM[a-, n-],
541   post---] :=>
542      4 Mp[Mom[a], MomM[b]] Mp[MomM[c], Mom[a]] Sp[pre, post] -
543      2 Mp[Mom[a], MomM[b]] Sp[pre, Sm[a], SmM[c], post] -
544      2 Mp[Mom[a], MomM[c]] Sp[pre, SmM[b], Sm[a], post] +
545      Mp2[MomM[a]] Sp[pre, SmM[b], SmM[a], post]}
546
547 (* Commute gamma matrices if it can cause a term to vanish by placing
548    two elements for the same massless momenta next to each other *)
548 CommuteSigmaMatriciesAway[expr_] :=
549 expr //. {Sp[pre---, Sm[m1-, h-], \[Sigma][l-], Sm[m1-, h-],
550   post---] :=> 2 Mom[m1] Sp[pre, Sm[m1, h], post],
551 Sp[pre---, SmM[m1-, h-], \[Sigma][l-], SmM[m1-, h-], post---] :=>

```

```

552     2 Mom[m1] Sp[pre, Sm[m1, h], post] -
553     Mp2[MomM[m1]] Sp[pre, \[Sigma][h], post],
554     Sp[pre_---, Sm[m1_, l_], \[Sigma][l_], spinor[m1_, h_---],
555         post_---] :> 2 Mom[m1] Sp[pre, spinor[m1, h], post],
556     Sp[pre_---, spinorbar[m1_, h_---], \[Sigma][l_], Sm[m1_, m_],
557         post_---] :> 2 Mom[m1] Sp[pre, spinorbar[m1, h], post]}
558
559 (* Dimension independant code for evaluating expressions numerically *)
560 SpinorExpression[d_, p_List, hs_---] :=
561 SpinorExpression[d, PadRight[p, d, 0], hs] /; Length[p] != d
562 SpinorExpression[d_, p_List, hs_---] :=
563 SpinorExpression[d, PadRight[p, d, 0], hs] /; Length[p] != d
564 AdjointSpinor[d_, s_, hs_--] :=
565 AdjointSign[d, hs] Transpose[s].AdjointMetric[d]
566
567 \[Gamma]Expression[2, -, 0] := {{1}}
568 \[Gamma]Expression[2, h_, -] := {{h}}
569 \[Gamma]Expression[d_, h_, i_] := ArrayFlatten[{{
570     \[Gamma]Expression[d - 2, h, i], 0}, {0, -
571     \[Gamma]Expression[d - 2, -h, i]}}] /;
572     d > 2 && EvenQ[d] && i < d - 2
573 \[Gamma]Expression[d_, s_, i_] :=
574 ArrayFlatten[{{0,
575     I IdentityMatrix[2^(d/2 - 2)]}, {I IdentityMatrix[2^(d/2 - 2)],
576     0}}] /; d > 2 && EvenQ[d] && i == d - 2
577 \[Gamma]Expression[d_, s_, i_] :=
578 ArrayFlatten[{{0,
579     IdentityMatrix[2^(d/2 - 2)]}, {-IdentityMatrix[2^(d/2 - 2)],
580     0}}] /; d > 2 && EvenQ[d] && i == d - 1
581
582 shift$$reps = {};
583 ReplaceShiftedSpinors[expr_] := expr //. shift$$reps
584
585 vectors$$momentums$$defined[-, -] := False
586 vectors$$spinors[tag_, p_] :=
587 Module[{spinors$},
588     spinors$[d_,
589         hs_--] := (spinors$[d, hs] =
590         SpinorExpression[d, vectors$$momentums[tag, p], hs]);
591     tag /: vectors$$spinors[tag, p] = spinors$ /;
592     vectors$$momentums$$defined[tag, p]
593

```

```

594 vectors$$spinors[tag_, p : CalculatedP[pexpr_]] :=
595 Module[{spinors$},
596   spinors$[d_,
597     hs_ ] := (spinors$[d, hs] =
598     SpinorExpression[d, pexpr // Ev[d, tag], hs]);
599   tag /: vectors$$spinors[tag, p] = spinors$
600
601 vectors$$momentums[tag_,
602   shift[{p_, q_}, z_, {h_}, {l_}][
603     p_] := (Mom[p] + z Sp[spinorbar[p, h], \[Gamma], spinor[q, l]]/2 //
604     Ev[4, tag]) /;
605   h == -1 && MatchQ[vectors$$momentums[tag, p], _List] &&
606     MatchQ[vectors$$momentums[tag, q], _List]
607 vectors$$momentums[tag_,
608   shift[{p_, q_}, z_, {h_}, {l_}][
609     q_] := (Mom[q] - z Sp[spinorbar[p, h], \[Gamma], spinor[q, l]]/2 //
610     Ev[4, tag]) /;
611   h == -1 && MatchQ[vectors$$momentums[tag, p], _List] &&
612     MatchQ[vectors$$momentums[tag, q], _List]
613 shift$$reps =
614   Join[shift$$reps, {Mom[shift[{p_, q_}, z_, {h_}, {l_}][p_] ]>
615     Mom[p] + z Sp[spinorbar[p, h], \[Gamma], spinor[q, l]]/2,
616     Mom[shift[{p_, q_}, z_, {h_}, {l_}][q_] ]>
617     Mom[q] - z Sp[spinorbar[p, h], \[Gamma], spinor[q, l]]/2,
618     Sm[shift[{p_, q_}, z_, {h_}, {l_}][p_], h_] ]>
619     Sm[p, h] + z Sp[spinor[q, l], spinorbar[p, h]],
620     Sm[shift[{p_, q_}, z_, {h_}, {l_}][p_], l_] ]>
621     Sm[p, l] + z Sp[spinor[p, h], spinorbar[q, l]],
622     Sm[shift[{p_, q_}, z_, {h_}, {l_}][q_], h_] ]>
623     Sm[q, h] - z Sp[spinor[q, l], spinorbar[p, h]],
624     Sm[shift[{p_, q_}, z_, {h_}, {l_}][q_], l_] ]>
625     Sm[q, l] - z Sp[spinor[p, h], spinorbar[q, l]]}];
626 vectors$$spinors[tag_, shift[{p_, q_}, z_, {h_}, {l_}][p_][4,
627   a_] := (spinor[p, a] + z spinor[q, a] // Ev[4, tag]) /;
628   h == -1 && a == 1 && MatchQ[vectors$$momentums[tag, p], _List] &&
629     MatchQ[vectors$$momentums[tag, q], _List]
630 vectors$$spinors[tag_, shift[{p_, q_}, z_, {h_}, {l_}][p_][4, a_] :=
631   (spinor[p, a] // Ev[4, tag]) /;
632   h == -1 && a == -1 && MatchQ[vectors$$momentums[tag, p], _List] &&
633     MatchQ[vectors$$momentums[tag, q], _List]
634 vectors$$spinors[tag_, shift[{p_, q_}, z_, {h_}, {l_}][q_][4,
635   a_] := (spinor[q, a] - z spinor[p, a] // Ev[4, tag]) /;

```

```

636 h == -1 && a == h && MatchQ[vectors$$momentums[tag, p], _List] &&
637 MatchQ[vectors$$momentums[tag, q], _List]
638 vectors$$spinors[tag_, shift[{p_, q_}, z_, {h_}, {l_}][q_][4,
639 a_] := (spinor[q, a] // Ev[4, tag]) /;
640 h == -1 && a == -h && MatchQ[vectors$$momentums[tag, p], _List] &&
641 MatchQ[vectors$$momentums[tag, q], _List]
642 shift$$reps =
643 Join[shift$$reps, {(ss : spinor | spinorbar)[
644 shift[{p_, q_}, z_, {h_}, {l_}][p_, h_] :=>
645 ss[p, h], (ss : spinor | spinorbar)[
646 shift[{p_, q_}, z_, {h_}, {l_}][p_, l_] :=>
647 ss[p, l] + z ss[q, l], (ss : spinor | spinorbar)[
648 shift[{p_, q_}, z_, {h_}, {l_}][q_, l_] :=>
649 ss[q, l], (ss : spinor | spinorbar)[
650 shift[{p_, q_}, z_, {h_}, {l_}][q_, h_] :=>
651 ss[q, h] - z ss[p, h]}}];
652
653 vectors$$momentums[tag_,
654 shift[{p_, q_}, z_, {h_, l_}, {m_, n_}][
655 p_] := (Mom[p] -
656 HelicitySign[h,
657 n] z Sp[spinorbar[p, h, l], \[Gamma], Sm[q], spinor[p, m, n]]/
658 2 // Ev[6, tag]) /;
659 h == -m && MatchQ[vectors$$momentums[tag, p], _List] &&
660 MatchQ[vectors$$momentums[tag, q], _List]
661 vectors$$momentums[tag_,
662 shift[{p_, q_}, z_, {h_, l_}, {m_, n_}][
663 q_] := (Mom[q] +
664 HelicitySign[h,
665 n] z Sp[spinorbar[p, h, l], \[Gamma], Sm[q], spinor[p, m, n]]/
666 2 // Ev[6, tag]) /;
667 h == -m && MatchQ[vectors$$momentums[tag, p], _List] &&
668 MatchQ[vectors$$momentums[tag, q], _List]
669 shift$$reps =
670 Join[shift$$reps, {Mom[
671 shift[{p_, q_}, z_, {h_, l_}, {m_, n_}][p_] :=>
672 Mom[p] -
673 HelicitySign[h,
674 n] z Sp[spinorbar[p, h, l], \[Gamma], Sm[q], spinor[p, m, n]]/
675 2, Mom[shift[{p_, q_}, z_, {h_, l_}, {m_, n_}][q_] :=>
676 Mom[q] +
677 HelicitySign[h,

```



```

678      n] z Sp[spinorbar[p, h, l], \[Gamma], Sm[q], spinor[p, m, n]]/
679      2,
680      Sm[shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][p-], h-] :>
681      Sm[p, h] +
682      z (HelicitySign[h, l] Sp[Sm[q], spinor[p, h, l],
683      spinorbar[p, m, n]] -
684      HelicitySign[m, n] Sp[spinor[p, m, n], spinorbar[p, h, l],
685      Sm[q]]),
686      Sm[shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][p-], m-] :>
687      Sm[p, m] +
688      z (-HelicitySign[h, l] Sp[spinor[p, h, l], spinorbar[p, m, n],
689      Sm[q]] +
690      HelicitySign[m, n] Sp[Sm[q], spinor[p, m, n],
691      spinorbar[p, h, l]]),
692      Sm[shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][q-], h-] :>
693      Sm[q, h] -
694      z (HelicitySign[h, l] Sp[Sm[q], spinor[p, h, l],
695      spinorbar[p, m, n]] -
696      HelicitySign[m, n] Sp[spinor[p, m, n], spinorbar[p, h, l],
697      Sm[q]]),
698      Sm[shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][q-], m-] :>
699      Sm[q, m] -
700      z (-HelicitySign[h, l] Sp[spinor[p, h, l], spinorbar[p, m, n],
701      Sm[q]] +
702      HelicitySign[m, n] Sp[Sm[q], spinor[p, m, n],
703      spinorbar[p, h, l]]);
704      vectors$$$spinors[tag-, shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][p-]][6,
705      a-, b-] := (spinor[p, a, b] -
706      z HelicitySign[a, n] Sm[q, m].spinor[p, m, n] // Ev[6, tag]) /;
707      m == -h && h == a && l == -b &&
708      MatchQ[vectors$$$momentums[tag, p], _List] &&
709      MatchQ[vectors$$$momentums[tag, q], _List]
710      vectors$$$spinors[tag-, shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][p-]][6,
711      a-, b-] := (spinor[p, a, b] -
712      z HelicitySign[a, l] Sm[q, h].spinor[p, h, l] // Ev[6, tag]) /;
713      m == -h && m == a && n == -b &&
714      MatchQ[vectors$$$momentums[tag, p], _List] &&
715      MatchQ[vectors$$$momentums[tag, q], _List]
716      vectors$$$spinors[tag-, shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][p-]][6,
717      a-, b-] := (spinor[p, a, b] // Ev[6, tag]) /;
718      m == -h && ((h == a && l == b) || (m == a && n == b)) &&
719      MatchQ[vectors$$$momentums[tag, p], _List] &&

```

```

720 MatchQ[vectors$$momentums[tag, q], _List]
721 vectors$$spinors[tag-, shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][q-]][6,
722 a-, b-] := (spinor[q, a, b] +
723 z HelicitySign[h, l] spinor[p, h, l] Sp[spinorbar[p, m, n],
724 spinor[q, a, b]] // Ev[6, tag]) /;
725 m == -h && h == a && MatchQ[vectors$$momentums[tag, p], _List] &&
726 MatchQ[vectors$$momentums[tag, q], _List]
727 vectors$$spinors[tag-, shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][q-]][6,
728 a-, b-] := (spinor[q, a, b] +
729 z HelicitySign[m, n] spinor[p, m, n] Sp[spinorbar[p, h, l],
730 spinor[q, a, b]] // Ev[6, tag]) /;
731 m == -h && m == a && MatchQ[vectors$$momentums[tag, p], _List] &&
732 MatchQ[vectors$$momentums[tag, q], _List]
733 shift$$reps =
734 Join[shift$$reps, {(ss : spinor | spinorbar)[
735 shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][p-], h-, l-] :=
736 ss[p, h, l], (ss : spinor | spinorbar)[
737 shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][p-], m-, n-] :=
738 ss[p, m, n],
739 spinor[shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][p-], h-, b-] :=
740 spinor[p, h, b] -
741 z HelicitySign[h, n] Sp[Sm[q, m], spinor[p, m, n]] /; b == -l,
742 spinor[shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][p-], m-, b-] :=
743 spinor[p, m, b] -
744 z HelicitySign[m, l] Sp[Sm[q, h], spinor[p, h, l]] /; b == -n,
745 spinorbar[shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][p-], h-, b-] :=
746 spinorbar[p, h, b] -
747 z HelicitySign[h, l] Sp[spinorbar[p, m, n], Sm[q, m]] /;
748 b == -l,
749 spinorbar[shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][p-], m-, b-] :=
750 spinorbar[p, m, b] -
751 z HelicitySign[m, n] Sp[spinorbar[p, h, l], Sm[q, h]] /;
752 b == -n, (ss : spinor | spinorbar)[
753 shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][q-], h-, b-] :=
754 ss[q, h, b] +
755 z If[TrueQ[ss == spinorbar], HelicitySign[h, b],
756 HelicitySign[h, l]] ss[p, h, l] Sp[spinorbar[p, m, n],
757 spinor[q, h, b]], (ss : spinor | spinorbar)[
758 shift[{p-, q-}, z-, {h-, l-}, {m-, n-}][q-], m-, b-] :=
759 ss[q, m, b] +
760 z If[TrueQ[ss == spinorbar], HelicitySign[m, b],
761 HelicitySign[m, n]] ss[p, m, n] Sp[spinorbar[p, h, l],

```

```

762     spinor[q, m, b]]]);
763
764 (* Function to declare momenta for names on the momenta set 'tag' *)
765 DeclareVectorMomentum[tag_, p_,
766   pp_List] := (tag /: vectors$$momentums$$defined[tag, p] = True;
767   tag /: vectors$$momentums[tag, p] = pp)
768 $CheckSpinorsConsistent = True;
769 SetAttributes[CheckSpinorsMomentum, HoldFirst]
770 CheckSpinorsMomentum::Inconsistent =
771   "The spinors '3' and '4' give a momentum of '5' which is not the \
772   same as the momenta '6' for '1' '2'.";
773 CheckSpinorsMomentum[tag_,
774   p_, {hs : (1 | -1) ...} -> {spp : {{-} ..}, spm : {{-} ..}}] :=
775   Message[CheckSpinorsMomentum::Inconsistent, tag, p, spp, spm,
776     CalculateMomenta[spp, spm, {hs}],
777     vectors$$momentums[tag, p]] /; $CheckSpinorsConsistent &&
778   Simplify[
779     CalculateMomenta[spp, spm, {hs}] != vectors$$momentums[tag, p]]
780 DefineSpinors[
781   spinors$, {hs : (1 | -1) ...} -> {spp : {{-} ..},
782     spm : {{-} ..}}] := (spinors$[-, hs, 1] = spp;
783   spinors$[-, hs, -1] = spm)
784 DeclareVectorMomentum[tag_, p_,
785   sp : ({hs : (1 | -1) ...} -> {spp : {{-} ..}, spm : {{-} ..})},
786   sps : ({(1 | -1) ...} -> {{{-} ..}, {{-} ..}) ...] := (tag /:
787     vectors$$momentums[tag,
788       p] := (tag /: vectors$$momentums[tag, p] =
789         CalculateMomenta[spp, spm, {hs}]);
790   CheckSpinorsMomentum[tag, p, #] & /@ {sp, sps};
791   Module[{spinors$}, DefineSpinors[spinors$, #] & /@ {sp, sps};
792     tag /: vectors$$spinors[tag, p] = spinors$; spinors$)
793 DeclareVectorMomentum[tag_, p_, spspre : ({(1 | -1) ...} -> -) ...,
794   sph : {(1 | -1) ...} -> {sp.Sp, other-} ,
795   spspost : ({(1 | -1) ...} -> -) ...] :=
796   DeclareVectorMomentum[tag, p, spspre, sph -> {other, sp[[1]]},
797     spspost] /; Length[sp] == 1
798 DeclareVectorMomentum[tag_, p_, spspre : ({(1 | -1) ...} -> -) ...,
799   sph : {(1 | -1) ...} -> {other-, sp.Sp} ,
800   spspost : ({(1 | -1) ...} -> -) ...] :=
801   DeclareVectorMomentum[tag, p, spspre, sph -> {sp[[1]], other},
802     spspost] /; Length[sp] == 1
803 DeclareVectorMomentum[tag_, p_,

```

```

804 sp : {{{(1 | -1) ...} -> ({{-} ..} | -Sp)) ..} :=
805 Module[{sp$},
806   DeclareVectorMomentum[tag, p,
807     Sequence @@ (#[[1,
808       1, ;; -2]] -> ({sp$[1],
809       sp$[-1]} // ((sp$[#[[1, -1]]] -> #[[2]]) & /@ #)) &) /@
810     GatherBy[sp, #[[1, ;; -2]] &]]]
811
812 (* Remove the declaration of a momenta for a specified 'tag' *)
813 UndeclareVectorMomentum[tag_,
814   p_] := (Remove[Evaluate[vectors$$spinors[tag, p]]];
815   tag /: vectors$$spinors[tag, p] =.;
816   tag /: vectors$$momentums[tag, p] =.;
817   tag /: vectors$$momentums$$defined[tag, p] =.);
818
819 (* Declare that a 'tag' inherits all momenta values defined for another
      tag *)
820 DeclareInheritingTag[tag_,
821   parent_] := (tag /: vectors$$momentums[tag, p_] :=
822     vectors$$momentums[parent, p];
823   tag /: vectors$$spinors[tag, p_] := vectors$$spinors[parent, p];
824   tag /: vectors$$momentums[tag, p_] :=
825     vectors$$momentums$$defined[parent, p];)
826
827 (* Evaluate expressions numerically *)
828 Ev[d_, moms_][expr_] :=
829   Module[{e = expr},
830     e //.{spinor[p_, hs_Integer] :>
831       vectors$$spinors[moms, p][d, hs] /;
832       MatchQ[vectors$$spinors[moms, p][d, hs], {--List}],
833     spinorbar[p_, hs_Integer] :>
834       AdjointSpinor[d, vectors$$spinors[moms, p][d, hs], hs] /;
835       MatchQ[vectors$$spinors[moms, p][d, hs], {--List}], (Sm | SmM)[
836       p_, h_] :> \[Gamma]Expression[d, h,
837       0] vectors$$momentums[moms, p][[1]] -
838       Sum[\[Gamma]Expression[d,
839         h, $$i] vectors$$momentums[moms, p][[$$i + 1]], {$$i, 1,
840         Length[vectors$$momentums[moms, p]] - 1}] /;
841       MatchQ[vectors$$momentums[moms, p], _List], (Mom | MomM)[p_] :>
842       PadRight[vectors$$momentums[moms, p], d, 0] /;
843       MatchQ[vectors$$momentums[moms, p], _List],
844     Sp[pre_, \[Sigma][h_], post_] :>

```

```

845      Table[Sp[pre, \[Gamma]Expression[d, h, \[Mu]], post], {\[Mu], 0,
846          d - 1}]]]
847
848 (* Implementations for functions that depend on the number of dimensions
      *)
849 ConvertHels[i_] := Sequence[-i] (* 4d *)
850
851 ConvertHels[i_, j_] := Sequence[i, -j] (* 6d *)
852
853 ConvertMidHelicities[sp_, 1] := FlipHelicities[sp]
854 ConvertMidHelicities[sp_, 2] := sp
855
856 SpFlipSign[{-}, {-}, n_] := HelicitySign[(-1)^(n + 1)] (* 4d *)
857
858 SpFlipSign[{-, h1_}, {-, h2_}, n_] :=
859 HelicitySign[(-1)^n, h1, h2] (* 6d *)
860
861 Sp[pre___, spinorbar[p_, hh_, h_], Sm[___], spinor[p_, hh_, h_],
862   post___] := 0
863 Sp[pre___, spinorbar[p_, hh_, h_], \[Sigma][_], spinor[p_, hh_, h_],
864   post___] := 0
865
866 SpinorExpression[4, {p0_, p1_, p2_, p3_}, 1] :=
867 If[TrueQ[Simplify[p0 + p1 == 0]],
868   If[TrueQ[p0 - p1 == 0],
869     Module[{sqrtp3 =
870       Sqrt[2 p3]}, {{(p3 + I p2)/Sqrt[2 p3]}, {(p3 - I p2)/
871       Sqrt[2 p3]}}],
872     Module[{sqrtpm =
873       Sqrt[p0 - p1]}, {{(p3 + I p2)/sqrtpm}, {sqrtpm}}]],
874   Module[{sqrtpp = Sqrt[p0 + p1]}, {{sqrtpp}, {(p3 - I p2)/sqrtpp}}]]]
875 SpinorExpression[4, {p0_, p1_, p2_, p3_}, -1] :=
876 If[TrueQ[Simplify[p0 + p1 == 0]],
877   If[TrueQ[p0 - p1 == 0],
878     Module[{sqrtp3 =
879       Sqrt[2 p3]}, {{(p3 + I p2)/Sqrt[2 p3]}, {(p3 - I p2)/
880       Sqrt[2 p3]}}],
881     Module[{sqrtpm =
882       Sqrt[p0 - p1]}, {sqrtpm}, {(p3 - I p2)/sqrtpm}}]],
883   Module[{sqrtpp = Sqrt[p0 + p1]}, {{(p3 + I p2)/sqrtpp}, {sqrtpp}}]]]
884 AdjointMetric[4] = {{0, -1}, {1, 0}};
885 AdjointSign[4, i_] := 1

```

```

886
887 CalculateMomenta[{{ppp-}, {ppmp-}}, {{pppm-}, {ppm-}}, {}] := {(ppm \
888 ppp + ppmp ppm)/2, (ppm ppp - ppmp ppm)/2,
889 I (ppm ppmp - ppp ppm)/2, (ppm ppmp + ppp ppm)/2}
890
891 SpinorExpression[6, {p0-, p1-, p2-, p3-, p4-, p5-}, 1, 1] :=
892 If[TrueQ[Simplify[p0 + p1 == 0]],
893 If[TrueQ[Simplify[p0 - p1 == 0]],
894 If[TrueQ[Simplify[p3 + I p2 == 0]],
895 If[TrueQ[
896 Simplify[
897 p3 - I p2 == 0]], {{(p5 + I p4)/
898 Sqrt[2 p5]}, {0}, {0}, {(p5 - I p4)/Sqrt[2 p5]}}],
899 Module[{sqrtppm =
900 Sqrt[p3 - I p2]}, {{sqrtppm}, {0}, {0}, {-(p5 - I p4)/
901 sqrtppm}}]],
902 Module[{sqrtppp =
903 Sqrt[p3 + I p2]}, {{sqrtppp}, {0}, {0}, {-(p5 - I p4)/
904 sqrtppp}}]],
905 Module[{sqrtpm =
906 Sqrt[p0 - p1]}, {{(p3 + I p2)/
907 sqrtpm}, {sqrtpm}, {0}, {-(p5 - I p4)/sqrtpm}}]],
908 Module[{sqrtpp =
909 Sqrt[p0 + p1]}, {{sqrtpp}, {(p3 - I p2)/sqrtpp}, {(p5 - I p4)/
910 sqrtpp}, {0}}]]
911 SpinorExpression[6, {p0-, p1-, p2-, p3-, p4-, p5-}, 1, -1] :=
912 If[TrueQ[Simplify[p0 + p1 == 0]],
913 If[TrueQ[Simplify[p0 - p1 == 0]],
914 If[TrueQ[Simplify[p3 + I p2 == 0]],
915 If[TrueQ[
916 Simplify[
917 p3 - I p2 == 0]], {{0}, {-(p5 + I p4)/
918 Sqrt[2 p5]}, {-(p5 - I p4)/Sqrt[2 p5]}, {0}},
919 Module[{sqrtppm =
920 Sqrt[p3 - I p2]}, {{0}, {-(p5 + I p4)/
921 sqrtppm}, {sqrtppm}, {0}}]],
922 Module[{sqrtppp =
923 Sqrt[p3 + I p2]}, {{0}, {-(p5 + I p4)/
924 sqrtppp}, {sqrtppp}, {0}}]],
925 Module[{sqrtpm =
926 Sqrt[p0 - p1]}, {{(p5 + I p4)/
927 sqrtpm}, {0}, {sqrtpm}, {(p3 - I p2)/sqrtpm}}]],

```

```

928 Module[{sqrtpp =
929   Sqrt[p0 + p1]}, {{0}, {-(p5 + I p4)/sqrtpp}, {(p3 + I p2)/
930   sqrtpp}, {sqrtpp}}]]
931 SpinorExpression[6, {p0-, p1-, p2-, p3-, p4-, p5-}, -1, 1] :=
932 If[TrueQ[Simplify[p0 + p1 == 0]],
933 If[TrueQ[Simplify[p0 - p1 == 0]],
934 If[TrueQ[Simplify[p3 + I p2 == 0]],
935 If[TrueQ[
936   Simplify[
937     p3 - I p2 == 0]], {{0}, {-(p5 + I p4)/
938     Sqrt[2 p5]}, {-(p5 - I p4)/Sqrt[2 p5]}, {0}},
939 Module[{sqrtppm =
940   Sqrt[p3 - I p2]}, {{0}, {-(p5 + I p4)/
941   sqrtppm}, {sqrtppm}, {0}}]],
942 Module[{sqrtppp =
943   Sqrt[p3 + I p2]}, {{0}, {-(p5 + I p4)/
944   sqrtppp}, {sqrtppp}, {0}}]],
945 Module[{sqrtpm =
946   Sqrt[p0 - p1]}, {{0}, {-(p5 + I p4)/sqrtpm}, {(p3 + I p2)/
947   sqrtpm}, {sqrtpm}}]],
948 Module[{sqrtpp =
949   Sqrt[p0 + p1]}, {{(p5 + I p4)/
950   sqrtpp}, {0}, {sqrtpp}, {(p3 - I p2)/sqrtpp}}]]
951 SpinorExpression[6, {p0-, p1-, p2-, p3-, p4-, p5-}, -1, -1] :=
952 If[TrueQ[Simplify[p0 + p1 == 0]],
953 If[TrueQ[Simplify[p0 - p1 == 0]],
954 If[TrueQ[Simplify[p3 + I p2 == 0]],
955 If[TrueQ[
956   Simplify[
957     p3 - I p2 == 0]], {{(p5 + I p4)/
958     Sqrt[2 p5]}, {0}, {0}, {(p5 - I p4)/Sqrt[2 p5]}},
959 Module[{sqrtppm =
960   Sqrt[p3 - I p2]}, {{sqrtppm}, {0}, {0}, {-(p5 - I p4)/
961   sqrtppm}}]],
962 Module[{sqrtppp =
963   Sqrt[p3 + I p2]}, {{sqrtppp}, {0}, {0}, {-(p5 - I p4)/
964   sqrtppp}}]],
965 Module[{sqrtpm =
966   Sqrt[p0 - p1]}, {{sqrtpm}, {(p3 - I p2)/sqrtpm}, {(p5 - I p4)/
967   sqrtpm}, {0}}]],
968 Module[{sqrtpp =
969   Sqrt[p0 + p1]}, {{(p3 + I p2)/

```

```

970      sqrtpp}, {sqrtpp}, {0}, {-(p5 - I p4)/sqrtpp}}]]
971 AdjointMetric [
972   6] = {{0, 0, 0, 1}, {0, 0, -1, 0}, {0, 1, 0, 0}, {-1, 0, 0, 0}};
973 AdjointSign[6, i_, j_] := HelicitySign[i j]
974
975 CalculateMomenta6dI[pp_, pm_, ppp_, ppm_, pdp_, pdm_,
976  h_] := {pm + pp, pp - pm, I (ppm - ppp), ppm + ppp, h I (pdm - pdp),
977  pdm + pdp}/2
978 CalculateMomenta[{{sp1_}, {sp2_}, {sp3_}, {sp4_}}, {{sm1_}, {sm2_}, \
979 {sm3_}, {sm4_}}, {h : 1}] :=
980 CalculateMomenta6dI[sm4 sp1 - sm1 sp4, sm3 sp2 - sm2 sp3,
981  sm3 sp1 - sm1 sp3, sm4 sp2 - sm2 sp4, -sm2 sp1 + sm1 sp2,
982  sm4 sp3 - sm3 sp4, h]
983 CalculateMomenta[{{sp3_}, {sp4_}, {sp1_}, {sp2_}}, {{sm3_}, {sm4_}, \
984 {sm1_}, {sm2_}}, {h : -1}] :=
985 CalculateMomenta6dI[sm4 sp1 - sm1 sp4, sm3 sp2 - sm2 sp3,
986  sm3 sp1 - sm1 sp3, sm4 sp2 - sm2 sp4, -sm2 sp1 + sm1 sp2,
987  sm4 sp3 - sm3 sp4, h]
988
989 (* The polarisation vector in any number of dimensions. The number of
    dimensions is set by the number of signs given as each of i_ and
    j_ *)
990 PolVec[p_, q_, i_, j_] :=
991 Sp[spinorbar[p, i], \[Gamma], Sm[q], spinor[p, j]]/2^(3/2)/
992 Mp[Mom[p], Mom[q]] /; Length[{i}] == Length[{j}]
993 PolVec[p_, q_, i_, j_, \[Mu]_] :=
994 Sp[spinorbar[p, i], \[Gamma][\[Mu]], Sm[q], spinor[p, j]]/2^(3/2)/
995 Mp[Mom[p], Mom[q]] /; Length[{i}] == Length[{j}]
996
997 (* Replacements for reducing 6 dimensional momenta to 4 dimensional
    momenta and the reverse *)
998 rep = {spinor[pp : p | q | n6p | n6m | shift[{p, q}, z][p | q], h_,
999  l_] :=> spinor[pp[a], h, l] -
1000  HelicitySign[h,
1001  l] (pp[5] -
1002  I HelicitySign[h, l] pp[4]) Sp[Sm[a, -h],
1003  spinor[pp[a], -h, l]]/2/Mp[Mom[a], Mom[pp[a]]],
1004  spinorbar[pp : p | q | n6p | n6m | shift[{p, q}, z][p | q], h_,
1005  l_] :=> spinorbar[pp[a], h, l] +
1006  HelicitySign[h,
1007  l] (pp[5] -
1008  I HelicitySign[h, l] pp[4]) Sp[spinorbar[pp[a], -h, l],

```



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1009      Sm[a, -h]]/2/Mp[Mom[a], Mom[pp[a]]],
1010  Mom[pp : p | q | n6p | n6m | shift[{p, q}, z][p | q]] :>
1011  Mom[pp[a]] + Mom[a] (pp[4]^2 + pp[5]^2)/2/Mp[Mom[pp[a]], Mom[a]] +
1012  Sum[h Sp[spinorbar[pp[a], 1, h], \[Sigma][1], Sm[a, -1],
1013      spinor[pp[a], -1, -h]] (pp[5] + I h pp[4]), {h, {1, -1}}]/4/
1014  Mp[Mom[pp[a]], Mom[a]],
1015  Sm[pp : p | q | n6p | n6m | shift[{p, q}, z][p | q], h_] :>
1016  Sp[Sm[pp[a], h]] +
1017  Sum[l HelicitySign [
1018      h] (pp[5] -
1019      I l HelicitySign[h] pp[4]) (Sp[spinor[pp[a], -h, 1],
1020      spinorbar[pp[a], h, -1], Sm[a, h]] -
1021      Sp[Sm[a, h], spinor[pp[a], h, -1],
1022      spinorbar[pp[a], -h, 1]] ), {l, {1, -1}}]/2/
1023  Mp[Mom[pp[a]], Mom[a]] + (pp[4]^2 + pp[5]^2) Sp[Sm[a, h]]/2/
1024  Mp[Mom[pp[a]], Mom[a]]];
1025  reph[l1_] := {spinor[pp : p | q | n6p | n6m | shift[{p, q}, z][p | q],
1026      h_, l_] :>
1027      spinor[pp[a], h, 1] -
1028      HelicitySign[h,
1029      l] (pp[5] -
1030      I HelicitySign[h, l] pp[4]) Sp[Sm[a, -h],
1031      spinor[pp[a], -h, 1]]/2/Mp[Mom[a], Mom[pp[a]]],
1032  spinorbar[pp : p | q | n6p | n6m | shift[{p, q}, z][p | q], h_,
1033      l_] :> spinorbar[pp[a], h, 1] +
1034      HelicitySign[h,
1035      l] (pp[5] -
1036      I HelicitySign[h, l] pp[4]) Sp[spinorbar[pp[a], -h, 1],
1037      Sm[a, -h]]/2/Mp[Mom[a], Mom[pp[a]]],
1038  Mom[pp : p | q | n6p | n6m | shift[{p, q}, z][p | q]] :>
1039  Mom[pp[a]] + Mom[a] (pp[4]^2 + pp[5]^2)/2/Mp[Mom[pp[a]], Mom[a]] +
1040  Sum[HelicitySign[h] Sp[spinorbar[pp[a], 1, h], \[Sigma][1],
1041      Sm[a, -1],
1042      spinor[pp[a], -1, -h]] (pp[5] +
1043      I HelicitySign[h] pp[4]), {h, {11, -11}}]/4/
1044  Mp[Mom[pp[a]], Mom[a]],
1045  Sm[pp : p | q | n6p | n6m | shift[{p, q}, z][p | q], h_] :>
1046  Sp[Sm[pp[a], h]] +
1047  Sum[HelicitySign[l] HelicitySign [
1048      h] (pp[5] -
1049      I HelicitySign[l] HelicitySign[h] pp[4]) (Sp[
1050      spinor[pp[a], -h, 1], spinorbar[pp[a], h, -1], Sm[a, h]] -

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1051      Sp[Sm[a, h], spinor[pp[a], h, -1],
1052      spinorbar[pp[a], -h, 1] ], {1, {11, -11}}/2/
1053      Mp[Mom[pp[a]], Mom[a]] + (pp[4]^2 + pp[5]^2) Sp[Sm[a, h]]/2/
1054      Mp[Mom[pp[a]], Mom[a]];
1055      unrep = {spinor[(pp : p | q | shift[{p, q}, z][p | q])[a], h-, 1-] :>
1056      spinor[pp, h, 1] +
1057      HelicitySign[h,
1058      1] (pp[5] -
1059      I HelicitySign[h, 1] pp[4]) Sp[Sm[a, -h], spinor[pp, -h, 1]]/
1060      2/Mp[Mom[a], Mom[pp]],
1061      spinorbar[(pp : p | q | shift[{p, q}, z][p | q])[a], h-, 1-] :>
1062      spinorbar[pp, h, 1] -
1063      HelicitySign[h,
1064      1] (pp[5] -
1065      I HelicitySign[h, 1] pp[4]) Sp[spinorbar[pp, -h, 1],
1066      Sm[a, -h]]/2/Mp[Mom[a], Mom[pp]],
1067      Mom[(pp : p | q | shift[{p, q}, z][p | q])[a]] :>
1068      Mom[pp] + Mom[a] (pp[4]^2 + pp[5]^2)/2/Mp[Mom[pp[a]], Mom[a]] -
1069      Sum[h Sp[spinorbar[pp, 1, h], \[Sigma][1], Sm[a, -1],
1070      spinor[pp, -1, -h]] (pp[5] + I h pp[4]), {h, {1, -1}}]/4/
1071      Mp[Mom[pp], Mom[a]],
1072      Sm[(pp : p | q | shift[{p, q}, z][p | q])[a], h-] :>
1073      Sp[Sm[pp, h]] -
1074      Sum[1 HelicitySign [
1075      h] (pp[5] -
1076      I 1 HelicitySign[h] pp[4]) (Sp[spinor[pp, -h, 1],
1077      spinorbar[pp, h, -1], Sm[a, h]] -
1078      Sp[Sm[a, h], spinor[pp, h, -1],
1079      spinorbar[pp, -h, 1] ]), {1, {1, -1}}]/2/
1080      Mp[Mom[pp], Mom[a]] + (pp[4]^2 + pp[5]^2) Sp[Sm[a, h]]/2/
1081      Mp[Mom[pp], Mom[a]]];
1082
1083      tomu = {(pp : p | q)[
1084      4] :> (I/2)*(Subsuperscript[\[Mu], pp, "-"] -
1085      Subsuperscript[\[Mu], pp, "+"]), (pp : p | q)[
1086      5] :> (Subsuperscript[\[Mu], pp, "-"] +
1087      Subsuperscript[\[Mu], pp, "+"])/2};
1088      tomuh[h-] := {(pp : p | q)[4] :> (I/2)*
1089      HelicitySign [
1090      h] (Subsuperscript[\[Mu], pp, -h] -
1091      Subsuperscript[\[Mu], pp, h]), (pp : p | q)[
1092      5] :> (Subsuperscript[\[Mu], pp, -h] +

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1093     Subsuperscript [\[Mu], pp, h])/2}
1094 tomuhp[h_] := {(pp : p | q)[4] :> (I/2)*
1095     HelicitySign[
1096     h[pp]] (Subsuperscript [\[Mu], pp, -h[pp]] -
1097     Subsuperscript [\[Mu], pp, h[pp]]), (pp : p | q)[
1098     5] :> (Subsuperscript [\[Mu], pp, -h[pp]] +
1099     Subsuperscript [\[Mu], pp, h[pp]])/2}
1100 frommu = {Subsuperscript [\[Mu], p-, "+" ] :> p[5] + I p[4],
1101     Subsuperscript [\[Mu], p-, "-" ] :> p[5] - I p[4],
1102     Subsuperscript [\[Mu], p-, h_] :> p[5] + I HelicitySign[h] p[4]}

```