

QCD splitting-function dependence on evolution variable

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Abstract: In this note we present the results of the study performed at IFJ PAN in Krakow in the context of the parton-shower-related research which suggests that QCD splitting function depends on the type of physical evolution variable. We sketch a method of including NLO corrections into a cascade (ladder) of a parton shower (PS). The method is based on the Curci-Furmanski-Petronzio classical paper and matching between PS and NLO distributions obtained from the Feynman diagrams. We note that one does not reproduce the classical CFP NLO kernels if the k_T ordering is used for the calculation. This seems to be a limitation of the definition of the CFP procedure and requires further investigation.

Parton-Shower-related activities at IFJ PAN

Let us give a short review of the results of the group at IFJ PAN related to the parton shower (PS) research.

- (A) In the first step we developed a series of works on evolving parton distribution functions (PDFs) using Markovian Monte Carlo (MMC) programs. We solved exactly the LO and NLO DGLAP evolution equations using the MMC methods with long term aim of developing new types of PS MCs. a) The first MMCs were evolving quark and gluon PDFs according to the LO and NLO DGLAP equations in $t = \ln \mu^2$ variable, with the running $\alpha_s(t)$ [1], [2]. b) Next we developed the MMC for CCFM (1-loop) gluonstrahlung with $t = \ln(k_T/(1-z))$ and $\alpha_s(k_T^2)$ [3]. c) Finally, the most sophisticated MMCs, with various choices of t and an argument of α_s , combining the full DGLAP and CCFM, together with the numerical results, were reported in [4]. d) A corresponding MMC program in C++ implementing variety of the above models was published in [5]. The PDFs $q(t, x)$ and $g(t, x)$ from MMCs agreed (within 3-digits) with non-MC calculations of the QCDnum16 and APChab programs. e) Let us mention also a more recent study of the MCFM modelling exploiting our MMC done in Ref. [6].
- (B) MMCs were designed to test numerically series of constrained MC (CMC) programs, implemented with the same evolution, but with a constrained predefined final x variable (an alternative to the backward evolution in the PS MC) aiming at strict (NLO level) control of the distributions of LO PS MC. a) We began with CMC and MMC modules for single ladder/shower, without a hard process, with exclusive LO kernels, optionally inclusive NLO kernels. CMC programs tested with MMCs were published in [7], [8]. b) Next, we combined two CMCs with a hard process matrix element (ME) into PSMC for Drell-Yan (DY) process

[9], [10]. Such an academic PS MC was instrumental in testing new ideas on the NLO corrections in the exclusive evolution included in all kernels in the initial state ladders/showers, and on the NLO corrections to the hard process (a simpler alternative to MC@NLO [11] and POWHEG [12] approaches) thanks to a perfect numerical and algebraical control over LO distributions.

- (C) The next project concerns the NLO corrections in an exclusive form to PS MC (one ladder):
- a) The first solution, albeit limited to the non-singlet evolution, was proposed and tested numerically in [13], [14].
 - b) The exclusive NLO kernels had to be recalculated in the Curci-Furmanski-Petronzio (CFP) [15] framework from scratch. The non-singlet 2-real kernels were recalculated in [16] and the non-singlet 1-real-1-virtual ones in the PhD thesis [17]. Also the singlet evolution kernels are now almost complete (unpublished).
 - c) Simplified and faster scheme was reported (numerical tests) in [18].
 - d) An even simpler and faster scheme of NLO-correcting was presented in [19].
 - e) A major problem was how to include consistently virtual corrections. The first solution has been formulated but not yet published. It exploits recalculated virtual corrections in the CFP scheme to the non-singlet kernels [20], [17].
- D) Finally, the last project is the implementation of the NLO corrections to the hard process in a new way, named KrkNLO method. It provides a simpler alternative to MC@NLO and POWHEG methods:
- a) Methodology for DY process was defined (without numerical tests) in [21].
 - b) Numerical tests of KrkNLO with the help of Double-CMC PS was done in [10].
 - c) Further improvements and introduction of PDFs in the MC factorisation scheme can be found in [22], where the the MC implementation is still done with the help of the not-so realistic Double-CMC PS.
 - d) Next, the KrkNLO method was implemented on top of the SHERPA [23] parton shower. Comparisons for the DY process with the fixed-order NLO calculations of MCFM [24] as well as with the NLO-PS matched calculations of MC@NLO and POWHEG were also done in [25] and in the conference note [26].
 - e) A complete definition of a new scheme for PDFs, called the MC scheme, necessary for the KrkNLO method, both for the DY processes and Higgs-boson production via gluon-gluon fusion was presented in [27] and in the conference note [28].
 - f) At last, the implementation of the KrkNLO method in the Herwig 7 MC generator [29] and a series of comparisons for Higgs production in gluon-gluon fusion with the programs: MCFM, MC@NLO, POWHEG and HNNLO [30] as well as with experimental data of the ATLAS Collaboration from the LHC Run 1 was published in [31].

In this note we briefly present the findings of Ref. [32]. This is however, a rather technical paper, so let us first explain the background and then summarise on the main results of Ref. [32]. Generally, the context is that of the N+NLO parton shower, with the NLO evolution kernels (PDFs) in the shower (ladder) and the NNLO hard process. As listed earlier, we have already defined and tested some important elements of the complete N+NLO parton shower.

What is NLO Parton Shower?

Let us start with a simpler question: What is LO PS? It is built using the LO-class evolution kernels and the LO PDFs within a certain collinear factorisation scheme. LO PS MC implements the LO DGLAP evolution for semi-inclusive distributions (structure functions, etc.). If the hard process is corrected to the NLO level (N+LO), LO PS MC desirably encapsulates all collinear/soft singularities of the NLO hard process ME in the exclusive form. In the N+LO schemes *one* parton originally generated within LO PS MC gets promoted to the hard process.

contribution of this graph to the NLO kernel is different than in the original CFP paper. This is, of course, rather a problem of CFP, not of PS MC, as according to the theorem proven in the CFP paper the dependence of the MS NLO kernel on the choice of the upper kinematic boundary Q is forbidden. The problem is absent for the gluonstrahlung (non-singlet) kernels shown in the right-hand-side plots of Fig. 1 as well as for all single-pole diagrams. Next, we found that the choice of the type of the “evolution time”, i.e. the Q -variable, is essential: for $Q = \text{virtuality}$, $Q = |\vec{k}_{T1} + \vec{k}_{T2}|$ and $Q = |\vec{k}_{T1}| + |\vec{k}_{T2}|$ the problem is absent. It shows up only for the diagrams with the final-state radiation (FSR) gluon splitting into a gluon pair calculated for $Q = \max(k_{T1}, k_{T2})$, combined with the relevant virtual diagram counter-partner. The analytical calculation shows that source of the effect is a mismatch between the upper kinematic limit in the virtual and real diagram contributions. Such a mismatch is usually absent, except for the $Q = \max(k_{T1}, k_{T2})$ case.

Let us discuss a list of possible solutions of this problem: A) The simplest explanation would be that the CFP scheme is not as general as expected, though it is rather unlikely. B) Is it then due to partial introduction of the dimensional regularisation in CFP, only for the collinear singularity, not for the light-cone variable? C) Perhaps it is due to the use of the axial gauge? D) Finally, maybe it cancels when combined with the NNLO hard process?

Summary and outlook

We have reviewed a methodology of including NLO corrections along the parton shower MC ladder. It has been tested numerically. The method is based on matching between PS MC and CFP diagrammatic NLO distributions. We have spotted a problem of not reproducing NLO evolution kernels in the calculations with k_T -ordering kinematics. It is essentially the problem of the CFP scheme itself and requires better understanding.

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