The treatment of the infrared region
in perturbative QCD

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Abstract

We discuss the contribution coming from the infrared region to NLO matrix elements and/or coefficient functions of hard QCD processes. Strictly speaking, this contribution is not known theoretically, since it is beyond perturbative QCD. For DGLAP evolution all the infrared contributions are collected in the phenomenological input parton distribution functions (PDFs), at some relatively low scale $Q_0$; functions which are obtained from a fit to the ‘global’ data. However dimensional regularization sometimes produces a non-zero result coming from the infrared region. Instead of this conventional regularization treatment, we argue that the proper procedure is to first subtract from the NLO matrix element the contribution already generated at the same order in $\alpha_s$ by the LO DGLAP splitting function convoluted with the LO matrix element. This prescription eliminates the logarithmic infrared divergence, giving a well-defined result which is consistent with the original idea that everything below $Q_0$ is collected in the PDF input. We quantify the difference between the proposed treatment and the conventional approach using low-mass Drell-Yan production and deep inelastic electron-proton scattering as examples; and discuss the potential impact on the ‘global’ PDF analyses. We present arguments to show that the difference cannot be regarded as simply the use of an alternative factorization scheme.
1 Introduction

While studying the perturbative QCD (pQCD) description\(^1\) of the Drell-Yan production of low-mass \(\mu^+\mu^-\) pairs, in preparation for the interpretation of the forthcoming LHC measurements\(^2\), we unearthed a puzzle in the conventional procedure to remove the infrared divergencies. The puzzle is not just confined to the description of Drell-Yan production, but occurs for other QCD processes, such as deep inelastic electron-proton scattering, etc. The resolution of the puzzle can lead to numerical corrections at low scales. As such, it has the potential to influence the global PDF analyses \([3]-[8]\).

To introduce the problem, recall that pQCD calculations are based on factorization theorems. The cross section of a hard process is calculated from the convolutions of parton distribution functions (PDFs) with the cross sections of the various hard subprocesses. The PDFs satisfy DGLAP evolution in \(\ln Q^2\) starting from an initial input at some low starting scale \(Q_0^2\). The input values are known from ‘global’ fits to all the available deep inelastic and related hard scattering data. Below \(Q_0\) the higher-order \(\alpha_s\) and higher-twist and other corrections, become too large for the use of pQCD.

In the collinear approximation, at leading order (LO), each additional power of small \(\alpha_s\) is compensated by the large logarithm of a hard scale (such as \(\ln Q^2\)). At next-to-leading order (NLO) we have an \(\alpha_s\) term in the hard matrix element and/or the DGLAP splitting function unaccompanied by a large log. At NNLO the \(\alpha_s^2\) terms do not have large logs, and so on.

There is no problem with this procedure at LO — due to strong \(k_t\)-ordering, when the DGLAP evolution is started from \(Q_0\), we do not encounter contributions from the \(k_t < Q_0\) domain. On the other hand, already inside the NLO cell there appears a contribution with \(k_t < Q_0\). How to treat the (infrared) contribution from this large distance domain is the subject of this note.

We discuss two alternative treatments of the infrared domain. First, we adopt a practical approach, which we shall call the ‘physical’ prescription, see Section 2. To avoid double counting, we have to subtract from the NLO diagram the \(\alpha_s\log\) contribution that has already been generated by LO DGLAP evolution. As we will show, this completely removes the infrared divergency from the matrix element. To be precise, after the subtraction, there remains a very small contribution coming from the \(k_t < Q_0\) domain of \(\mathcal{O}(Q_0^2/Q_0^2)\), where \(\mu_F\) is the factorization scale, which we will discuss in Section 2.

The second treatment of the infrared divergency, which we will call the ‘conventional’ prescription, is based on dimensional regularisation, see Section 3. Here just the \(1/\epsilon\) pole, which corresponds to the anomalous dimension generated by LO DGLAP evolution, is subtracted from the NLO contribution. In this case, after the subtraction, we are left with a contribution from the \(k_t < Q_0\) region which does not vanish as \(Q_0^2 \rightarrow 0\). So, surprisingly, the two treatments of the infrared domain yield different results.

\(^1\)A preliminary study can be found in \([1]\).
\(^2\)See, for example, the preliminary LHCB data \([2]\).
Sections 2 and 3 introduce the two treatments in more detail, using the NLO coefficient function of Drell-Yan production as an example. Then in Section 4 we quantify the difference of the contributions from the infrared domain taking again the NLO coefficient function in Drell-Yan production and, in addition, the $\gamma^* g$ coefficient function, $C_g$, of deep inelastic scattering, as topical examples. Section 5 (as well as the latter part of Section 4) discusses the origin of the difference.

## 2 Calculation at NLO – the physical approach

Schematically, we may write the cross section in the form

$$
\frac{d\sigma}{d^3p} = \int dx_1 dx_2 \text{PDF}(x_1, \mu_F) |\mathcal{M}(p; \mu_F, \mu_R)|^2 \text{PDF}(x_2, \mu_F),
$$

(1)

where a sum over the various pairs of PDFs is implied. The matrix elements squared, $|\mathcal{M}|^2$, describe the cross sections of the elementary partonic subprocesses. However, the problem is that using (1) we do not know the factorization scale $\mu_F$ at which the PDFs are measured. Moreover, in the low $x$ region, the PDFs strongly depend on the choice of $\mu_F$.

In general, after the summation of all perturbative orders, the final result should not depend on the choice of $\mu_F$ that is used to separate the incoming PDFs from the hard matrix element. Thus we need to account for the NLO, NNLO,...corrections. Contributions with low virtuality, $q^2 < \mu_F^2$, of the incoming partons are included in the PDFs, while those with $q^2 > \mu_F^2$ are assigned to the matrix element.

Let us start with the LO expression for the cross section. In the collinear approach, the cross section has the form

$$
\sigma(\mu_F) = \text{PDF}(\mu_F) \otimes C_{\text{LO}} \otimes \text{PDF}(\mu_F),
$$

(2)

The effect of varying the scale from $m$ to $\mu_F$, in both the left and right PDFs, can be expressed, to first order in $\alpha_s$, as

$$
\sigma(\mu_F) = \text{PDF}(m) \otimes \left( C_{\text{LO}} + \frac{\alpha_s}{2\pi} \ln \left( \frac{\mu_F^2}{m^2} \right) (P_{\text{left}} C_{\text{LO}} + C_{\text{LO}} P_{\text{right}}) \right) \otimes \text{PDF}(m),
$$

(3)

where the splitting functions $P_{\text{right}}$ and $P_{\text{left}}$ act on the right and left PDFs respectively. Recall that in calculating the $\alpha_s$ correction in (3), the integral over the transverse momentum (virtuality) of the parton in the LO DGLAP evolution was approximated by the pure logarithmic $dk^2/k^2$ form. That is, in the collinear approach, the Leading Log Approximation (LLA) is used.

Now we turn to the expression for the cross section at NLO. First, we note that the original Feynman integrals corresponding to the NLO matrix element, $C_{\text{NLO}}$, do not depend on $\mu_F$. However, we will see below how a scale dependence enters. At NLO we may write

$$
\sigma(\mu_F) = \text{PDF}(\mu_F) \otimes (C_{\text{LO}} + \alpha_s C_{\text{corr}}) \otimes \text{PDF}(\mu_F),
$$

(4)
where we include the NLO correction to the coefficient function. However to avoid double counting we must subtract the DGLAP-generated $\alpha_s$ term of (3). This cancels the infrared singularity, and gives a non-singular result which can be safely integrated in normal 4-dimensional space.

2.1 Example: Drell-Yan production at NLO

We illustrate the general procedure in terms of the specific example of Drell-Yan production of a $\mu^+\mu^-$ pair of mass $M$. At LO the subprocess is $q\bar{q} \rightarrow \gamma^*$, while at NLO we have the $2 \rightarrow 2$ subprocesses $q\bar{q} \rightarrow g\gamma^*$ and $gq \rightarrow q\gamma^*$ shown in diagrams (a,b) of Fig. 1. These NLO contributions are now calculated with better, than LLA, accuracy. However part of these contributions are already included, to LLA accuracy, in the second term in (3), where the splitting functions

$$P_{\text{right}} = P_{qq} + P_{qg}, \quad P_{\text{left}} = P_{\bar{q}q} + P_{\bar{q}g},$$

act on the right and left PDFs respectively. So this LLA part should be subtracted from $C_{\text{NLO}}$.

For Drell-Yan production at low $x$, the majority of quarks/antiquarks are produced via the low-$x$ gluon-to-quark splitting, $P_{qq}$. That is, the most important NLO subprocess is $gq \rightarrow q\gamma^*$. The corresponding cross section, for the production of a $\mu^+\mu^-$ pair of mass $M$ and rapidity $Y$, has the schematic form

$$\frac{d\sigma}{dM^2 dy dt} = \int dx_1 dx_2 \, \text{PDF}_i(x_1) \frac{d\hat{\sigma}_{ij}}{dt} \, \text{PDF}_j(x_2) \, \delta(x_1^\gamma x_2^\gamma s - M^2) \, \delta \left( \frac{1}{2} \ln \frac{x_1^\gamma}{x_2^\gamma} - Y \right),$$

where $x_1^\gamma = x_i$ in the LO case, but $x_i^\gamma$ are the true momentum fractions ($x^+$ and $x^-$) carried by the photon at NLO. The subprocess cross sections have the form

$$\hat{\sigma}_{ij} = \hat{\sigma}_{ij}^{\text{LO}} + \hat{\sigma}_{ij}^{\text{NLO}} + \ldots,$$

3 We may equally well have incoming $\bar{q}$'s in $P_{\text{right}}$ and incoming $q$'s in $P_{\text{left}}$. 

Figure 1: Diagrams (a,b) show NLO subprocesses for Drell-Yan production resulting from the splitting of the upper PDF (or ‘right’ PDF, in the notation of (3)).
where, for the main NLO subprocess, we have

$$\frac{d\hat{\sigma}_{qg}}{dt} = \frac{1}{9} \frac{\alpha_s^2}{\hat{s}^2} \left[ -\frac{s}{t} - \frac{t}{\hat{s}} - \frac{2M^2u}{\hat{s}t} \right].$$

(8)

Since $M^2 = z\hat{s}$ and $u = M^2 - \hat{s} - t$, this gives

$$\frac{d\hat{\sigma}(gg \rightarrow q\gamma^*)}{d|t|} = \frac{\alpha_s^2 z}{9M^2 |t|} \left[ ((1-z)^2 + z^2) + \frac{z^2 t^2}{M^4} - 2z^2 \frac{t}{M^2} \right].$$

(9)

In order to calculate the inclusive cross section $d\sigma/dM^2$, it seems that we have to integrate over $t$ starting from $t = 0$. If this were necessary, then we would face an infrared divergency.

2.2 Treatment of the infrared region in the physical approach

When dealing with PDFs we avoid the problems of confinement and interactions at large distances. We start with some phenomenological input at a relatively large $Q_0$, that is $Q_0^2 \gg \Lambda_{\text{QCD}}^2$, and consider just the evolution of the PDFs with $Q^2$ increasing from $Q_0^2$ to the factorization scale $\mu_F^2$. Everything below $Q_0$ is absorbed in the phenomenological input PDF. Accepting this logic, it is natural to replace the lower limit $t = 0$ by the same $Q_0^2$, and never to consider the contribution from low virtualities, $k^2 < Q_0^2$.

The subtraction of the LO DGLAP contribution from the NLO $d\hat{\sigma}/dt$ completely eliminates the $1/t$ singularity as $t \to 0$ by introducing a $\Theta(|t| - \mu_F^2)$ function in the first term of (9). Let us explain, more explicitly, how this happens. We denote the $gg \rightarrow q\gamma^* \mathcal{O}(\alpha_s)$ contribution by $\hat{\sigma}^{(1)}$. The result of the explicit calculation of the $gg \rightarrow q\gamma^*$ cross section is shown in (9). On the other hand, this contribution may be written as the term generated by LO DGLAP evolution and the remaining NLO part, that is

$$d\hat{\sigma}^{(1)} = d\hat{\sigma}_{\text{rem}}^{\text{NLO}} + \sigma_{qg}^{\text{LO}} \otimes \frac{\alpha_s}{2\pi} P_{qg}^{\text{LO}} dt.$$

(10)

The last term reads

$$\sigma_{qg}^{\text{LO}} \otimes \frac{\alpha_s}{2\pi} P_{qg}^{\text{LO}} dt = \frac{\alpha_s^2 z}{9M^2} \frac{dt}{|t|} \frac{(1-z)^2 + z^2}{t} \Theta(\mu_F^2 - |t|),$$

(11)

where here DGLAP evolution has accounted for all virtualities $|t| = k^2 < \mu_F^2$; and where the contribution of $k^2 < Q_0^2$ is hidden in the phenomenological input PDF. After the subtraction of this LO DGLAP generated term, the remaining contribution of (9) is

$$d\hat{\sigma}_{\text{rem}}^{\text{NLO}} \frac{dt}{|t|} = \frac{\alpha_s^2 z}{9M^2} \frac{dt}{|t|} \left[ (1-z)^2 + z^2 \Theta(|t| - \mu_F^2) + \frac{z^2 t^2}{M^4} - 2z^2 \frac{t}{M^2} \right].$$

(12)

which has no singularity as $t \to 0$.

Recall that we need to subtract the DGLAP generated contribution of (11) in order to avoid the double counting. This subtraction is unambiguously defined. It is done in terms of $t$ (and
not $k_t$) since the original DGLAP evolution was written as an evolution in $k^2 = -t$, and not as an evolution in $k_t$.

A small uncertainty is still present coming from the treatment of the non-singular terms in \([12]\). We return to the choice of the lower limit of integration. It is not evident whether we should integrate from $t = 0$ or from $|t| = Q_0^2$. From the ladder diagram, Fig. 2(a), which is of exactly the same form as that in LO DGLAP evolution in which everything below $Q_0^2$ is absorbed in the input, we are led to integrate from $|t| = Q_0^2$. However, the NLO hard subprocess cross section also includes non-ladder diagrams, Fig. 2(b) with an $s$-channel quark, for which the $|t| < Q_0^2$ domain does not correspond to low virtuality. Thus the true uncertainty due to the non-singular terms is $\mathcal{O}(Q_0^2/\mu_F^2)$, which may be neglected for $\mu_F \gg Q_0^2$.

![Ladder and non-ladder diagrams](image)

**Figure 2:** (a) Ladder and (b) non-ladder diagrams.

### 3 Calculation at NLO – conventional approach

We now consider the conventional prescription for the evaluation of

$$
\frac{d\hat{\sigma}(gq \to q\gamma^*)}{d|t|} = \frac{\alpha_s^2\alpha_s}{9M^2} \frac{1}{|t|} \left[ \left((1-z)^2 + z^2\right) + z^2 \frac{t^2}{M^4} - 2z^2 \frac{t}{M^2} \right].
$$

Recall, that to calculate the inclusive cross section $d\sigma/dM^2$, we appear to have to integrate over $t$ starting from $t = 0$, and that we face an infrared divergency\(^4\). In the conventional approach the integral is regularized by introducing the $4+2\epsilon$ dimensional space \([9, 10, 11]\). Then the contribution from very small $t$ produces a $1/\epsilon$ pole, which is absorbed into the incoming PDF. However, in $4+2\epsilon$ dimensional space, the number of gluon states with transverse polarisation is $2 + 2\epsilon$, giving $\epsilon$ in the numerator of the matrix element. Besides this, there is an $\epsilon$ dependence coming from the decomposition of the phase space factor, like $(1 - z)^{-\epsilon}$. So finally we have an $\epsilon/\epsilon$ term, which produces a non-zero result as $\epsilon \to 0$.

\(^4\)In perturbation theory this divergency is avoided by the (small) current quark mass, $m_q$. However, in practice, the light quark mass is neglected.
Simultaneously, we have to consider the same diagram generated by LO DGLAP evolution, which gives a $1/\epsilon$ pole. The $1/\epsilon$ poles cancel, but we are left with the $\epsilon/\epsilon$ term, which produces a non-zero result as $\epsilon \to 0$. Unlike the physical approach, this non-zero term does not vanish as $O(Q_0^2/\mu_F^2)$, and does not vanish for $Q_0^2 \ll \mu_F^2$.

Thus, surprisingly, we find that the physical and conventional prescriptions yield different results. We quantify this difference in the next Section, and then in Section 5 we discuss its origin.

Sometimes the exact treatment of the infrared singularity does not matter. In particular, the emission of a soft gluon has a logarithmic infrared divergence which, according to the Bloch-Nordsieck theorem, is exactly cancelled between the real emission contribution and the virtual loop correction. However, in NLO Drell-Yan or in the $\gamma g$ coefficient function for the structure function $F_2$ of deep inelastic scattering, for example, the difference in the treatment of the infrared singularity does matter.

4 Quantitative estimates

To illustrate the numerical differences of the two treatments of the infrared contribution, we study, as examples, Drell-Yan production and the structure function $F_2$ of deep inelastic scattering.

4.1 Drell-Yan

The Drell-Yan cross section calculated using the Vrap code [12] and MSTW08 NLO PDFs [3] is shown by the continuous curve in Fig. 3 for the LHC energy $\sqrt{s} = 7$ TeV and a mass $M = 6$ GeV of the produced $\mu^+\mu^-$ pair. As usual, the factorization scale is chosen to be $\mu_F = M$. The long-dashed line is the remaining contribution of the NLO $gq \rightarrow q\gamma^*$ subprocess calculated in the framework of dimensional regularisation. We see that it turns out to be negative, since, for $\mu_F = M$, the contribution formally generated by LO evolution, (11), is larger than the whole $O(\alpha_s)$ cross section.

On the other hand, if we were to adopt the ‘physical’ approach and subtract the explicit DGLAP-generated term (as discussed in Section 3), then we obtain the result shown by the short-dashed curve, which is significantly different to that obtained using dimensional regularisation. Moreover, we see the uncertainty coming from the $|t| < Q_0^2$ domain is practically invisible in this figure — essentially identical results are obtained by integrating the last two (non-singular) terms in [...] in (12) with a lower limit $|t| = Q_0^2 = 1$ GeV$^2$ (dotted curve) and with $t = 0$ (short-dashed curve).

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5 This is the justification for the $+_{\text{prescription}}$ used for the $1/(1 - z)$ terms in the DGLAP splitting functions.
4.2 Deep inelastic scattering

An analogous situation is shown in Fig. 4 and 5 for the NLO gluon contribution to $F_2$ in DIS. The calculation is very similar to that for the Drell-Yan case since these two subprocesses are closely related by crossing symmetry. As before, the explicit subtraction of the term generated by the LO DGLAP evolution leads to a result for the $\gamma^*g$ coefficient function,

$$C_g = T_R \left( [(1-z)^2 + z^2] \ln \frac{1}{z} + 6z(1-z) - 1 \right),$$

that is different to that obtained by the conventional prescription,

$$C_g = T_R \left( [(1-z)^2 + z^2] \ln \frac{1-z}{z} + 8z(1-z) - 1 \right).$$

Here the notation of [11] has been used. The additional $\ln(1-z)$ singularity reflects the threshold singularity; as $z \to 1$ the phase space available for the quark and antiquark is proportional to $(1-z)$. In $\epsilon$ regularization this threshold singularity produces the factor

$$(1-z)^{-\epsilon} = 1 - \epsilon \ln(1-z) + O(\epsilon^2),$$

that is different to that obtained by the conventional prescription,

$$C_g = T_R \left( [(1-z)^2 + z^2] \ln \frac{1}{z} + 6z(1-z) - 1 \right).$$

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$$(1-z)^{-\epsilon} = 1 - \epsilon \ln(1-z) + O(\epsilon^2),$$

that is different to that obtained by the conventional prescription,
which, on being multiplied by the $1/\epsilon$ pole, gives a constant $\ln(1-z)$ term, that does not vanish as $\epsilon \to 0$. Unlike the conventional treatment, the physical prescription has, at most, only a very small $\mathcal{O}(Q^2_0/\mu_F^2)$ infrared contribution coming from the non-singular terms in the infrared region.

To gain insight into the difference between the ‘conventional’ and ‘physical’ treatments of the infrared divergency, we discuss the $t$-dependence of the $\gamma^*g \to q\bar{q}$ matrix element. Adopting the ‘physical’ procedure we start with the explicit perturbative QCD result and subtract the LO DGLAP-generated contribution. The result, after the photon polarizations have been convoluted with $g_{\mu\nu}$, may be written in the form

$$|M|^2 = |M_t|^2 + |M_u|^2,$$

where

$$|M_t|^2 = [(1-z)^2 + z^2]Q^2_{z t} \Theta(|t| - \mu_F^2) - 1.$$  

Similarly for $M_u$ with $u$ interchanged with $t$. The result is non-singular as $t \to 0$. Numerically

\[\text{Integrating (18) over } t, \text{ and adding the term which originates from } F_L \text{ (where there is no infrared divergency), leads to (14)}.\]
Figure 5: The fractional correction to $F_2$ arising from using the ‘physical’, rather the ‘conventional’, treatment of the infrared divergency in the $\gamma^*g$ coefficient function, shown as function of $Q^2$ for two values of $x$, and for two different sets of NLO PDFs, namely MSTW08 [3] and CT10 [4].

the non-singular contribution coming from the $0 \leq |t| \leq Q_0^2$ domain is not important. After integration over $t$, this contribution vanishes as $O(Q_0^2/\mu_F^2)$ as $Q_0 \to 0$. Indeed, the ‘physical’ prescription eliminates the logarithmic infrared divergence, giving a well-defined result which is consistent with the original idea that everything below $Q_0$ is collected in the PDF input.

On the other hand, the $1/t$ singularity in $|\mathcal{M}_t|^2$, in the ‘conventional’ treatment, is first regularized by working in $4+2\epsilon$ space. After the subtraction of the analogous DGLAP-initiated contribution, the $1/\epsilon$ term is cancelled. However, the resulting $\infty - \infty$ subtraction leaves some non-zero remaining contribution, which can be viewed as effective non-local vertices, which act as a $\delta(t)$ contribution. Recall that analogous vertices arising from an ultra-violet divergence, play the role of local counter terms. Dimensional regularization provides us with the possibility to introduce such counter terms in a way which does not spoil the renormalizability of the theory. However, the analogous effective vertices of infrared-origin are unacceptable. They are contrary to the confinement property of QCD. Due to the confinement property it is not possible to continue the $1/t$ behaviour below some low scale of $O(\Lambda_{QCD})$. Recall that in the global PDF analyses this low $t$ contribution is included phenomenologically in the input PDFs at the starting scale $Q_0$, which is taken to be larger than $\Lambda_{QCD}$.

\footnote{For the analogous Drell-Yan process we saw in Fig. 3 that the difference between the results obtained by integrating over the regions $|t| > 0$ and $|t| > Q_0^2 = 1$ GeV$^2$ were invisible on the plot. The same insensitivity of the results, when using the ‘physical’ treatment of infrared domain, applies also to the short-dashed curve in the $F_2$ plot of Fig 4.}
5 Discussion

Our approach has some similarities with the ‘subtraction method’ [13, 14] where to regularize
the infrared divergency one subtracts from the real cross section of \( n + 1 \) parton production,
\( d\sigma^{R}_{n+1} \), a term \( d\sigma^{A} \), which has exactly the same divergency. Simultaneously the \( d\sigma^{A} \) term is
added to the virtual loop correction for the \( n \) parton cross section, \( d\sigma^{V}_{n} \). After this both the
real and the virtual contribution have no infrared divergency, and so may be calculated in the
normal 4 dimensional space (without the \( \epsilon \) decomposition).

In comparison with this ‘subtraction method’, where the subtraction term is introduced
artificially, here we subtract the contribution \( \text{really} \) generated by DGLAP evolution. On one
hand, this guarantees the absence of double counting, while on the other hand, it eliminates
the infrared divergency. Note, also, that in our particular examples, like the NLO correction
to \( F_{2} \) due to photon-gluon fusion, or the NLO correction to the Drell-Yan cross section due to
\( gq \rightarrow q\gamma^{*} \) subprocess, there is no analogous LO contribution caused by gluons. Therefore there
is no NLO virtual correction, like \( d\sigma^{V}_{n} \), in the ‘subtraction method’.

Let us return to the comparison of the two treatments of the infrared domain discussed
in this paper: that is the ‘conventional’ \( \epsilon \) regularisation subtraction and the ‘physical’ direct
DGLAP subtraction methods. Why do the results differ? The reason is subtle. In the ‘physical’
approach we do not need to assume that perturbative QCD is valid for virtualities \( k^{2} < Q_{0}^{2} \),
and, in particular, in the confinement domain. We evolve starting from some phenomenological
‘input’ distribution at \( \mu^{2} = Q_{0}^{2} \). By considering different choices of \( Q_{0}^{2} \) we verify that the
contribution coming from the \( k^{2} < Q_{0}^{2} \) domain is small, and vanishes as \( Q_{0}^{2} \rightarrow 0 \). On the
other hand, the conventional dimensional \( \epsilon \) regularisation procedure continues the \( 1/t \) singular
behaviour below \( \Lambda_{QCD}^{2} \), giving finally a non-vanishing infrared contribution, which may be
viewed as a \( \delta(t) \) term during the \( t \) integration. However, the part coming from low virtualities,
\( k^{2} < Q_{0}^{2} \), has no reason to have perturbative form, and, moreover, does not vanish as \( Q_{0}^{2} \rightarrow 0 \).

We elaborate this in a little more detail. Dimensional regularization is clearly a valuable
tool in the ultraviolet region. But, here, we are concerned about its use in the infrared domain.
The hope was that for infrared stable quantities (like \( F_{2} \) the result will not depend on the
contribution from the large distance domain; the real and virtual loop contributions will cancel
each other. Sometimes this is true. In particular, due to the Bloch-Nordsieck theorem, we
can use the so-called ‘plus prescription’ which eliminates the \( 1/(1 - z) \) singularity in splitting
functions.

However, there are examples where ‘conventional’ dimensional regularization gives a non-
zero infrared contribution of the form of \( \epsilon/\epsilon \), which is the correct mathematical result of a
particular perturbative QCD diagram calculation, but which makes no physical sense for real
QCD with confinement. Here we have presented two examples (the NLO \( \gamma^{*}g \) coefficient function
in Drell-Yan production and \( C_{g} \) in DIS) and have shown that the subtraction of the LO DGLAP-
generated contribution completely eliminates the infrared singularity; the cancellation occurs
between terms in \( C^{NLO} \) and in the convolution the \( C^{LO} \otimes P^{LO} \). Recall that this subtraction
is necessary to avoid double counting. After this, the integral can be calculated in normal
4-dimensional space, and, surprisingly, as demonstrated above, the results differ from those
found in the ‘conventional’ approach.

Similarly, using the conventional $\epsilon$ regularisation procedure, the NLO splitting functions,
in particular $P_{gg}^{NLO}(z)$, of DGLAP evolution may also contain contributions of infrared origin
which are inconsistent with confinement. On the other hand, in the ‘physical’ approach, the
absence of the infrared singularity is due to the cancellation of the contribution to the NLO
splitting function arising from the direct calculation of the NLO Feynman diagram and the
contribution already generated by LO DGLAP evolution. That is, it is due to cancellations of
the contributions occurring in $P^{NLO}$ and in $P^{LO} \otimes P^{LO}$. A more detailed discussion is given in
the Appendix. The derivation of these new infrared-corrected NLO splitting functions is more
complicated than the treatment of the coefficient functions that we have used as examples, but
it is still possible.\footnote{We are at present tackling this problem.}

In summary, the problem cases for the ‘conventional’ treatment of the infrared divergence are
the NLO coefficient and splitting functions, in particular the $C_{g}$ and $P_{gg}$ functions, which leave
‘erroneous’ non-singular contributions. The main effect will be on the behaviour of the gluon at
low scales, a domain where the present global parton analyses show unexpected ‘valence-like’
$x$ behaviour of the gluon; see, for example, \cite{3, 4, 5}. In order to carry about such a ‘global’
investigation it will be necessary to compute, not only the coefficient functions, but also the
NLO splitting functions using the ‘physical’ treatment of the infrared domain.

One might think, say, in the DIS example that we considered, that the infrared corrections to
the coefficient and splitting functions could compensate each other, and that the transition from
the ‘conventional’ coefficient function (which includes a non-vanishing infrared contribution
inconsistent with colour confinement) to that obtained in a ‘consistent’ treatment, may be
effectively accounted for by a re-definition of the PDFs, and thus considered as an alternative
‘factorization scheme’. However such an alternative scheme, which may correspond to new
coefficient functions, prescribes a new set of splitting functions to describe the evolution of the
re-defined PDFs, which are, in general, different to the infrared-corrected splitting functions.

Even without explicit knowledge of the NLO splitting functions corresponding to the ‘physical’
treatment of the infrared domain, there are several ways to see that compensation is not
possible and that the results cannot be regarded as simply a scheme change. A hint, that this
is so, is the presence of the additional $\ln(1 - z)$ singularity in the ‘conventional’ $C_{g}$ of (15).
An argument is to consider the non-singlet channel. Here, due to flavour conservation, the
PDF normalization is fixed, and we have no possibility to re-define the non-singlet PDF via
the admixture with some gluon or other quark flavour contribution. In other words, there is
no chance to compensate the correction to NLO coefficient function by the re-definition of the
(non-singlet) parton density and/or by the correction to the NLO splitting function, which is
exponentiated, that is iterated many times already at NLO, especially at small $x$. It is not a
NNLO effect.
Alternatively, to be more explicit, we may consider the quark contribution to the $gg$ splitting function. This is very similar to the coefficient function of $\gamma^* g$ fusion. The part of the splitting function of (inconsistent) infrared origin gives a contribution, $\Delta \gamma$, to the anomalous dimensions, $\gamma$, — it therefore exponentiates during the evolution, which is especially long in the low $x$ region. In terms of the anomalous dimensions, the inconsistency $\Delta \gamma$ is strongly enhanced in the low $x$ region by the small value of the analytic continuation of the Mellin moment $\omega$, recall $\Delta \gamma \sim \text{const.}/\omega$. Compensation is clearly not possible over a long interval of $Q^2$.

Another good, related, example is evolution in QED (where there is no confinement) and where the electron density is well defined. Here the evolution was first considered by Gribov and Lipatov \cite{15} in terms of the Leading Log Approximation (rather than the ‘factorization theorem’ approach), which allows one to trace each step of the evolution and the contribution of each Feynman diagram explicitly. As there is no confinement, the electron and photon fields are well defined. Thus the only possibility to re-define the ‘partons’ is a field renormalization $Z(q^2)$ factor which does not depend on the momentum fraction $z$. Of course, one may consider some mixture of electron and photon PDFs, but we discuss the original electron and photons. For these distributions the only way to avoid double counting is the ‘physical’ approach. We emphasize that it is useful to first consider the problem of double counting in QED where everything is well defined. After this it is straightforward to discuss what happens in QCD.

A more general observation is that in our approach we work with partons defined by OPE operators, and the corresponding Feynman diagrams. For these quarks and gluons we evaluate the Feynman diagrams in both the ‘conventional’ and the ‘physical’ approach. In either case, we first have to calculate the splitting and the coefficient functions for the normal (OPE) quarks and gluons (for which we have Feynman rules and where each line in the diagrams has definite quantum numbers). Only after this may we define the new $q'$ and $g'$, and new splitting (and coefficient) functions corresponding to an alternative factorization scheme. So differences in the infrared treatments cannot be attributed to a scheme change.

We conclude that differences in parton behaviour that follow from an analysis using a ‘physical’ treatment of the infrared domain cannot be reproduced by a factorization scheme transition from the conventional PDF analyses. (The differences are expected to be mainly in the gluon at low scales.) The only way to account for the appropriate treatment of the infrared region is to perform a new global analysis using a complete set of corrected coefficient and splitting functions.

**Appendix: Infrared contributions to NLO splitting functions**

In the axial gauge the only infrared singularity of a splitting function is due to the ladder-type (box) diagram (plus a self-energy contribution $\propto \delta(1 - z) \int_0^1 dz P(z)$ which can be determined from this diagram in usual way based on the flavour and energy-momentum conservation laws).
The singularity has a logarithmic form $dk^2/k^2$ and is exactly equal to the contribution $P^{\text{LO}} \otimes P^{\text{LO}}$ generated by LO DGLAP evolution, since there is no other dimensionful parameter apart from $Q^2$. To avoid double counting, we have to subtract the LO DGLAP-generated contribution from the NLO splitting $P^{\text{NLO}}$. This subtraction eliminates the infrared singularity of $P^{\text{NLO}}$. However the remaining part may still contain some non-singular contribution of infrared (large distance) origin.

Recall that DGLAP evolution is performed starting from some relatively large virtuality $Q_0^2 > \Lambda_{\text{QCD}}^2$. Everything below $Q_0$ is collected in the phenomenological ‘input’ parton distributions, PDF($Q_0^2$). It would be best to calculate the NLO box diagram also starting the integral over $k^2$ from $Q_0^2$; that is to consider $\int_{Q_0^2} dk^2/k^2$. On the other hand, in such a case we will introduce a new dimensionful parameter which will strongly complicate the logarithmic DGLAP evolution. If we assume that the starting scale is much smaller than the final factorization scale $\mu^2_F = Q^2$, that is $Q_0^2 \ll Q^2$, then we may consider the non-singular contribution from the region with $k^2 < Q_0^2$ as a power, $O(Q_0^2/Q^2)$, correction, and neglect it, together with the other power corrections. That is to consider the $\int_0 dk^2/k^2$.

The problem is that within the conventional approach we deal with singular integrands. Using the dimensional, $\epsilon$, regularization, this results in a $1/\epsilon$ pole plus some constant terms (with respect to $\epsilon$) of infrared, $\epsilon/\epsilon$, origin. Finally we subtract the ‘anomalous dimension’ function\textsuperscript{9} generated by the LO DGLAP evolution; that is we eliminate the $1/\epsilon$ pole. However, this prescription does not cancel the constant terms (with respect to $\epsilon$) of infrared, $\epsilon/\epsilon$, origin.

If, on the other hand, we adopt the ‘physical’ approach and integrate the non-singular expression given by the box diagram minus the DGLAP-generated $P^{\text{LO}} \otimes P^{\text{LO}}$ contribution (or put some infrared cutoff $q_0$ on the integral $\int_{q_0^2} dk^2/k^2$) we will never obtain such terms of infrared origin, even for $q_0 \to 0$. We thus conclude that the above LO DGLAP-generated subtraction will give the correct NLO splitting functions, and that the term coming from very large distances (greater than $O(1/\Lambda_{\text{QCD}})$) in the ‘conventional’ procedure should not be there.

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