

A covariant representation of the Ball-Chiu vertex

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Abstract

In nonabelian gauge theory the three-gluon vertex function contains important structural information, in particular on infrared divergences, and is also an essential ingredient in the Schwinger-Dyson equations. Much effort has gone into analyzing its general structure, and at the one-loop level also a number of explicit computations have been done, using various approaches. Here we use the string-inspired formalism to unify the calculations of the scalar, spinor and gluon loop contributions to the one-loop vertex, leading to an extremely compact representation in all cases. The vertex is computed fully off-shell and in dimensionally continued form, so that it can be used as a building block for higher-loop calculations. We find that the Bern-Kosower loop replacement rules, originally derived for the on-shell case, hold off-shell as well. We explain the relation of the structure of this representation to the low-energy effective action, and establish the precise connection with the standard Ball-Chiu decomposition of the vertex. This allows us also to predict that the vanishing of the completely antisymmetric coefficient function S of this decomposition is not a one-loop accident, but persists at higher loop orders. The sum rule found by Binger and Brodsky, which leads to the vanishing of the one-loop vertex in $N = 4$ SYM theory, in the present approach relates to worldline supersymmetry.

Keywords: Three-gluon vertex, Ball-Chiu decomposition, string-inspired formalism, Bern-Kosower formalism

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1. Introduction

The one-particle-irreducible (‘1PI’) off-shell three-gluon Green’s function (in the following simply called “three-vertex” or “vertex”) is a basic object of interest in nonabelian gauge theory and quantum chromodynamics. It contains important structural information, in particular on infrared divergences (see, e.g., [1]) and refs. therein) and is a main ingredient of the Schwinger-Dyson equations. In perturbation theory it can, computed explicitly to a certain loop order, in principle be used as a convenient building block for higher-loop calculations.

However, explicit calculations of the three-vertex have so far been essentially restricted to the one-loop level [2, 3, 4, 5, 6, 7] (at two loops, the three-gluon vertex has been obtained so far only for some very special momentum configurations [8, 9, 10]). In this paper we will recalculate, in a simple and unifying way, the scalar, spinor and gluon loop contributions to the one-loop three-vertex (with “gluon” we mean any nonabelian gauge boson). In fig. 1 for definiteness we show the fermion loop contribution.

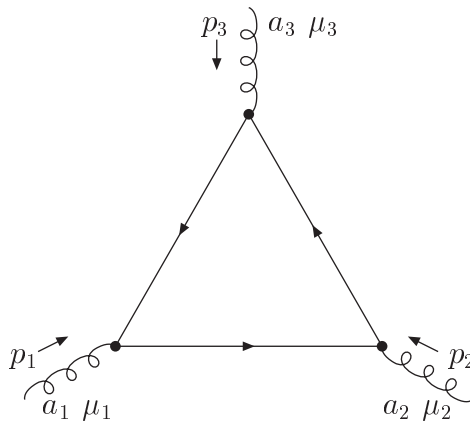


Figure 1: Three-gluon vertex.

Following the notation of [5, 6], we write

$$\Gamma_{\mu_1\mu_2\mu_3}^{a_1a_2a_3}(p_1, p_2, p_3) = -igf^{a_1a_2a_3}\Gamma_{\mu_1\mu_2\mu_3}(p_1, p_2, p_3). \quad (1.1)$$

The gluon momenta are ingoing and $p_1 + p_2 + p_3 = 0$. There are actually two diagrams differing by the two inequivalent orderings of the three gluons along the loops (or equivalently by a change of the fermion line orientation). Those diagrams add to produce a factor of two.

By an analysis of the nonabelian gauge Ward identities, Ball and Chiu [4] in 1980 found a form factor decomposition of this vertex which is valid at any order in perturbation theory, and also independent of whether the particle in the loop is a scalar, fermion or gluon, with the only restriction that a covariant gauge be used. It can be written as

$$\begin{aligned}
\Gamma_{\mu_1\mu_2\mu_3}(p_1, p_2, p_3) = & A(p_1^2, p_2^2; p_3^2)g_{\mu_1\mu_2}(p_1 - p_2)_{\mu_3} + B(p_1^2, p_2^2; p_3^2)g_{\mu_1\mu_2}(p_1 + p_2)_{\mu_3} \\
& + C(p_1^2, p_2^2; p_3^2) \left[p_{1\mu_2}p_{2\mu_1} - p_1 \cdot p_2 g_{\mu_1\mu_2} \right] (p_1 - p_2)_{\mu_3} \\
& + \frac{1}{3}S(p_1^2, p_2^2, p_3^2) \left[p_{1\mu_3}p_{2\mu_1}p_{3\mu_2} + p_{1\mu_2}p_{2\mu_3}p_{3\mu_1} \right] \\
& + F(p_1^2, p_2^2; p_3^2) \left[p_{1\mu_2}p_{2\mu_1} - p_1 \cdot p_2 g_{\mu_1\mu_2} \right] \left[p_1 \cdot p_3 p_{2\mu_3} - p_2 \cdot p_3 p_{1\mu_3} \right] \\
& + H(p_1^2, p_2^2, p_3^2) \left\{ -g_{\mu_1\mu_2} \left[p_{1\mu_3}p_2 \cdot p_3 - p_{2\mu_3}p_1 \cdot p_3 \right] \right. \\
& \quad \left. + \frac{1}{3} \left[p_{1\mu_3}p_{2\mu_1}p_{3\mu_2} - p_{1\mu_2}p_{2\mu_3}p_{3\mu_1} \right] \right\} \\
& + \left\{ \text{cyclic permutations of } (p_1, \mu_1), (p_2, \mu_2), (p_3, \mu_3) \right\}.
\end{aligned} \tag{1.2}$$

Here the functions A , C and F are symmetric in the first two arguments, the function B is antisymmetric in the first two arguments, H is totally symmetric and S totally antisymmetric with respect to interchange of any pair of arguments. A different decomposition of the three-gluon vertex was proposed in [11].

Although we wish to study the off-shell vertex, with our calculation method it will be convenient to contract it with polarization vectors $\varepsilon_{1,2,3}$. Those vectors are arbitrary and serve book-keeping purposes only. Thus we will use (1.2) in the form

$$\begin{aligned}
\varepsilon_1^{\mu_1} \varepsilon_2^{\mu_2} \varepsilon_3^{\mu_3} \Gamma_{\mu_1\mu_2\mu_3}(p_1, p_2, p_3) \equiv & \left\{ A(p_1^2, p_2^2; p_3^2)T_A + B(p_1^2, p_2^2; p_3^2)T_B + C(p_1^2, p_2^2; p_3^2)T_C \right. \\
& \left. + F(p_1^2, p_2^2; p_3^2)T_F \right\} + \{2 \text{ cyclic permutations}\} \\
& + H(p_1^2, p_2^2, p_3^2)T_H + S(p_1^2, p_2^2, p_3^2)T_S.
\end{aligned} \tag{1.3}$$

Note that the tensors T_F and T_H are totally transversal, i.e., they give zero when any ε_i is replaced by p_i .

For the gluon loop in Feynman gauge and at the one-loop level, Ball and Chiu also calculated the coefficient functions A to H , up to their constant terms in the ϵ -expansion; here it turned out that S actually vanishes. Later Davydychev, Osland and Tarasov [5] computed this gluon loop contribution vertex more generally for an arbitrary covariant gauge, and in arbitrary spacetime dimension.

The quark loop contribution to the vertex was first calculated for massless quarks and in the symmetric limit $p_1^2 = p_2^2 = p_3^2$ by Celmaster and Gonsalves [2] and Pascual and Tarrach [3]. For general off-shell momenta the massless quark loop contribution was obtained in [5], and the massive quark loop one in [6] (again in arbitrary spacetime dimension).

However, this is not yet the whole story, since for the gluon loop contribution to the vertex there are subtle issues with gauge dependence. When calculated in the standard formalism using any covariant gauge, it satisfies rather complicated Slavnov-Taylor identities involving not only the gluon propagator, but also the ghost propagator and the gluon-ghost-ghost vertex (see, e.g., [12, 4, 7]). The scalar and fermion loop contributions, on the other hand, satisfy the simple QED-like Ward identity

$$\begin{aligned} \Gamma(\varepsilon_3 \rightarrow p_3) &= -(p_1^2 \varepsilon_1 \cdot \varepsilon_2 - p_1 \cdot \varepsilon_1 p_1 \cdot \varepsilon_2) \left(1 + \Pi(p_1^2)\right) \\ &\quad + (p_2^2 \varepsilon_1 \cdot \varepsilon_2 - p_2 \cdot \varepsilon_1 p_2 \cdot \varepsilon_2) \left(1 + \Pi(p_2^2)\right) \end{aligned} \tag{1.4}$$

where $\Pi(p^2)$ is the corresponding vacuum polarization function. Having the same simple Ward identity also for the gluon loop case is possible, but requires more sophisticated techniques. It can be achieved using either the background field method (‘BFM’) [13, 14, 15] with Feynman gauge for the quantum field, or the pinch technique [16, 17]. Although very different, those two methods turn out to lead to precisely the same Green’s functions [18, 19]. The corresponding three-gluon vertex, also called the “gauge-invariant vertex”, was studied by Freedman et al. [20] with an emphasis on its conformal properties. Binger and Brodsky [7] explicitly calculated it in the transversality-based Ball-Chiu decomposition (1.2), as well as in a different basis related to current conservation. They also calculated the scalar loop contribution, which enabled them to find certain sum rules between the massless scalar, spinor and gluon loop contributions. In particular, with all particles massless and in the adjoint representation they obtain the identity

$$3\Gamma_{\text{scalar}} + 2\Gamma_{\text{spinor}} + \Gamma_{\text{gluon}} = 0 \quad (1.5)$$

(here it is understood that the contribution of the gluon includes the one of its ghost). This identity is related to supersymmetry, and implies that the off-shell three-gluon amplitude in $N = 4$ Super-Yang-Mills theory vanishes (a statement which, contrary to the vanishing of the gluon propagator in that theory, does not follow from UV finiteness in an obvious way).

In the present work, we will recalculate the scalar, spinor and gluon loop contributions to this “gauge-invariant” three-gluon vertex using the “string-inspired” formalism along the lines of [21, 22, 23, 24, 25, 26, 27] (for a review, see [28]). Our starting point is the “Bern-Kosower master formula” [21, 22, 23]:

$$\begin{aligned} & \Gamma_{\text{scalar}}^{a_1 \dots a_N} [p_1, \varepsilon_1; \dots; p_N, \varepsilon_N] \\ &= (-ig)^N \text{tr}(T^{a_1} \dots T^{a_N}) (2\pi)^D i \delta(\sum p_i) \int_0^\infty dT (4\pi T)^{-D/2} e^{-m^2 T} \\ & \times \int_0^T d\tau_1 \int_0^{\tau_1} d\tau_2 \dots \int_0^{\tau_{N-2}} d\tau_{N-1} \\ & \times \exp \left\{ \sum_{i,j=1}^N \left[\frac{1}{2} G_{Bij} p_i \cdot p_j - i \dot{G}_{Bij} \varepsilon_i \cdot p_j + \frac{1}{2} \ddot{G}_{Bij} \varepsilon_i \cdot \varepsilon_j \right] \right\} \Big|_{\text{lin}(\varepsilon_1 \dots \varepsilon_N)}. \end{aligned} \quad (1.6)$$

As it stands, this formula represents the color-ordered contribution to the 1PI N - gluon amplitude due to a (complex) scalar loop of mass m , calculated in D spacetime dimensions. The i th gluon carries the momentum p_i , polarization ε_i and a gauge group generator T^{a_i} in some representation. T is the total proper-time length of the loop, and τ_i is the position in proper-time along the loop of gluon i . One integration is redundant and has been eliminated by setting $\tau_N = 0$. The derivation of this formula involved a formal exponentiation, which needs to be undone by expanding out the exponential factor and keeping only the terms linear in each of the N polarization vectors. The color-ordering means that one still has to sum over all $(N-1)!$ inequivalent orderings of the gluons along the loop to get the full amplitude. $G_{Bij} \equiv G_B(\tau_i, \tau_j)$ denotes the “bosonic” worldline Green’s function, defined by

$$G_B(\tau_1, \tau_2) = |\tau_1 - \tau_2| - \frac{(\tau_1 - \tau_2)^2}{T}, \quad (1.7)$$

and dots generally denote a derivative acting on the first variable. Explicitly,

$$\begin{aligned}\dot{G}_B(\tau_1, \tau_2) &= \text{sign}(\tau_1 - \tau_2) - 2\frac{(\tau_1 - \tau_2)}{T}, \\ \ddot{G}_B(\tau_1, \tau_2) &= 2\delta(\tau_1 - \tau_2) - \frac{2}{T}.\end{aligned}\tag{1.8}$$

The master formula (1.6) was originally derived from string theory [21, 22, 23], starting from a representation of the N -gluon amplitude for the heterotic string and analyzing its field theory limit. Here the object of interest was, however, the full N -gluon amplitude on-shell, rather than the 1PI amplitude off-shell. Thus on one hand on-shell conditions were used from the beginning, already at the string level; on the other hand the fact that the distinction between reducible and irreducible diagrams emerges only in the field theory limit made it possible to establish certain formal rules that allow one to reconstruct, from the formula for the 1PI amplitude (1.6), also all the missing reducible contributions to the full on-shell matrix element. Bern and Kosower were moreover able to derive simple “loop replacement rules”, based on worldsheet supersymmetry, that allow one to obtain from (1.6) also integral representations for the spinor and gluon loop contributions to the full on-shell N -gluon amplitudes [21, 22, 23]. We need not discuss these “Bern-Kosower rules” here in full, but it is important to note that they all involve integration-by-parts (‘IBP’) in an essential way. Namely, performing the expansion of the exponential factor in (1.6) will yield an integrand $\sim P_N e^{(\cdot)}$, where we abbreviated

$$e^{(\cdot)} := \exp\left\{\frac{1}{2}\sum_{i,j=1}^N G_{Bij} p_i \cdot p_j\right\},\tag{1.9}$$

and P_N is a polynomial in $\dot{G}_{Bij}, \ddot{G}_{Bij}$ and the kinematic invariants. It is possible to remove all second derivatives \ddot{G}_{Bij} appearing in P_N by suitable integrations by parts, leading to a new integrand $\sim Q_N e^{(\cdot)}$ which is the real starting point for the application of the Bern-Kosower rules. Relevant for the following will be only the “loop replacement rules”. The rule for passing from the scalar to the spinor loop is easy to state in general: look in Q_N for “ τ -cycles”, that is, products of \dot{G}_{Bij} ’s whose indices form a closed chain. A τ -cycle can thus be written as $\dot{G}_{B_{i_1 i_2}} \dot{G}_{B_{i_2 i_3}} \cdots \dot{G}_{B_{i_n i_1}}$ (to put it into this form may require the use of the antisymmetry of \dot{G}_B , e.g. $\dot{G}_{B_{12}} \dot{G}_{B_{12}} =$

$-\dot{G}_{B_{12}}\dot{G}_{B_{21}}$). Then, apart from a global factor of -2 correcting for degrees of freedom and statistics, the integrand for the spinor loop case can be obtained from the one for the scalar loop simply by simultaneously replacing every τ -cycle appearing in Q_N by

$$\dot{G}_{B_{i_1 i_2}}\dot{G}_{B_{i_2 i_3}}\cdots\dot{G}_{B_{i_n i_1}} \rightarrow \dot{G}_{B_{i_1 i_2}}\dot{G}_{B_{i_2 i_3}}\cdots\dot{G}_{B_{i_n i_1}} - G_{F_{i_1 i_2}}G_{F_{i_2 i_3}}\cdots G_{F_{i_n i_1}}, \quad (1.10)$$

where $G_{F_{12}} \equiv \text{sign}(\tau_1 - \tau_2)$ denotes the ‘fermionic’ worldline Green’s function. The rule for passing from the scalar to the gluon loop is similar but somewhat more complicated, and will be dealt with in chapter 4 below.

This formalism was used for the first calculation of the one-loop on-shell QCD five gluon amplitudes [29], but not further employed for such on-shell multi-gluon amplitude calculations due to the emergence of other extremely powerful methods for the computation of one-loop on-shell amplitudes such as generalized unitarity; see, e.g., [30].

In the present paper, we will instead start an effort to exploit the Bern-Kosower formalism as a tool for the calculation of the one-loop N -gluon vertex. Thus we will need an off-shell extension of this formalism, which is provided by its worldline path integral formulation due to Strassler [24, 25]. Here the starting point is the following path integral representation of the nonabelian one-loop effective action due to a scalar loop [24] (this generalizes Feynman’s famous 1950 formula for scalar QED [31]),

$$\Gamma_{\text{scalar}}[A] = \int_0^\infty \frac{dT}{T} e^{-m^2 T} \text{tr} \int \mathcal{D}x \exp \left[- \int_0^T d\tau \left(\frac{1}{4} \dot{x}^2 + ig \dot{x} \cdot A \right) \right], \quad (1.11)$$

where the integral $\int \mathcal{D}x$ is over the space of all closed trajectories in space-time with periodicity T in proper-time, $x^\mu(T) = x^\mu(0)$.

Although this approach fell somewhat short of yielding a rederivation of the full set of Bern-Kosower rules including the non-1PI contributions, it provides a simple way to see that the master formula (1.6) as it stands is valid off-shell as a formula for the 1PI part of the N -gluon Green’s function (see [24, 28]). In [25], Strassler moreover started a systematic investigation of the IBP procedure, and discovered that it bears an interesting connection to gauge invariance. Namely, it so turns out that, once all $\ddot{G}_{B_{ij}}$ ’s have been removed and all terms contributing to a given ‘ τ -cycle’ $\dot{G}_{B_{i_1 i_2}}\dot{G}_{B_{i_2 i_3}}\cdots\dot{G}_{B_{i_n i_1}}$ been combined, the sum of their Lorentz factors can be written as a ‘Lorentz cycle’ Z_n , defined by

$$\begin{aligned}
Z_2(ij) &\equiv \frac{1}{2} \text{tr}(f_i f_j), \\
Z_n(i_1 i_2 \dots i_n) &\equiv \text{tr}\left(\prod_{j=1}^n f_{i_j}\right) \quad (n \geq 3),
\end{aligned}
\tag{1.12}$$

where

$$f_i^{\mu\nu} \equiv p_i^\mu \varepsilon_i^\nu - \varepsilon_i^\mu p_i^\nu \tag{1.13}$$

is the momentum space form of the abelian field strength tensor. Z_n generalizes the transversal projector which is familiar from the two-point case. However, in [25] no systematic way was found to perform the partial integrations at arbitrary N , nor how to preserve the permutation symmetry. This issue was taken up again in [27], where a definite and computerizable IBP algorithm was given which works for any N and preserves the full permutation symmetry. This algorithm is still not satisfactory from the point of view of gauge invariance, though. A given term in the integrand after the IBP in general has not only cycle factors, but also a leftover, called “tail”, and the algorithm arranges into field strength tensors only the polarization vectors contained in the cycles, not the ones in the tails. Only very recently an extension of the algorithm of [27] was found which, for any N and preserving the permutation invariance, achieves this “covariantization” for all the polarization vectors, including the ones in tails [32].

This in some sense completes the investigation started in [25]. It also suggests that, with this optimized IBP at hand, the string-inspired formalism might become a powerful tool for the computation of the N -vertex. This is for three reasons: (i) The covariantization means that the bulk integrand after the IBP is manifestly transversal, so that any nontransversality must come from boundary terms. Thus the IBP procedure itself should generate a transversality-based form factor decomposition similar to the Ball-Chiu one (1.2). (ii) Like the Ball-Chiu one, this decomposition will respect the cyclic invariance (which is the remnant of the permutation invariance after the color ordering). (iii) The work of [24] also suggests that the “loop-replacement” part of the Bern-Kosower rules may hold off-shell for the 1PI amplitudes, which would reduce the calculational effort very significantly.

Here we will recalculate the three-gluon vertex along the above lines, and find these expectations to be fully justified. The organization of the paper is as follows:

In chapter 2 we will start with the scalar loop contribution to the vertex. We perform the IBP using the old algorithm of [27] as well as the improved one of [32]. With both choices we obtain a very compact integral representation for the vertex, however the new algorithm has the advantage that all non-transversality is pushed into the boundary (two-point) terms.

In chapters 3 and 4 we show that the “loop replacement rules” indeed hold for the three-gluon vertex. As is well-known, the Dirac fermion possesses an $N = 1$ supersymmetric worldline path integral representation [33, 34, 35, 36, 37, 38, 39], and the gluon an $N = 2$ supersymmetric one [33, 35, 40, 24, 41]. Analogously to the original string-based derivation of those rules [21, 22, 23], where worldsheet SUSY was identified as the underlying symmetry, in the worldline approach the same rules can be related to this worldline SUSY [24, 28].

In chapter 5 we summarize and unify our results for the scalar, spinor and gluon loop, and in chapter 6 we establish their exact relation to the Ball-Chiu decomposition (1.2).

The 1PI vertices hold the same information as the effective action. Nevertheless, contrary to the QED case where there is no essential difference between the calculation of the effective action and of the off-shell N -photon amplitudes, in the nonabelian case the effective action is mathematically an intrinsically more natural object. This is because it can be written in terms of full field strength tensors

$$F_{\mu\nu} \equiv F_{\mu\nu}^a T^a = (\partial_\mu A_\nu^a - \partial_\nu A_\mu^a) T^a + ig[A_\mu^b T^b, A_\nu^c T^c] \quad (1.14)$$

whereas upon Fourier transformation those inevitably get split up into their “abelian parts” $f_{\mu\nu}^a := \partial_\mu A_\nu^a - \partial_\nu A_\mu^a$ and the commutator terms. This suggests that the analysis of the structure of the 1PI vertices should benefit from a comparison with the low energy expansion of the effective action, and indeed we will show in chapter 7 for the three-point case that in the present formalism, due to the systematic generation of “abelian” field strength tensors and commutator terms by the IBP, it is possible to keep the relation between the effective action and the vertex very transparent.

Our conclusions are given in chapter 8. In particular, we give there a general argument showing that the off-shell validity of the loop replacement rules extends to the N -vertex. Appendix A lists our conventions.

2. The scalar loop case

We proceed to the calculation of the (off-shell, 1PI) three-gluon amplitude for a scalar loop. It will be instructive to consider the two-point case first.

For $N = 2$ we get from the master formula (1.6), after expanding out the exponential (in the following we generally omit the global factor $(2\pi)^4 i \delta(\sum p_i)$ for energy-momentum conservation),

$$\begin{aligned} \Gamma^{a_1 a_2}[p_1, \varepsilon_1; p_2, \varepsilon_2] &= (-ig)^2 \text{tr}(T^{a_1} T^{a_2}) \int_0^\infty dT (4\pi T)^{-D/2} e^{-m^2 T} \\ &\quad \times \int_0^T d\tau_1 (-i)^2 P_2 e^{G_{B12} p_1 \cdot p_2} \end{aligned} \quad (2.1)$$

where

$$P_2 = \dot{G}_{B12} \varepsilon_1 \cdot p_2 \dot{G}_{B21} \varepsilon_2 \cdot p_1 - \ddot{G}_{B12} \varepsilon_1 \cdot \varepsilon_2. \quad (2.2)$$

By an IBP of the term involving \ddot{G}_{B12} we can remove the second derivative and transform P_2 into Q_2 ,

$$\begin{aligned} Q_2 &= \dot{G}_{B12} \dot{G}_{B21} (\varepsilon_1 \cdot p_2 \varepsilon_2 \cdot p_1 - \varepsilon_1 \cdot \varepsilon_2 p_1 \cdot p_2) \\ &= \frac{1}{2} \dot{G}_{B12} \dot{G}_{B21} \text{tr}(f_1 f_2) = \dot{G}_{B12} \dot{G}_{B21} Z_2(12). \end{aligned} \quad (2.3)$$

Thus the IBP has allowed us to absorb the polarization vectors into the ‘‘abelian’’ field strength tensors f_i , defined in (1.13), thereby making the transversality of the two-point function manifest.

At the three-point level, the expansion of (1.6) yields

$$\begin{aligned} \Gamma^{a_1 a_2 a_3}[p_1, \varepsilon_1; p_2, \varepsilon_2; p_3, \varepsilon_3] &= (-ig)^3 \text{tr}(T^{a_1} T^{a_2} T^{a_3}) \int_0^\infty dT (4\pi T)^{-D/2} e^{-m^2 T} \\ &\quad \times \int_0^T d\tau_1 \int_0^{\tau_1} d\tau_2 (-i)^3 P_3 e^{(\cdot)} \end{aligned} \quad (2.4)$$

where

$$\begin{aligned}
P_3 = & \dot{G}_{B1i\varepsilon_1} \cdot p_i \dot{G}_{B2j\varepsilon_2} \cdot p_j \dot{G}_{B3k\varepsilon_3} \cdot p_k - \ddot{G}_{B12\varepsilon_1} \cdot \varepsilon_2 \dot{G}_{B3k\varepsilon_3} \cdot p_k \\
& - \ddot{G}_{B13\varepsilon_1} \cdot \varepsilon_3 \dot{G}_{B2j\varepsilon_2} \cdot p_j - \ddot{G}_{B23\varepsilon_2} \cdot \varepsilon_3 \dot{G}_{B1i\varepsilon_1} \cdot p_i
\end{aligned} \tag{2.5}$$

and we have introduced the convention that repeated indices i, j, k, \dots are to be summed from 1 to $N = 3$. To remove, e.g., the term involving $\ddot{G}_{B12}\dot{G}_{B31}$ in the second term of P_3 , we add the total derivative

$$-\frac{\partial}{\partial \tau_2} \left(\dot{G}_{B12\varepsilon_1} \cdot \varepsilon_2 \dot{G}_{B31\varepsilon_3} \cdot p_1 e^{(G_{B12}p_1 \cdot p_2 + G_{B13}p_1 \cdot p_3 + G_{23}p_2 \cdot p_3)} \right). \tag{2.6}$$

Adding five more similar total derivative terms removes all the \ddot{G}_B 's. Decomposing the new integrand according to its ‘‘cycle content’’, P_3 gets replaced by $Q_3 = Q_3^3 + Q_3^2$, where

$$\begin{aligned}
Q_3^3 &= \dot{G}_{B12} \dot{G}_{B23} \dot{G}_{B31} Z_3(123), \\
Q_3^2 &= \dot{G}_{B12} \dot{G}_{B21} Z_2(12) \dot{G}_{B3k\varepsilon_3} \cdot p_k + \dot{G}_{B13} \dot{G}_{B31} Z_2(13) \dot{G}_{B2j\varepsilon_2} \cdot p_j \\
&\quad + \dot{G}_{B23} \dot{G}_{B32} Z_2(23) \dot{G}_{B1i\varepsilon_1} \cdot p_i.
\end{aligned} \tag{2.7}$$

Note that Q_3^3 contains a τ -cycle of length three and Q_3^2 of length two, as indicated by the upper indices, and that each τ -cycle appears together with the corresponding ‘‘Lorentz-cycle’’, as advertised in the introduction. The terms of Q_3^2 have, apart from the cycle, also a ‘‘one-tail’’, defined by

$$T_1(a) := \varepsilon_a \cdot p_i \dot{G}_{Bai}. \tag{2.8}$$

Although the form of the integrand reached in (2.7) is already suitable for the application of the Bern-Kosower rules, it is natural to ask whether the polarization vectors appearing in the tails can also somehow be completed to field strength tensors. Now in this three-point case there are already various chains of integrations-by-part that can be used to remove all the \ddot{G}_B 's, but if one assumes that the corresponding total derivative terms are added with constant coefficients (i.e., they involve no functions of momentum or polarization other than the ones already present in the term which one wishes to modify), then it is easy to convince oneself that they all lead to

the same Q_3 of (2.7). This applies, in particular, to the “old” IBP procedure proposed in [27], where this Q_3 is obtained by a chain of IBPs different from the above. Thus a more general type of IBP’s is called for if one wishes to achieve this “covariantization of the tails”, and in a companion paper [32] it will be shown how, using total derivative terms with coefficients that *do* depend on momenta and polarizations, this can indeed be done for arbitrary N . Here we need not discuss this matter in more depth, since for $N = 3$ the solution of this problem is still very simple. Consider the first term in Q_3^2 above, eq. (2.7). Choose a momentum vector r_3 such that $r_3 \cdot p_3 \neq 0$, and add the total derivative

$$-\frac{r_3 \cdot \varepsilon_3}{r_3 \cdot p_3} Z_2(12) \frac{\partial}{\partial \tau_3} \left(\dot{G}_{B12} \dot{G}_{B21} e^{(\cdot)} \right). \quad (2.9)$$

The addition of this term to the first term in Q_3^2 , and of similar terms to the second and third one, transforms Q_3^2 into

$$\begin{aligned} R_3^2 := & \dot{G}_{B12} \dot{G}_{B21} Z_2(12) \dot{G}_{B3k} \frac{r_3 \cdot f_3 \cdot p_k}{r_3 \cdot p_3} + \dot{G}_{B13} \dot{G}_{B31} Z_2(13) \dot{G}_{B2j} \frac{r_2 \cdot f_2 \cdot p_j}{r_2 \cdot p_2} \\ & \dot{G}_{B23} \dot{G}_{B32} Z_2(23) \dot{G}_{B1i} \frac{r_1 \cdot f_1 \cdot p_i}{r_1 \cdot p_1}. \end{aligned} \quad (2.10)$$

Thus now all polarization vectors have been absorbed into tensors f_i , leading to manifest transversality. This IBP procedure can be systematized to obtain closed-form integral representations of the Scalar and Spinor QED N - photon amplitudes that are manifestly gauge invariant at the integrand level [32].

Here, however, we are in the nonabelian case, where the color-induced restriction of the parameter integrations to ordered sectors leads to the appearance of boundary terms in the IBP [24, 25]. Let us look again at our total derivative term (2.6). In the abelian case it would be integrated over the whole circle, and the result would be zero, since the worldline Green’s function $G_B(\tau_1, \tau_2)$ has the appropriate periodicity properties to make the two boundary terms cancel. Here instead we find a nonzero result:

$$-\dot{G}_{B12} \varepsilon_1 \cdot \varepsilon_2 \dot{G}_{B31} \varepsilon_3 \cdot p_1 e^{(\cdot)} \Big|_{\tau_2=\tau_3}^{\tau_2=\tau_1} = 0 + \dot{G}_{B13} \varepsilon_1 \cdot \varepsilon_2 \dot{G}_{B31} \varepsilon_3 \cdot p_1 e^{G_{B13} p_1 \cdot (p_2 + p_3)}. \quad (2.11)$$

Now, in the three-point case there are already two inequivalent orderings, say, (123) and (132); thus the full amplitude will also have a part $\Gamma^{a_1 a_3 a_2}$ with color trace $\text{tr}(T^{a_1} T^{a_3} T^{a_2})$, and the same total derivative term will contribute to it a boundary term

$$-\dot{G}_{B12}\varepsilon_1 \cdot \varepsilon_2 \dot{G}_{B31}\varepsilon_3 \cdot p_1 e^{(\cdot)} \Big|_{\tau_2=\tau_1}^{\tau_2=\tau_3} = -\dot{G}_{B13}\varepsilon_1 \cdot \varepsilon_2 \dot{G}_{B31}\varepsilon_3 \cdot p_1 e^{G_{B13}p_1 \cdot (p_2+p_3)} - 0. \quad (2.12)$$

These two boundary terms would cancel in the abelian case, but now instead combine to produce a color commutator $\text{tr}(T^{a_1}[T^{a_2}, T^{a_3}])$. Moreover, among the other five similar total derivative terms needed to convert P_3 into Q_3 there is one that differs from (2.6) only by the interchange $2 \leftrightarrow 3$. With some relabeling of integration variables, we can combine the two boundary terms generated by that term with the two above to the structure

$$\text{tr}(T^{a_1}[T^{a_2}, T^{a_3}])\varepsilon_3 \cdot f_1 \cdot \varepsilon_2 \dot{G}_{B12}\dot{G}_{B21} e^{G_{B12}p_1 \cdot (p_2+p_3)}. \quad (2.13)$$

Comparing with (2.1) and (2.3) we note that this term has a parameter integral identical to the one of the two-gluon amplitude, except for the replacement of p_2 by $p_2 + p_3$. In terms of the effective action, from (1.14) and (2.13) its role is evidently to provide a piece needed to extend the ‘‘abelian’’ Maxwell term $\text{tr}(f_{\mu\nu}f^{\mu\nu})$ to the full nonabelian one $\text{tr}(F_{\mu\nu}F^{\mu\nu})$. We will discuss this in more detail in chapter 7 below.

To summarize so far, we can decompose the three-point amplitude for the scalar loop as (here and in the following we will often suppress the superscript ‘‘ $a_1 a_2 a_3$ ’’)

$$\Gamma_{\text{scalar}} = \frac{g^3}{(4\pi)^{\frac{D}{2}}} (\Gamma_{\text{scalar}}^3 + \Gamma_{\text{scalar}}^2 + \Gamma_{\text{scalar}}^{\text{bt}}) \quad (2.14)$$

where (note the global factor of $(-i)^6 = -1$)

$$\begin{aligned} \Gamma_{\text{scalar}}^3 &= -\text{tr}(T^{a_1} T^{a_2} T^{a_3}) \int_0^\infty \frac{dT}{T^{\frac{D}{2}}} e^{-m^2 T} \int_0^T d\tau_1 \int_0^{\tau_1} d\tau_2 Q_3^3 e^{(\cdot)} \\ &\quad -\text{tr}(T^{a_1} T^{a_3} T^{a_2}) \int_0^\infty \frac{dT}{T^{\frac{D}{2}}} e^{-m^2 T} \int_0^T d\tau_1 \int_0^{\tau_1} d\tau_3 Q_3^3 e^{(\cdot)}, \\ \Gamma_{\text{scalar}}^2 &= \Gamma_{\text{scalar}}^3(Q_3^3 \rightarrow Q_3^2), \end{aligned}$$

$$\begin{aligned}
\Gamma_{\text{scalar}}^{\text{bt}} &= -\text{tr}(T^{a_1}[T^{a_2}, T^{a_3}]) \int_0^\infty \frac{dT}{T^{\frac{D}{2}}} e^{-m^2 T} \int_0^T d\tau_1 \dot{G}_{B12} \dot{G}_{B21} \\
&\quad \times \left[\varepsilon_3 \cdot f_1 \cdot \varepsilon_2 e^{G_{B12} p_1 \cdot (p_2 + p_3)} + \varepsilon_1 \cdot f_2 \cdot \varepsilon_3 e^{G_{B12} p_2 \cdot (p_1 + p_3)} \right. \\
&\quad \left. + \varepsilon_2 \cdot f_3 \cdot \varepsilon_1 e^{G_{B12} p_3 \cdot (p_1 + p_2)} \right]
\end{aligned} \tag{2.15}$$

(here and in the following it is understood that always the last integration is eliminated by setting its integration variable equal to zero; e.g., for the ordering $\tau_1 > \tau_3 > \tau_2$ we set $\tau_2 = 0$).

Alternatively, we can replace Q_3^2 by R_3^2 in the Γ_{scalar}^2 part, but then we have to also add to $\Gamma_{\text{scalar}}^{\text{bt}}$ a term $\tilde{\Gamma}_{\text{scalar}}^{\text{bt}}$ containing the further boundary contributions coming from the total derivative terms of the type (2.9). Collecting those, one finds

$$\begin{aligned}
\tilde{\Gamma}_{\text{scalar}}^{\text{bt}} &= \frac{1}{2} \text{tr}(T^{a_1}[T^{a_2}, T^{a_3}]) \int_0^\infty \frac{dT}{T^{\frac{D}{2}}} e^{-m^2 T} \int_0^T d\tau_1 \dot{G}_{B12} \dot{G}_{B21} \\
&\quad \times \left\{ \left[\text{tr}(f_1 f_2) \rho_3 - \text{tr}(f_3 f_1) \rho_2 \right] e^{G_{B12} p_1 \cdot (p_2 + p_3)} + \left[\text{tr}(f_2 f_3) \rho_1 - \text{tr}(f_1 f_2) \rho_3 \right] e^{G_{B12} p_2 \cdot (p_1 + p_3)} \right. \\
&\quad \left. + \left[\text{tr}(f_3 f_1) \rho_2 - \text{tr}(f_2 f_3) \rho_1 \right] e^{G_{B12} p_3 \cdot (p_1 + p_2)} \right\}
\end{aligned} \tag{2.16}$$

where we have now abbreviated $\rho_i := r_i \cdot \varepsilon_i / r_i \cdot p_i$.

3. The spinor loop case

For the spinor loop case, it will be convenient to use the worldline super formalism [36, 37, 38, 42, 43, 44, 45]. In this formalism, one defines for each gluon leg a Grassmann variable θ_i , $\theta_i^2 = 0$, and also considers the polarization vectors ε_i as being Grassmann. Thus all $\varepsilon_j^\nu, \theta_k$, and $d\theta_l$ anticommute with each other. One further introduces the superderivative

$$D = \frac{\partial}{\partial \theta} - \theta \frac{\partial}{\partial \tau} \tag{3.1}$$

and the super proper-time distance

$$\hat{\tau}_{ij} := \tau_i - \tau_j + \theta_i \theta_j. \tag{3.2}$$

Then the Bern-Kosower master formula can be generalized to the case of a Dirac fermion loop case as follows [42, 28]:

$$\begin{aligned}
\Gamma_{\text{spinor}}^{a_1 \dots a_N} [p_1, \varepsilon_1; \dots; p_N, \varepsilon_N] &= -2(-ig)^N \text{tr}(T^{a_1} \dots T^{a_N}) \int_0^\infty \frac{dT}{T} (4\pi T)^{-\frac{D}{2}} e^{-m^2 T} \\
&\times \prod_{k=1}^N \int_0^T d\tau_k \int d\theta_k \delta\left(\frac{\tau_N}{T}\right) \vartheta(\hat{\tau}_{1N}) \prod_{l=1}^{N-1} \vartheta(\hat{\tau}_{l(l+1)}) \\
&\times \exp \left\{ \sum_{i,j=1}^N \left[\frac{1}{2} \hat{G}_{ij} p_i \cdot p_j + i D_i \hat{G}_{ij} \varepsilon_i \cdot p_j + \frac{1}{2} D_i D_j \hat{G}_{ij} \varepsilon_i \cdot \varepsilon_j \right] \right\} \Bigg|_{\text{lin}(\varepsilon_1 \dots \varepsilon_N)}.
\end{aligned} \tag{3.3}$$

Here ϑ is the Heaviside step function and the Green's functions G_B and G_F appear now combined into the super Green's function

$$\hat{G}(\tau_i, \theta_i; \tau_j, \theta_j) \equiv G_B(\tau_i, \tau_j) + \theta_i \theta_j G_F(\tau_i, \tau_j). \tag{3.4}$$

The overall sign of (3.3) refers to the standard ordering of the polarization vectors $\varepsilon_1 \varepsilon_2 \dots \varepsilon_N$. Next, note that

$$\vartheta(\hat{\tau}_{ij}) = \vartheta(\tau_i - \tau_j) + \theta_i \theta_j \delta(\tau_i - \tau_j). \tag{3.5}$$

The terms arising in the expansion of the spinor loop master formula (3.3) can thus be divided into three types: (i) terms that were there already for the scalar loop, (ii) new terms not involving any of the delta functions appearing in (3.5) and (iii) terms that do involve one such delta function (not more than one is possible in the three-point case).

Concerning the type (ii) terms, those are known already from the abelian case, and it was shown in [28] by a direct combinatorial argument, starting from the abelian version of (3.3), that they can be taken into account correctly by the ‘‘loop replacement rule’’ eq. (1.10). Thus in our three-point case their effect is to change each of the ‘‘bulk terms’’ $\Gamma_{\text{scalar}}^{2,3}$ of (2.15) to a corresponding $\Gamma_{\text{spinor}}^{2,3}$ differing from its scalar loop counterpart only by a change of $Q_3^{2,3}$ to $\hat{Q}_3^{2,3}$, where

$$\begin{aligned}
\hat{Q}_3^3 &= (\dot{G}_{B12} \dot{G}_{B23} \dot{G}_{B31} - G_{F12} G_{F23} G_{F31}) Z_3(123), \\
\hat{Q}_3^2 &= (\dot{G}_{B12} \dot{G}_{B21} - G_{F12} G_{F21}) Z_2(12) \dot{G}_{B3k} \varepsilon_3 \cdot p_k + \text{two permutations}.
\end{aligned} \tag{3.6}$$

Terms of type (iii) are specific to the nonabelian case. They would cancel between adjacent ordered sectors in the abelian case, but now produce color commutators; thus it is natural to think of them as a fermionic counterpart to the boundary terms encountered in the scalar loop calculation. Moreover, it is straightforward to check that they still conform to the same replacement rule. Taking also the global factor of -2 into account, our final result for the spinor case becomes

$$\Gamma_{\text{spinor}} = -2 \frac{g^3}{(4\pi)^{\frac{D}{2}}} (\Gamma_{\text{spinor}}^3 + \Gamma_{\text{spinor}}^2 + \Gamma_{\text{spinor}}^{\text{bt}}) \quad (3.7)$$

where

$$\begin{aligned} \Gamma_{\text{spinor}}^3 &= -\text{tr}(T^{a_1} T^{a_2} T^{a_3}) \int_0^\infty \frac{dT}{T^{\frac{D}{2}}} e^{-m^2 T} \int_0^T d\tau_1 \int_0^{\tau_1} d\tau_2 \hat{Q}_3^3|_{\tau_3=0} e^{(\cdot)} \\ &\quad -\text{tr}(T^{a_1} T^{a_3} T^{a_2}) \int_0^\infty \frac{dT}{T^{\frac{D}{2}}} e^{-m^2 T} \int_0^T d\tau_1 \int_0^{\tau_1} d\tau_3 \hat{Q}_3^3|_{\tau_2=0} e^{(\cdot)}, \\ \Gamma_{\text{spinor}}^2 &= \Gamma_{\text{spinor}}^3(\hat{Q}_3^3 \rightarrow \hat{Q}_3^2), \\ \Gamma_{\text{spinor}}^{\text{bt}} &= -\text{tr}(T^{a_1} [T^{a_2}, T^{a_3}]) \int_0^\infty \frac{dT}{T^{\frac{D}{2}}} e^{-m^2 T} \int_0^T d\tau_1 (\dot{G}_{B12} \dot{G}_{B21} - G_{F12} G_{F21}) \\ &\quad \times \left[\varepsilon_3 \cdot f_1 \cdot \varepsilon_2 e^{G_{B12} p_1 \cdot (p_2 + p_3)} + \varepsilon_1 \cdot f_2 \cdot \varepsilon_3 e^{G_{B12} p_2 \cdot (p_1 + p_3)} + \varepsilon_2 \cdot f_3 \cdot \varepsilon_1 e^{G_{B12} p_3 \cdot (p_1 + p_2)} \right]. \end{aligned} \quad (3.8)$$

The alternative form of the scalar loop result, involving R_3^2 instead of Q_3^2 and the additional boundary contribution $\tilde{\Gamma}_{\text{scalar}}^{\text{bt}}$, can also be generalized to the spinor loop case simply by an application of the replacement rule (1.10).

4. The gluon loop case

As was already mentioned, the case of the gluon loop is intrinsically more subtle, because here one has the issue of gauge (in) dependence not only for the background field but also for the loop particle. The preferred way of fixing the corresponding ambiguity for the three-vertex leads to the ‘‘gauge invariant vertex’’, which obeys the simple Ward identity (1.4). This version of the vertex is generated by the BFM with Feynman gauge for the quantum part, and it so happens that the only generalization of the worldline path integral representation (1.11) of the effective action to the gluon-loop case

presently known is just based on the BMF with quantum Feynman gauge, and thus the right starting point for a calculation of the “gauge invariant vertex”.

This representation was developed in [24, 45] in component fields, and reformulated in terms of worldline superfields in [44]. Again we will take a user’s approach here and proceed directly to the relevant master formula; the interested reader may consult [28] for more details. This master formula for the (color-ordered) contribution to the off-shell 1PI N -gluon amplitude due to a gluon loop reads

$$\begin{aligned}
\Gamma_{\text{gluon}}^{a_1 \dots a_N} [p_1, \varepsilon_1; \dots; p_N, \varepsilon_N] &= -\frac{(-ig)^N}{4} \text{tr}(T^{a_1} \dots T^{a_N}) \lim_{C \rightarrow \infty} \int_0^\infty \frac{dT}{T} (4\pi T)^{-\frac{D}{2}} e^{-CT} \\
&\times \prod_{k=1}^N \int_0^T d\tau_k \int d\theta_k \delta\left(\frac{\tau_N}{T}\right) \vartheta(\hat{\tau}_{1N}) \prod_{l=1}^{N-1} \vartheta(\hat{\tau}_{l(l+1)}) \sum_{p=P,A} \sigma_p Z_p \\
&\times \exp \left\{ \sum_{i,j=1}^N \left[\frac{1}{2} \hat{G}_{p,ij}^C p_i \cdot p_j + i D_i \hat{G}_{p,ij}^C \varepsilon_i \cdot p_j + \frac{1}{2} D_i D_j \hat{G}_{p,ij}^C \varepsilon_i \cdot \varepsilon_j \right] \right\} \Big|_{\text{lin}(\varepsilon_1 \dots \varepsilon_N)}.
\end{aligned} \tag{4.1}$$

Here the generators T^a are now fixed to be in the adjoint representation. We have defined $\sigma_P = 1$, $\sigma_A = -1$ (corresponding to periodic ($p = P$) and antiperiodic ($p = A$) boundary conditions in the original path integral), and

$$\begin{aligned}
Z_A &= (2 \cosh[CT/2])^4, \\
Z_P &= (2 \sinh[CT/2])^4,
\end{aligned} \tag{4.2}$$

$$\hat{G}_{P,A}^C(\tau_1, \theta_1; \tau_2, \theta_2) = G_B(\tau_1, \tau_2) + \theta_1 \theta_2 G_{P,A}^C(\tau_1, \tau_2), \tag{4.3}$$

where

$$\begin{aligned}
G_P^C(\tau_1, \tau_2) &= 2 \text{sign}(\tau_1 - \tau_2) \frac{\sinh[C(\frac{T}{2} - |\tau_1 - \tau_2|)]}{\sinh[CT/2]}, \\
G_A^C(\tau_1, \tau_2) &= 2 \text{sign}(\tau_1 - \tau_2) \frac{\cosh[C(\frac{T}{2} - |\tau_1 - \tau_2|)]}{\cosh[CT/2]}.
\end{aligned} \tag{4.4}$$

The limit $C \rightarrow \infty$ and sum $\sum_{p=P,A}$ serve the purpose to remove unwanted degrees of freedom circulating in the loop. Now, note that at fixed C, p the gluon loop master formula (4.1) is isomorphic to the spinor loop one (3.3). For our three-point case, this formal analogy implies that we can generalize (3.7) to

$$\Gamma_{\text{gluon}} = -\frac{1}{4} \frac{g^3}{(4\pi)^{D/2}} \lim_{C \rightarrow \infty} \sum_{p=P,A} \sigma_p \left(\Gamma_{\text{gluon}}^3(C, p) + \Gamma_{\text{gluon}}^2(C, p) + \Gamma_{\text{gluon}}^{\text{bt}}(C, p) \right) \quad (4.5)$$

where $\Gamma_{\text{gluon}}^{(\cdot)}(C, p)$ differs from the corresponding $\Gamma_{\text{spinor}}^{(\cdot)}$ in (3.8) only by a replacement of m^2 by C , G_{Fij} by $G_{p,ij}^C$, and the insertion of Z_p under the T integral.

It remains to analyze the limit $C \rightarrow \infty$ and the sum over boundary conditions; however, this has already been done in complete generality in [24, 45, 28]. For the three different types of terms appearing in (4.5), the general rules found there give

$$\begin{aligned} \lim_{C \rightarrow \infty} e^{-CT} \sum_{p=P,A} \sigma_p Z_p &= -8, \\ \lim_{C \rightarrow \infty} e^{-CT} \sum_{p=P,A} \sigma_p Z_p G_{p,ij}^C G_{p,ji}^C &= 16, \\ \lim_{C \rightarrow \infty} e^{-CT} \sum_{p=P,A} \sigma_p Z_p G_{p,12}^C G_{p,23}^C G_{p,31}^C &= 16. \end{aligned} \quad (4.6)$$

From (3.8), (4.5) and (4.6) we now get our final result for the gluon loop,

$$\Gamma_{\text{gluon}} = 2 \frac{g^3}{(4\pi)^{D/2}} (\Gamma_{\text{gluon}}^3 + \Gamma_{\text{gluon}}^2 + \Gamma_{\text{gluon}}^{\text{bt}}) \quad (4.7)$$

where, in terms of the spinor loop results of (3.8),

$$\begin{aligned} \Gamma_{\text{gluon}}^3 &= \Gamma_{\text{spinor}}^3(G_{F12} G_{F23} G_{F31} \rightarrow -2), \\ \Gamma_{\text{gluon}}^2 &= \Gamma_{\text{spinor}}^2(G_{F12} G_{F21} \rightarrow -2), \\ \Gamma_{\text{gluon}}^{\text{bt}} &= \Gamma_{\text{spinor}}^{\text{bt}}(G_{F12} G_{F21} \rightarrow -2). \end{aligned} \quad (4.8)$$

However, we must not forget the ghost loop contribution, which is necessary for the subtraction of the unphysical degrees of freedom of the gluon in the loop, and not contained in (4.1). It is equal to the scalar loop vertex (2.14), but has to be taken with the opposite sign:

$$\Gamma_{\text{ghost}} = -\Gamma_{\text{scalar}}. \quad (4.9)$$

Finally, in the gluon loop case, too, we have the option of using R_3^2 instead of Q_3^2 with an additional boundary contribution $\tilde{\Gamma}_{\text{gluon}}^{\text{bt}}$, and it is easy to check that this form of the result still relates to the corresponding one for the spinor loop result by (4.8).

5. Summary

We will now summarize our results for the scalar, spinor, and gluon loop cases. We will combine the three cases using a subscript $s = 0, \frac{1}{2}, 1$, where $\Gamma_0 = \Gamma_{\text{scalar}}$, $\Gamma_{\frac{1}{2}} = \Gamma_{\text{spinor}}$, $\Gamma_1 = \Gamma_{\text{gluon}} + \Gamma_{\text{ghost}}$. It will also be useful to rewrite the multiple τ_i - integrals in terms of the more standard Feynman/Schwinger parameter integrals. First, as usual we rescale $\tau_i = Tu_i, i = 1, 2, 3$, after which the T integral can be done trivially. Then we change from $u_1, u_2, u_3 = 0$ to $\alpha_1, \alpha_2, \alpha_3$ via

$$\begin{aligned} u_1 &= \alpha_2 + \alpha_3, \\ u_2 &= \alpha_3, \end{aligned} \quad (5.1)$$

with $\alpha_1 + \alpha_2 + \alpha_3 = 1$. Introducing the six parameter integrals

$$\begin{aligned} I_{3,B}^D(p_1^2, p_2^2, p_3^2) &= \int_0^1 d\alpha_1 d\alpha_2 d\alpha_3 \delta(1 - \alpha_1 - \alpha_2 - \alpha_3) \frac{(1 - 2\alpha_1)(1 - 2\alpha_2)(1 - 2\alpha_3)}{\left(m^2 + \alpha_1\alpha_2 p_1^2 + \alpha_2\alpha_3 p_2^2 + \alpha_1\alpha_3 p_3^2\right)^{3 - \frac{D}{2}}}, \\ I_{3,F}^D(p_1^2, p_2^2, p_3^2) &= - \int_0^1 d\alpha_1 d\alpha_2 d\alpha_3 \delta(1 - \alpha_1 - \alpha_2 - \alpha_3) \frac{1}{\left(m^2 + \alpha_1\alpha_2 p_1^2 + \alpha_2\alpha_3 p_2^2 + \alpha_1\alpha_3 p_3^2\right)^{3 - \frac{D}{2}}}, \\ I_{2,B}^D(p_1^2, p_2^2, p_3^2) &= \int_0^1 d\alpha_1 d\alpha_2 d\alpha_3 \delta(1 - \alpha_1 - \alpha_2 - \alpha_3) \frac{(1 - 2\alpha_2)^2(1 - 2\alpha_1)}{\left(m^2 + \alpha_1\alpha_2 p_1^2 + \alpha_2\alpha_3 p_2^2 + \alpha_1\alpha_3 p_3^2\right)^{3 - \frac{D}{2}}}, \end{aligned}$$

$$\begin{aligned}
I_{2,F}^D(p_1^2, p_2^2, p_3^2) &= \int_0^1 d\alpha_1 d\alpha_2 d\alpha_3 \delta(1 - \alpha_1 - \alpha_2 - \alpha_3) \frac{1 - 2\alpha_1}{\left(m^2 + \alpha_1 \alpha_2 p_1^2 + \alpha_2 \alpha_3 p_2^2 + \alpha_1 \alpha_3 p_3^2\right)^{3 - \frac{D}{2}}}, \\
I_{\text{bt},B}^D(p^2) &= \int_0^1 d\alpha \frac{(1 - 2\alpha)^2}{\left(m^2 + \alpha(1 - \alpha)p^2\right)^{2 - \frac{D}{2}}}, \\
I_{\text{bt},F}^D(p^2) &= \int_0^1 d\alpha \frac{1}{\left(m^2 + \alpha(1 - \alpha)p^2\right)^{2 - \frac{D}{2}}}
\end{aligned} \tag{5.2}$$

we can then write:

$$\Gamma_s = d_s \frac{g^3}{(4\pi)^{\frac{D}{2}}} (\Gamma_s^3 + \Gamma_s^2 + \Gamma_s^{\text{bt}}). \tag{5.3}$$

Here $d_0 = d_1 = 1, d_{\frac{1}{2}} = -2$ and

$$\begin{aligned}
\Gamma_0^3 &= -\Gamma \left(3 - \frac{D}{2}\right) \text{tr}(T^{a_1} T^{a_2} T^{a_3}) \text{tr}(f_1 f_2 f_3) I_{3,B}^D(p_1^2, p_2^2, p_3^2) + (a_2 \leftrightarrow a_3, f_2 \leftrightarrow f_3, p_2 \leftrightarrow p_3), \\
\Gamma_0^2 &= \frac{1}{2} \Gamma \left(3 - \frac{D}{2}\right) \text{tr}(T^{a_1} T^{a_2} T^{a_3}) \left[\text{tr}(f_1 f_2) (\varepsilon_3 \cdot p_1 I_{2,B}^D(p_1^2, p_2^2, p_3^2) - \varepsilon_3 \cdot p_2 I_{2,B}^D(p_2^2, p_1^2, p_3^2)) \right. \\
&\quad \left. + \text{tr}(f_2 f_3) (\varepsilon_1 \cdot p_2 I_{2,B}^D(p_2^2, p_3^2, p_1^2) - \varepsilon_1 \cdot p_3 I_{2,B}^D(p_3^2, p_2^2, p_1^2)) \right. \\
&\quad \left. + \text{tr}(f_3 f_1) (\varepsilon_2 \cdot p_3 I_{2,B}^D(p_3^2, p_1^2, p_2^2) - \varepsilon_2 \cdot p_1 I_{2,B}^D(p_1^2, p_3^2, p_2^2)) \right] \\
&\quad + (a_2 \leftrightarrow a_3, f_2 \leftrightarrow f_3, \varepsilon_2 \leftrightarrow \varepsilon_3, p_2 \leftrightarrow p_3), \\
\Gamma_0^{\text{bt}} &= \Gamma \left(2 - \frac{D}{2}\right) \text{tr}(T^{a_1} [T^{a_2}, T^{a_3}]) \left[\varepsilon_3 \cdot f_1 \cdot \varepsilon_2 I_{\text{bt},B}^D(p_1^2) + \varepsilon_1 \cdot f_2 \cdot \varepsilon_3 I_{\text{bt},B}^D(p_2^2) \right. \\
&\quad \left. + \varepsilon_2 \cdot f_3 \cdot \varepsilon_1 I_{\text{bt},B}^D(p_3^2) \right],
\end{aligned} \tag{5.4}$$

$$\begin{aligned}
\Gamma_{\frac{1}{2}}^3 &= -\Gamma \left(3 - \frac{D}{2}\right) \text{tr}(T^{a_1} T^{a_2} T^{a_3}) \text{tr}(f_1 f_2 f_3) (I_{3,B}^D(p_1^2, p_2^2, p_3^2) - I_{3,F}^D(p_1^2, p_2^2, p_3^2)) + (2 \leftrightarrow 3), \\
\Gamma_{\frac{1}{2}}^2 &= \frac{1}{2} \Gamma \left(3 - \frac{D}{2}\right) \text{tr}(T^{a_1} T^{a_2} T^{a_3}) \left[\text{tr}(f_1 f_2) \left(\varepsilon_3 \cdot p_1 (I_{2,B}^D(p_1^2, p_2^2, p_3^2) - I_{2,F}^D(p_1^2, p_2^2, p_3^2)) \right. \right. \\
&\quad \left. \left. - \varepsilon_3 \cdot p_2 (I_{2,B}^D(p_2^2, p_1^2, p_3^2) - I_{2,F}^D(p_2^2, p_1^2, p_3^2)) \right) + 2 \text{ perm.} \right]
\end{aligned}$$

$$\begin{aligned}
& +(2 \leftrightarrow 3), \\
\Gamma_{\frac{1}{2}}^{\text{bt}} &= \Gamma \left(2 - \frac{D}{2} \right) \text{tr}(T^{a_1}[T^{a_2}, T^{a_3}]) \left[\varepsilon_3 \cdot f_1 \cdot \varepsilon_2 (I_{\text{bt},B}^D(p_1^2) - I_{\text{bt},F}^D(p_1^2)) + 2 \text{ perm.} \right], \\
\end{aligned} \tag{5.5}$$

$$\begin{aligned}
\Gamma_1^3 &= -\Gamma \left(3 - \frac{D}{2} \right) \text{tr}(T^{a_1}T^{a_2}T^{a_3}) \text{tr}(f_1 f_2 f_3) (I_{3,B}^D(p_1^2, p_2^2, p_3^2) - 4I_{3,F}^D(p_1^2, p_2^2, p_3^2)) + (2 \leftrightarrow 3), \\
\Gamma_1^2 &= \frac{1}{2} \Gamma \left(3 - \frac{D}{2} \right) \text{tr}(T^{a_1}T^{a_2}T^{a_3}) \left[\text{tr}(f_1 f_2) \left(\varepsilon_3 \cdot p_1 (I_{2,B}^D(p_1^2, p_2^2, p_3^2) - 4I_{2,F}^D(p_1^2, p_2^2, p_3^2)) \right. \right. \\
&\quad \left. \left. - \varepsilon_3 \cdot p_2 (I_{2,B}^D(p_2^2, p_1^2, p_3^2) - 4I_{2,F}^D(p_2^2, p_1^2, p_3^2)) \right) + 2 \text{ perm.} \right] \\
& +(2 \leftrightarrow 3), \\
\Gamma_1^{\text{bt}} &= \Gamma \left(2 - \frac{D}{2} \right) \text{tr}(T^{a_1}[T^{a_2}, T^{a_3}]) \left[\varepsilon_3 \cdot f_1 \cdot \varepsilon_2 (I_{\text{bt},B}^D(p_1^2) - 4I_{\text{bt},F}^D(p_1^2)) + 2 \text{ perm.} \right]. \\
\end{aligned} \tag{5.6}$$

Note that the ‘‘loop replacement rules’’ now have assumed the following form:

$$\Gamma_{\frac{1}{2}}^{(\cdot)} = \Gamma_0^{(\cdot)} \left(I_{(\cdot),B}^D \rightarrow I_{(\cdot),B}^D - I_{(\cdot),F}^D \right), \tag{5.7}$$

$$\Gamma_1^{(\cdot)} = \Gamma_0^{(\cdot)} \left(I_{(\cdot),B}^D \rightarrow I_{(\cdot),B}^D - 4I_{(\cdot),F}^D \right). \tag{5.8}$$

Further, it can be easily checked that, for the bulk terms $\Gamma_{(\cdot)}^{3,2}$, the terms with the interchange $(2 \leftrightarrow 3)$ just provide the other half of a color commutator $[T^{a_2}, T^{a_3}]$, so that for them, as for the boundary terms, the color structure factors out in a $\text{tr}(T^{a_1}[T^{a_2}, T^{a_3}])$. Therefore we can now use

$$\text{tr}(T^{a_1}[T^{a_2}, T^{a_3}]) = iC(r) f^{a_1 a_2 a_3} \tag{5.9}$$

to get the expected proportionality to $f^{a_1 a_2 a_3}$. Thus we can write

$$\begin{aligned}
\Gamma_s^{a_1 a_2 a_3} &= d_s \frac{g^3}{(4\pi)^{\frac{D}{2}}} \text{tr}(T^{a_1}[T^{a_2}, T^{a_3}]) (\gamma_s^3 + \gamma_s^2 + \gamma_s^{\text{bt}}) \\
&= i f^{a_1 a_2 a_3} C(r) d_s \frac{g^3}{(4\pi)^{\frac{D}{2}}} (\gamma_s^3 + \gamma_s^2 + \gamma_s^{\text{bt}}) \\
\end{aligned} \tag{5.10}$$

with

$$\begin{aligned}
\gamma_0^3 &= -\Gamma\left(3 - \frac{D}{2}\right) \text{tr}(f_1 f_2 f_3) I_{3,B}^D(p_1^2, p_2^2, p_3^2), \\
\gamma_0^2 &= \frac{1}{2}\Gamma\left(3 - \frac{D}{2}\right) \left[\text{tr}(f_1 f_2) \left(\varepsilon_3 \cdot p_1 I_{2,B}^D(p_1^2, p_2^2, p_3^2) - \varepsilon_3 \cdot p_2 I_{2,B}^D(p_2^2, p_1^2, p_3^2) \right) \right. \\
&\quad \left. + \text{tr}(f_2 f_3) \left(\varepsilon_1 \cdot p_2 I_{2,B}^D(p_2^2, p_3^2, p_1^2) - \varepsilon_1 \cdot p_3 I_{2,B}^D(p_3^2, p_2^2, p_1^2) \right) \right. \\
&\quad \left. + \text{tr}(f_3 f_1) \left(\varepsilon_2 \cdot p_3 I_{2,B}^D(p_3^2, p_1^2, p_2^2) - \varepsilon_2 \cdot p_1 I_{2,B}^D(p_1^2, p_3^2, p_2^2) \right) \right], \\
\gamma_0^{\text{bt}} &= \Gamma\left(2 - \frac{D}{2}\right) \left[\varepsilon_3 \cdot f_1 \cdot \varepsilon_2 I_{\text{bt},B}^D(p_1^2) + \varepsilon_1 \cdot f_2 \cdot \varepsilon_3 I_{\text{bt},B}^D(p_2^2) + \varepsilon_2 \cdot f_3 \cdot \varepsilon_1 I_{\text{bt},B}^D(p_3^2) \right]
\end{aligned} \tag{5.11}$$

and the $\gamma_{\frac{1}{2},1}^{(\cdot)}$'s obtained from the $\gamma_0^{(\cdot)}$'s by the rule (5.7) resp. (5.8).

In the version where Q_3^2 is traded for R_3^2 , γ_0^2 gets replaced by

$$\tilde{\gamma}_0^2 = \frac{1}{2}\Gamma\left(3 - \frac{D}{2}\right) \left\{ \text{tr}(f_1 f_2) \left[\frac{r_3 \cdot f_3 \cdot p_1}{r_3 \cdot p_3} I_{2,B}^D(p_1^2, p_2^2, p_3^2) - \frac{r_3 \cdot f_3 \cdot p_2}{r_3 \cdot p_3} I_{2,B}^D(p_2^2, p_1^2, p_3^2) \right] + 2 \text{ perm.} \right\} \tag{5.12}$$

and one has the additional boundary contribution

$$\tilde{\gamma}_0^{\text{bt}} = -\frac{1}{2}\Gamma\left(2 - \frac{D}{2}\right) \left\{ \left[\text{tr}(f_1 f_2) \rho_3 - \text{tr}(f_3 f_1) \rho_2 \right] I_{\text{bt},B}^D(p_1^2) + 2 \text{ perm.} \right\}. \tag{5.13}$$

The rules (5.7) and (5.8) continue to hold.

6. Comparison with previous results

We now study the connection between our results for the three-gluon vertex and previous work. Since our treatment of the gluon loop case is equivalent to the use of the BFM with quantum Feynman gauge, we expect the Binger-Brodsky relation (1.5) to hold; and indeed this relation here follows immediately from the replacement rules (5.7) and (5.8).

For the same reason, the QED-like Ward identity (1.4) should be fulfilled not only for the scalar and spinor, but also for the gluon loop case.

Here it is advantageous to use the version where R_3^2 is used instead of Q_3^2 . Since R_3^2 is transversal, the Ward identity then involves only the boundary terms $\gamma_0^{\text{bt}}, \tilde{\gamma}_0^{\text{bt}}$, and can be easily verified using (5.11),(5.13) together with integral formulas given in [28] that express the three vacuum polarization tensors $\Pi_{0,\frac{1}{2},1}$ in terms of the integrals $I_{2,B}^D(p^2)$ and $I_{2,F}^D(p^2)$, with the same replacement rules (5.7),(5.8) holding:

$$\begin{aligned}
\Pi_0(p^2) &= -C(r) \frac{g^2}{(4\pi)^{D/2}} \Gamma\left(2 - \frac{D}{2}\right) I_{\text{bt},B}^D(p^2), \\
\Pi_{\frac{1}{2}}(p^2) &= -2\Pi_0(p^2) \left(I_{\text{bt},B}^D \rightarrow I_{\text{bt},B}^D - I_{\text{bt},F}^D \right), \\
\Pi_1(p^2) &= \Pi_0(p^2) \left(I_{\text{bt},B}^D \rightarrow I_{\text{bt},B}^D - 4I_{\text{bt},F}^D \right).
\end{aligned} \tag{6.1}$$

Next, we proceed to the less straightforward task of relating our representation to the Ball-Chiu decomposition. As usual we start with the scalar case. Comparing our final result (5.10),(5.11) with (1.2), (1.3) we first note that $T_H = \text{tr}(f_1 f_2 f_3)$. Thus we must identify

$$H(p_1^2, p_2^2, p_3^2) = C(r) \frac{d_0 g^2}{(4\pi)^{D/2}} \Gamma\left(3 - \frac{D}{2}\right) I_{3,B}^D(p_1^2, p_2^2, p_3^2) \tag{6.2}$$

which is indeed totally symmetric in its arguments.

Further, it is also easy to recognize the functions A and B functions as symmetric and antisymmetric combinations of the functions contained in γ_0^{bt} :

$$\begin{aligned}
A(p_1^2, p_2^2, p_3^2) &= -C(r) \frac{d_0 g^2}{2(4\pi)^{D/2}} \Gamma\left(2 - \frac{D}{2}\right) \left[I_{\text{bt},B}^D(p_1^2) + I_{\text{bt},B}^D(p_2^2) \right], \\
T_A &= \varepsilon_1 \cdot \varepsilon_2 (p_1 \cdot \varepsilon_3 - p_2 \cdot \varepsilon_3)
\end{aligned} \tag{6.3}$$

and

$$\begin{aligned}
B(p_1^2, p_2^2, p_3^2) &= -C(r) \frac{d_0 g^2}{2(4\pi)^{D/2}} \Gamma\left(2 - \frac{D}{2}\right) \left[I_{\text{bt},B}^D(p_1^2) - I_{\text{bt},B}^D(p_2^2) \right], \\
T_B &= \varepsilon_1 \cdot \varepsilon_2 (p_1 \cdot \varepsilon_3 + p_2 \cdot \varepsilon_3).
\end{aligned} \tag{6.4}$$

Coming to the structure F , the fact that T_F is transversal suggests that we should again use the transversal structure $\tilde{\gamma}_0^2$ rather than γ_0^2 ; the question is, how to choose the still undetermined vectors r_i ? By inspection one finds that, with the cyclic choice $r_1 = p_2 - p_3, r_2 = p_3 - p_1, r_3 = p_1 - p_2$, and using the antisymmetry of f_i , e.g. the first term in braces in (5.12) turns into

$$\text{tr}(f_1 f_2) \frac{p_1 \cdot f_3 \cdot p_2}{(p_1 - p_2) \cdot p_3} \left[I_{2,B}^D(p_1^2, p_2^2, p_3^2) - I_{2,B}^D(p_2^2, p_1^2, p_3^2) \right]. \quad (6.5)$$

Noting that

$$T_F = \frac{1}{2} \text{tr}(f_1 f_2) p_1 \cdot f_3 \cdot p_2 \quad (6.6)$$

we are led to set

$$F(p_1^2, p_2^2, p_3^2) = C(r) \frac{d_0 g^2}{(4\pi)^{D/2}} \Gamma\left(3 - \frac{D}{2}\right) \frac{I_{2,B}^D(p_1^2, p_2^2, p_3^2) - I_{2,B}^D(p_2^2, p_1^2, p_3^2)}{p_1^2 - p_2^2} \quad (6.7)$$

where we have also used momentum conservation to rewrite

$$(p_1 - p_2) \cdot p_3 = p_2^2 - p_1^2. \quad (6.8)$$

Thus the remaining structure C must match $\tilde{\gamma}_0^{\text{bt}}$, and indeed one has

$$T_C = \frac{1}{2} \text{tr}(f_1 f_2) (p_1 - p_2) \cdot \varepsilon_3 = -\frac{1}{2} \text{tr}(f_1 f_2) \rho_3 (p_1^2 - p_2^2) \quad (6.9)$$

leading to the identification

$$C(p_1^2, p_2^2, p_3^2) = -C(r) \frac{d_0 g^2}{(4\pi)^{D/2}} \Gamma\left(2 - \frac{D}{2}\right) \frac{I_{\text{bt},B}^D(p_1^2) - I_{\text{bt},B}^D(p_2^2)}{p_1^2 - p_2^2}. \quad (6.10)$$

Note that F and C are indeed symmetric functions in the first two arguments, and that C is actually independent of p_3^2 . Also, A is the only one of the functions having an UV divergence (since in B the expression in square brackets is $O(\epsilon)$), and B and C are simply related by [6]

$$2B = (p_1^2 - p_2^2)C. \quad (6.11)$$

Passing from the scalar to the spinor and gluon loop cases using (5.7) and (5.8) will obviously not change anything essential in this analysis.

For the (massive) spinor loop case we have also verified the above correspondences explicitly, using the formulas for the functions A to H given in [6] (to be precise, we have done this check for A, B, C with arbitrary momentum, for F specializing to $p_3^2 = 0$ and for H specializing to $p_1^2 = p_2^2 = 0$). This provides also a check on the much more involved calculations of [6].

7. Comparison with the effective action

It will be instructive to compare our results for the three-point amplitude with the low energy expansion of the one-loop QCD effective action induced by a loop particle of mass m . The general form of this expansion is

$$\Gamma_{\text{scalar}}[F] = \int_0^\infty \frac{dT}{T} \frac{e^{-m^2 T}}{(4\pi T)^{D/2}} \text{tr} \int dx_0 \sum_{n=2}^{\infty} \frac{(-T)^n}{n!} O_n[F], \quad (7.1)$$

where $O_n(F)$ is a Lorentz and gauge invariant expression of mass dimension $2n$. For the scalar loop, in [46, 47] this expansion was obtained to order $O(T^6)$. To see the relation with our form factor decomposition, it will be sufficient to consider the $n = 2$ and $n = 3$ terms:

$$\begin{aligned} O_2 &= -\frac{1}{6} g^2 F_{\mu\nu} F_{\mu\nu}, \\ O_3 &= -\frac{2}{15} i g^3 F_{\kappa\lambda} F_{\lambda\mu} F_{\mu\kappa} - \frac{1}{20} g^2 D_\lambda F_{\mu\nu} D^\lambda F^{\mu\nu}. \end{aligned} \quad (7.2)$$

Here changing from the scalar to the spinor or gluon loop will change only the coefficients in the expansion (7.1), not its structure. Comparing with, e.g., (5.11) we easily recognize the correspondences

$$\begin{aligned} \gamma_{(\cdot)}^3 &\leftrightarrow F_\kappa^\lambda F_\lambda^\mu F_\mu^\kappa = f_\kappa^\lambda f_\lambda^\mu f_\mu^\kappa + \text{higher point terms}, \\ \gamma_{(\cdot)}^2 &\leftrightarrow (\partial + ig \underbrace{A}) F (\partial + ig A) F, \\ \gamma_{(\cdot)}^{\text{bt}} &\leftrightarrow (f + ig \underbrace{[A, A]}) (f + ig [A, A]). \end{aligned} \quad (7.3)$$

Thus all the pieces of our form factor decomposition have a simple meaning in terms of the effective action. Note that commutator terms are always generated by boundary terms in the IBP, and that our three-point results allow us to predict certain terms in the higher-point gluon amplitudes using the knowledge that any “abelian” field strength tensor in the nonabelian effective action must appear as part of the full nonabelian field strength tensor including the commutator term. Note also that the tensor structure multiplying the function S in the Ball-Chiu decomposition does not correspond to anything in the expansion (7.1), which is one way of understanding why S turned out to be zero in the calculations of [4, 6]. Since the structure of the effective action is loop-independent, this observation allows us also to predict that the vanishing of S is not a one-loop accident, and will be found to persist at higher loop orders.

8. Conclusions

We have recalculated here the one-loop QCD three-gluon vertex for the scalar, fermion and gluon loop cases in a unifying way, achieving a compact result involving only six different parameter integrals. We have established the precise relation of this result to the standard Ball-Chiu decomposition, and also verified this relation for the massive spinor loop case using the explicit results of [6].

As was mentioned already in the introduction, even in a four-dimensional calculation the use of the vertex as a building block for higher loop calculations will in most cases make it necessary to know its D - dimensional continuation. For that reason we have kept the full D - dependence as much as possible. In fact, our result for the scalar loop is complete in this sense and holds for arbitrary D . And also in the spinor loop case the only place where we have used $D = 4$ is in the normalization of the path integral, which however corresponds to the usual fixing of $\text{tr}_\gamma \mathbb{1} = 4$ in dimensional regularization. It is only in the gluon loop case where $D = 4$ has been used in a nontrivial way, namely already in the derivation of (4.1). Here a true extension to other space-time dimensions would require some more work. For the purpose of dimensional regularization it is, however, sufficient to note that our result for the gluon loop case corresponds to the dimensional reduction variant of dimensional regularization proposed in [23].

In this calculation, three main advantages of our approach have emerged.

First, the IBP procedure generates the standard transversality-based Ball-Chiu decomposition of the vertex almost automatically, bypassing the usual tedious analysis of the nonabelian Ward identities. Let us recapitulate

how this happens: for the bulk terms, in the IBP all polarization vectors get absorbed into “abelian” field strength tensors $f_i^{\mu\nu}$, and thus become transversal. In the abelian case, there would be no boundary contributions, and one would have achieved manifest transversality at the integrand level. In the nonabelian case there are boundary terms, and those combine into commutator terms that carry all non-transversality, and generally contribute to the covariantization of some lower-point bulk term.

Second, this emergence of field strength tensors in the IBP allows one to maintain a close relation between the momentum space amplitudes and the low energy effective action, and thus to profit from the superior organization of the latter with respect to gauge invariance. This has led us to predict that the vanishing of the coefficient function S of the Ball-Chiu decomposition will be found to persist beyond one-loop.

Third, the spinor and gluon loop contributions can be obtained from the scalar loop one trivially using the off-shell extended Bern-Kosower “loop replacement rules” (5.7),(5.8) and. The gluon loop result corresponds to a field theory calculation in the BFM with quantum Feynman gauge, and thus to the preferred “gauge-invariant vertex” which fulfills the simple Ward identity (1.4) and the SUSY-related identity (1.5). The latter here appears as a simple consequence of the replacement rules and thus relates to worldline SUSY.

Concerning the last point, in the calculation presented here we have verified, rather than assumed, the validity of this off-shell extension. Had we taken the validity of the replacement rules off-shell for granted from the beginning, our method of calculating the three-point vertex would have been even much more efficient; in fact, incomparably more efficient than the combined effort of [4, 6, 7] that was necessary to arrive at an explicit result for the scalar, spinor and gluon loop contributions to the three-gluon vertex with standard field theory methods. Before applying it to higher-point vertices, it will thus be important to show the validity of this off-shell extension in general. With hindsight, this can be done as follows: it is sufficient to show the validity of the replacement rules for the effective action. Let us consider the spinor loop case first. Here it was shown in [28] that the replacement rule in the abelian case holds for the off-shell N - photon amplitudes. Thus it holds also for the abelian effective action. The nonabelian effective action in the low energy expansion can be decomposed into terms that are Lorentz scalars built from covariant derivatives and field strength tensors. Each such term in general will, after Fourier transformation, contribute to momentum space functions with various different numbers of legs; e.g., the term $\text{tr}(D_\mu F_{\alpha\beta} D^\mu F^{\alpha\beta})$ will contribute to the N - point functions with N

between two and six. Generally, each such term in the nonabelian effective Lagrangian has a “core” term, which has a counterpart already in the abelian case (in the example this would be $\partial_\mu f_{\alpha\beta} \partial^\mu f^{\alpha\beta}$) and a number of “covariantizing” terms that all involve commutators, and belong to amplitudes with more legs than the core term. For the core term the IBP leads from bulk term to bulk term and is formally identical to the abelian case, so that the replacement rule holds. But a core term in the effective action appears combined with all its covariantizing terms, all sharing the same coefficient. The replacement rule induces a change of this coefficient defined through a cycle $\dot{G}_{Bi_1i_2} \dot{G}_{Bi_2i_3} \cdots \dot{G}_{Bi_ni_1}$ which is multiplied by a $\text{tr}(f_{i_1} f_{i_2} \cdots f_{i_n})$ that in the effective action corresponds to a $\text{tr}(f^n)$. Consistency is therefore possible only if the same change of coefficient applies also to all the terms where one or several of the factors $f^{\mu\nu}$ are replaced by a $[A, A]$ term, or where a ∂ acting on some F is replaced by a $[A, F]$. This settles the spinor loop case. For the gluon loop case, it is sufficient to remember that the corresponding master formula before taking the limit $C \rightarrow \infty$ and the projector sum $\sum_{p=P,A}$, eq. (4.1), is still isomorphic to the spinor loop one (3.3).

Based on this general validity of the off-shell replacement rules and the general IBP algorithm developed in [32] we anticipate that with the method presented here a first explicit calculation of the four-point vertex should be well in reach.

Less straightforward but very interesting would be the extension of the formalism presented here to the gravitational case. The one-loop three-graviton vertices have already been extensively studied for their conformal properties (see [52] and refs. therein). As far as external gravitons are concerned, suitable string-inspired representations exist already for the (off-shell) one-loop N - graviton amplitudes with a scalar or spinor loop [48, 49] as well as for a photon loop [50, 51]. However, a suitable IBP procedure still remains to be developed.

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Appendix A. Summary of Conventions

We work with the $(-+++)$ metric. The non-abelian covariant derivative is $D_\mu \equiv \partial_\mu + igA_\mu^a T^a$, with $[T^a, T^b] = if^{abc}T^c$. The adjoint representation is given by $(T^a)^{bc} = -if^{abc}$. The normalization of the generators is $\text{tr}(T^a T^b) = C(r)\delta^{ab}$, where for $SU(N)$ one has $C(N) = \frac{1}{2}$ for the fundamental and $C(G) = N$ for the adjoint representation.

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