

Numerical evaluation of loop corrections

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Together with Sebastian Becker and Stefan Weinzierl
[New in the group: Daniel Götz and Christopher Schwan]

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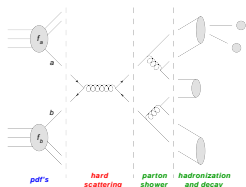
QCD @ LHC 2011, August 25th 2011
St Andrews, Scotland



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 - Experimenter's wish list: [\[Les Houches wish list\]](#)
 $pp \rightarrow VV + jets, H + 2jets, t\bar{t}b\bar{b}, t\bar{t} + 2jets, VVb\bar{b}, VV + 2jets, V + 3jets, VVV.$
 - Due to the large QCD background we need to find the famous "needle in the haystack".
- ⇒ A detailed understanding of multi-parton QCD final states is unavoidable!



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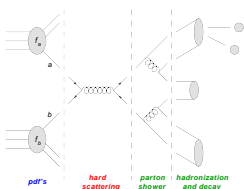


$$d\sigma_{pp \rightarrow x_{had.}}(s, \mu_F^2, \mu_R^2) \propto f_a(x_a; \mu_F^2) f_b(x_b; \mu_F^2) \otimes d\hat{\sigma}_{ab \rightarrow x_{partonic}}(\hat{s}; \{p_x\}, \mu_F^2, \mu_R^2) \otimes [{}^{\prime}PS/Had.{}^{\prime}]$$

- During the workshop we saw that for a reliable prediction a good description of all the parts is necessary, where we will focus on the hard scattering.
- $d\hat{\sigma}_{ab \rightarrow x_{partonic}}(\hat{s}; \{p_x\}, \mu_F^2, \mu_R^2) \propto |\mathcal{A}_{ab \rightarrow x_{partonic}}(\hat{s}; \{p_x\}, \mu_F^2, \mu_R^2)|^2$
- Many jets at the LHC: Want $\mathcal{A}_{ab \rightarrow x_{partonic}} \equiv \mathcal{A}_n$ for large n , w/ $n = 2 + \#x!$
- Want $|\mathcal{A}_n|^2$ at NLO in α_s due to a large scale dependence at LO and more accurate jet descriptions! See talk by Joey Huston on Tuesday.

Introduction - Physics at the LHC

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The NLO pieces to a LO contribution $\mathcal{A}_n^{(0)}$

$$|\mathcal{A}_n|^2 = |\mathcal{A}_n^{(0)}|^2 + 2\text{Re}(\mathcal{A}_n^{(0)*} \mathcal{A}_n^{(1)}) \sim \left| \text{Diagram 1} \right|^2 + 2 \text{Re} \left(\text{Diagram 2} \times \text{Diagram 3} \right) \quad \left| \mathcal{A}_{n+1} \right|^2 = |\mathcal{A}_{n+1}^{(0)}|^2 \sim \left| \text{Diagram 4} \right|^2$$

Our goal

Evaluate $|\mathcal{A}_n|^2$ for large n to NLO accuracy in α_s [especially the virtual piece] in a fully numerically Monte Carlo (MC) framework! [With the intention for real application, at the moment: $e^+e^- \rightarrow jets$ and $pp \rightarrow V + jets$]

Problem

$2\text{Re}(\mathcal{A}_n^{(0)*} \mathcal{A}_n^{(1)})$ and $|\mathcal{A}_{n+1}^{(0)}|^2$ contain singularities:

- In $\mathcal{A}_{n+1}^{(0)}$ due to the **unresolved 1-particle phase space integration** (soft and collinear).
- In $\mathcal{A}_n^{(1)}$ due to the **loop integration** (soft, collinear and ultraviolet).

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Tackle the real emission: $\mathcal{A}_{n+1}^{(0)}$

- Subtract suitably chosen dipole terms $\mathcal{D}_{[j,k]}$ in order to get a finite integrand [Catani, Seymour].
- This procedure is well known and exists in various improvements [Dittmaier et al., Czakon et al., Gehrmann et al., ...] and variations like residue subtraction [Frixione et al., ...], antenna subtraction [Kosower, Glover et al., ...], ...

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Tackle the virtual contributions: $\mathcal{A}_n^{(1)}$

- Feynman graph approach: [Passarino & Veltman, Denner & Dittmaier, ...]
 - Each single Feynman diagram has to be considered! The complexity grows factorially with the number of legs!
- Unitarity based methods: [BDK, OPP, Anastasiou et al., Berger et al., Ellis, Giele, Kunszt, Melnikov, Zanderighi, ...]
 - Can write any one-loop amplitude as linear combination of a [small] set of master integrals: $\mathcal{A}_n^{(1)} = \sum_j c_j I_j + \mathcal{R}$.
 - The c_j are rather involved functions of external momenta and helicities. Rational terms \mathcal{R} have to be considered.
 - In practice the c_j are determined numerically for each phase space point. Several evaluations are needed for a certain precision, which takes time!

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However, I was missing the following: An alternative, fully numerical method to evaluate the virtual piece. Based on ideas which were first investigated by [Nagy & Soper et al.] we would like to explore this alternative.

Fully numerical MC solution

- **The error of a MC does not depend on the dimensionality of the integration region** and grows just about as $1/\sqrt{N}$, with N the number of integrand evaluations.
- **Advantage:** The $(3n - 4)$ -dimensional **phase-space integral** and the 4-dimensional **loop integral can be performed together in one single MC evaluation** at $(\{p_1, p_2, \dots, p_n\}, k)$, where the p_j are the external momenta and k the loop momentum. No need to evaluate the inner loop integral separately per phase-space point! No extra cost!
- **However: A fully numerical [MC] integration has to be performed in $D = 4$.** Instabilities in the integrand due to infrared (IR) and ultraviolet (UV) divergences have to be taken care of first.

Introduction - Fully numerical method to solve $\mathcal{A}_n^{(1)}$

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Numerical roadmap to $\mathcal{A}_n^{(1)}$

[Becker, Reuschle, Weinzierl - JHEP 1012:013,2010]

- 1) For a numerically stable loop integration of $\mathcal{A}_n^{(1)}$ in $D=4$ we need to subtract the IR and UV divergences first.
 - ⇒ **Extend the subtraction method to the virtual part of the NLO calculation!** Slides 5, 6; Backup Slides
 - ⇒ **Devise [local] virtual subtraction terms!** Slides (8), 9 - 12
- 2) The construction of these subtraction terms depends on a fixed cyclic ordering of the external legs.
 - ⇒ **Work with color decomposition and color ordered Feynman rules!** Slides (6), 7
 - ⇒ **Use partial amplitudes rather than a pure Feynman diagrammatic approach!** Slides (6), 7
 - This reduces the complexity down to about exponential growth with the number of external legs.
- 3) The loop integrand [as well as the total UV subtraction term] may be constructed recursively.
 - ⇒ **Use Berends-Giele type recursion relations on color ordered one-loop off-shell currents!** Slides 13, (14), 15)
- 4) Some of the loop propagators still go on-shell for certain values of the loop momentum.
 - ⇒ **Find a suitable and numerically stable deformation of the integration contour into the complex plane!**
 - Won't go into detail here. Backup Slides

Final destination: A scheme in order to combine subtraction terms and contour deformation into one compatible method!

$$\langle O \rangle^{\text{LO}} + \langle O \rangle^{\text{NLO}} = \int_n O_n d\sigma^{\text{B}} + \int_{n+1} O_{n+1} d\sigma^{\text{R}} + \int_{n[+\text{loop}]} O_n d\sigma^{\text{V}} + \int_n O_n d\sigma^{\text{C}}$$

- $d\sigma^{\text{B}}$: Born level; $d\sigma^{\text{R}}$: Real emission; $d\sigma^{\text{V}}$: Virtual contribution; $d\sigma^{\text{C}}$: Initial state collinear subtraction term.
- Each of the **NLO** terms is separately divergent and only their sum is finite.
- However, for a numerical integration each term needs to be finite.
- Introduce additional terms to subtract the divergencies.
 - **Real emission**: Subtraction of suitable [dipole] terms $d\sigma^{\text{A}}$. Already known.
 - **Virtual contribution**: Subtract suitably chosen virtual subtraction terms $d\sigma^{\text{A}'}$ at the loop integrand level.
 - **Enables a fully numerical loop integration!**

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Wish list for $d\sigma^{\text{A}'}$

$$\langle O \rangle^{\text{NLO}} = \int_{n+1} (O_{n+1} d\sigma^{\text{R}} - O_n d\sigma^{\text{A}}) + \int_{n+\text{loop}} (O_n d\sigma^{\text{V}} - O_n d\sigma^{\text{A}'}) + \int_n (O_n \int_{\text{loop}} d\sigma^{\text{A}'} + O_n d\sigma^{\text{C}} + O_n \int_1 d\sigma^{\text{A}})$$

- Introduce additional terms $d\sigma^{\text{A}'}$ to render $(d\sigma^{\text{V}} - d\sigma^{\text{A}'})$ finite at the loop level.
- Integration of $d\sigma^{\text{A}'}$ yields simple analytic results, which cancel the poles of $d\sigma^{\text{C}} + \int_1 d\sigma^{\text{A}}$.
- The terms in brackets are finite:
 - The **subtracted real term** and the **subtracted virtual term(!)** are finite and can be integrated numerically!
 - The **finite remaining term** exhibits a simple analytical structure!

On the amplitude level $d\sigma^V \propto 2\text{Re}(\mathcal{A}_n^{(0)*} \mathcal{A}_n^{(1)}) d\phi_n$

- $\mathcal{A}_n^{(1)} = \mathcal{A}_{\text{bare}}^{(1)} + \mathcal{A}_{\text{CT}}^{(1)} = (\mathcal{A}_{\text{bare}}^{(1)} - \mathcal{A}_{\text{UV}}^{(1)} - \mathcal{A}_{\text{IR,soft}}^{(1)} - \mathcal{A}_{\text{IR,coll}}^{(1)}) + (\mathcal{A}_{\text{CT}}^{(1)} + \mathcal{A}_{\text{UV}}^{(1)} + \mathcal{A}_{\text{IR,soft}}^{(1)} + \mathcal{A}_{\text{IR,coll}}^{(1)})$.
- $\mathcal{A}_n^{(1)}$ is the finite renormalized amplitude. All IR and UV singularities are contained in the bare amplitude $\mathcal{A}_{\text{bare}}^{(1)}$.
- Define integrands $\mathcal{G}_x^{(1)}$ inside the amplitudes via $\mathcal{A}_x^{(1)} \equiv \int \frac{d^D k}{(2\pi)^D} \mathcal{G}_x^{(1)}$, $x = \text{bare, uv, soft, coll}$.
- They match exactly the singular behavior of $\mathcal{G}_{\text{bare}}^{(1)}$ in the divergent points of the integration region and are easily integrable analytically.
- $\mathcal{A}_n^{(1)} = (\int \{ \mathcal{G}_{\text{bare}}^{(1)} - \mathcal{G}_{\text{UV}}^{(1)} - \mathcal{G}_{\text{IR,soft}}^{(1)} - \mathcal{G}_{\text{IR,coll}}^{(1)} \}) + (\mathcal{A}_{\text{CT}}^{(1)} + \mathcal{A}_{\text{UV}}^{(1)} + \mathcal{A}_{\text{IR,soft}}^{(1)} + \mathcal{A}_{\text{IR,coll}}^{(1)})$.

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In the second bracket the UV subtraction term cancels [analytically] against the UV counterterm from renormalization, whereas the soft and collinear subtraction terms cancel [analytically] against the dipole contributions from real radiation.

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Cyclic ordering

Our virtual subtraction terms [as well as the contour deformation] depend on a fixed cyclic ordering of the external legs in the amplitude!

- Work with color ordered partial amplitudes. [Next slide]
- Define [partial] integrands $\mathcal{G}_x^{(1)}$ inside the partial amplitudes via $A_x^{(1)} \equiv \int \frac{d^D k}{(2\pi)^D} \mathcal{G}_x^{(1)}$, $x = \text{bare, uv, soft, coll}$, henceforth be denoted as "subtraction terms". [Note non-caligraphic letters]
- $\mathcal{G}_{\text{bare}}^{(1)}$ and $\mathcal{G}_{\text{UV}}^{(1)}$ will be constructed recursively.
- $\mathcal{G}_{\text{soft}}^{(1)}$ and $\mathcal{G}_{\text{coll}}^{(1)}$ will be formulated directly on the amplitude level.

Intermezzo: Color decomposition

Factorizing **color information** and **kinematic information** yields [simpler] color stripped amplitudes with a fixed cyclic ordering of the external legs. Gaining color information is simply a combinatorial issue.

Example: N-gluon amplitude in the color-flow decomposition

$$\dots g^2 = 4\pi\alpha_s$$

$$\mathcal{A}_n^{(1)}(1, \dots, n) = \left(\frac{g}{\sqrt{2}}\right)^{n-2} \left[\sum_{\sigma \in S_n/Z_n} N_c \delta_{i\sigma_1 j\sigma_2} \dots \delta_{i\sigma_n j\sigma_1} A_{n,0}^{(1)}(g_{\sigma_1}, \dots, g_{\sigma_n}) \right. \\ \left. + \sum_{\substack{\sigma \in S_n / (Z_m \times Z_{n-m}) \\ \text{For all partitions } m > 0}} \delta_{i\sigma_1 j\sigma_2} \dots \delta_{i\sigma_m j\sigma_1} \delta_{i\sigma_{m+1} j\sigma_{m+2}} \dots \delta_{i\sigma_n j\sigma_{m+1}} A_{n,m}^{(1)}(g_{\sigma_1}, \dots, g_{\sigma_m}; g_{\sigma_{m+1}}, \dots, g_{\sigma_n}) \right]$$

- Description of $\mathcal{A}_n^{(1)}$ in terms of color ordered one-loop partial amplitudes $A_{n,m}^{(1)}$. No color information to be considered.
- Subleading [in color] one-loop partial amplitudes $A_{n,m \neq 0}^{(1)}$ can be related to leading one-loop partial amplitudes $A_{n,0}^{(1)}$.
- Similar decompositions available for processes with m quark-pairs and n gluons. One has to classify all possible color structures in the permutation sum.

Properties

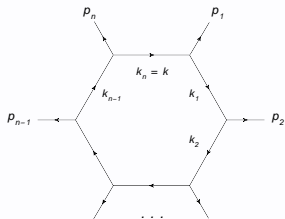
- **Color ordering:** Representation of all graphs with the same number and the same fixed cyclic ordering of the external legs by a partial amplitude. This ensures a fixed [fully ordered] structure of the loop propagators.
- **Particle content:** One-loop partial amplitudes may be further decomposed into **primitive amplitudes**, classified by the quark or gluon content in the loop: $A_{n,0}^{(1)} = A_{n,0,lc}^{(1)} + \frac{n_f}{N_c} A_{n,0,nf}^{(1)}$.
→ In amplitudes with **mixed quark/gluon content** in the loop the **routing of the fermion lines through the loop** matters.
- Partial/primitive amplitudes are **gauge invariant**, which is important for our method, and obey momentum conservation.
- [Mangano & Parke, Maltoni et al., Weinzierl, Bern, Dixon, Kosower, ...]

Intermezzo: One-loop integration

In order to construct [local] virtual subtraction terms we have to **understand the singular structure** of our one-loop integrand.

Using **partial amplitudes** ensures a **fixed sequential propagator structure**, with **only n different loop propagators** to consider. This enables a meaningful classification of the divergent regions in regard to automatization.

Typical one-loop diagram



$$\int \frac{d^4 k}{(2\pi)^4} P_a(k) \prod_{i=1}^n \frac{1}{k_i^2 - m_i^2 + i\delta}, \quad k_i \equiv k - q_i.$$

$P_a(k)$ a polynomial of degree a in k and $q_i \equiv p_1 + \dots + p_i$.

Divergent regions

... with m_i the masses in the loop and p_i the external momenta

Soft infrared divergencies for $k \sim q_i$, if $p_i^2 = m_{i-1}^2$, $m_i = 0$, $p_{i+1}^2 = m_{i+1}^2$.
Massless particle exchanged between two on-shell particles & $k_i \rightarrow 0$

$$k \rightarrow q_i \Rightarrow \begin{cases} k_{i-1}^2 - m_{i-1}^2 & \rightarrow & p_i^2 - m_{i-1}^2 & = & 0 \\ k_i^2 - m_i^2 & \rightarrow & 0 - m_i^2 & = & 0 \\ k_{i+1}^2 - m_{i+1}^2 & \rightarrow & p_{i+1}^2 - m_{i+1}^2 & = & 0 \end{cases}$$

Collinear infrared divergencies for $k \sim q_i - xp_i$, if $p_i^2 = 0$, $m_{i-1} = 0$, $m_i = 0$.
 $x \in [0, 1]$ Massless external on-shell particle attached to two massless propagators & $k_i || p_i$

$$k \rightarrow q_i - xp_i \Rightarrow \begin{cases} k_{i-1}^2 - m_{i-1}^2 & \rightarrow & (1-x)^2 p_i^2 - m_{i-1}^2 & = & 0 \\ k_i^2 - m_i^2 & \rightarrow & x^2 p_i^2 - m_i^2 & = & 0 \end{cases}$$

Ultraviolet divergencies for $k \rightarrow \infty$, if $4 + a - 2n \geq 0$.

Regularization

- In dimensional regularization we use a D -dimensional integral, with $D = 4 - 2\epsilon$ and $|\epsilon| \ll 1$. The result is analytically known and can be expanded around ϵ , which yields terms $\propto 1/\epsilon$ and $\propto 1/\epsilon^2$.
- For integrals with large n the traditional analytic calculation is cumbersome. We choose a numerical method which, however, has to be applied in $D = 4$.

Subtraction terms - IR subtraction I

As we saw the infrared divergences are related to **soft and collinear partons in the loop**. An amplitude with soft or collinear divergences must have **at least one gluon line in the loop**.

The soft and collinear contributions to a given one-loop partial amplitude are the **soft and collinear contributions to all associated one-loop graphs G** with the same number and fixed cyclic ordering of external legs.

Consider $m_i = 0$ and let k_i^2 be a **propagator related to a gluon in the loop** of graph G . The single one-loop graphs G are not to be confused with the [total] subtraction terms $G_{UV,coll,soft}$.

Soft subtraction functions \sim soft source terms

There is a soft singularity when the loop momentum k approximates q_i , or in other words when $k_i \rightarrow 0$. We define a soft subtraction function [Nagy & Soper, ...]:

$$S_i^{soft}(G, p_1, \dots, p_n) = \frac{\lim_{k \rightarrow q_i} \{ k_{i-1}^2 k_i^2 k_{i+1}^2 G(k, p_1, \dots, p_n) \}}{k_{i-1}^2 k_i^2 k_{i+1}^2}$$

How to: 1) "Multiplying out" the dangerous propagators. 2) Taking the soft limit. 3) "Dividing" the proper structure "back in".

Collinear subtraction functions \sim collinear source terms

There is a collinear singularity when the loop momentum k approximates $q_i - xp_i$, w/ $x \in [0, 1]$, or in other words $k_i || p_i$. We define a collinear subtraction function [Nagy & Soper, ...]:

$$S_i^{coll}(G, p_1, \dots, p_n) = \frac{\lim_{k \rightarrow q_i - xp_i} \{ k_{i-1}^2 k_i^2 G(k, p_1, \dots, p_n) \}}{k_{i-1}^2 k_i^2} g_{UV}(k_{i-1}^2, k_i^2)$$

- Introduce a factor $g_{UV}(k_{i-1}^2, k_i^2)$ to avoid possible UV divergences in the collinear subtraction term.
- $g_{UV}(k_{i-1}^2, k_i^2) = 1$ in the collinear region and suppresses our collinear term with additional $\sim O(1/k)$ in the UV limit.
- For $x = 0, 1$ the collinear singularity runs into the soft singularity. To avoid double counting we have to subtract these singularities in a suitable way.

Subtraction terms - IR subtraction II

As we saw on the previous slide, each subtraction term $S_i^{\text{soft}}(G, p_1, \dots, p_n)$ or $S_i^{\text{coll}}(G, p_1, \dots, p_n)$ depends on a single one-loop graph G . [Not to be confused with the subtraction terms $G_{\text{uv, coll, soft}}$]

Now, in simple terms: **Summing up the subtraction terms for all graphs** with a **gluon at position i** in the loop, further **summing over all conceivable positions i** in the loop and **using gauge invariance** to simplify the results in the end yields simple total subtraction terms in local form, proportional to the Born level.

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Total [local] IR subtraction terms

New!

$$G_{\text{soft}}^{(1)} = 4i \sum_{j \in I_g} \frac{p_j p_{j+1}}{k_{j-1}^2 k_j^2 k_{j+1}^2} A_j^{(0)} \quad \text{and} \quad G_{\text{coll}}^{(1)} = -2i \sum_{j \in I_g} \left(\frac{S_j g_{UV}(k_{j-1}^2, k_j^2)}{k_{j-1}^2 k_j^2} + \frac{S_{j+1} g_{UV}(k_j^2, k_{j+1}^2)}{k_j^2 k_{j+1}^2} \right) A_j^{(0)}.$$

I_g is the set of gluons in the loop and the propagator corresponding to $j \in I_g$ in the loop belongs to a gluon.

$S_j = 1$ if the outgoing line j corresponds to a quark, $S_j = 1/2$ if it corresponds to a gluon.

- Formulated directly on the amplitude level! Match the soft and collinear limit of the amplitude on integrand level!
- Simple and fast! Ideal for numerical implementation!
- Yield simple analytical results upon integration! Proportional to the Born level amplitudes!

[Assadsolimani, Becker, Weinzierl - Phys.Rev.D81:094002,2010]

[Assadsolimani, Becker, Reuschle, Weinzierl - Nucl.Phys.Proc.Suppl.205-206:224-229,2010]

[Becker, Reuschle, Weinzierl - JHEP 1012:013,2010]

$$S_\epsilon^{-1} \mu_s^{2\epsilon} \int \frac{d^D k}{(2\pi)^D} G_{\text{soft}}^{(1)} = \frac{-1}{(4\pi)^2} \frac{\exp(\epsilon \gamma_E)}{\Gamma(1-\epsilon)} \sum_{j \in I_g} \frac{2}{\epsilon^2} \left(\frac{-2p_j p_{j+1}}{\mu_s^2} \right)^{-\epsilon} A_j^{(0)} \quad \& \quad S_\epsilon^{-1} \mu_s^{2\epsilon} \int \frac{d^D k}{(2\pi)^D} G_{\text{coll}}^{(1)} = \frac{-1}{(4\pi)^2} \frac{\exp(\epsilon \gamma_E)}{\Gamma(1-\epsilon)} \sum_{j \in I_g} \frac{2}{\epsilon} (S_j + S_{j+1}) \left(\frac{\mu_{UV}}{\mu_s} \right)^{-\epsilon} A_j^{(0)}$$

$S_\epsilon \equiv (4\pi)^\epsilon \exp(-\epsilon \gamma_E)$ the typical volume factor in dimensional regularization.

Remark: These are the subtraction terms for the case $m_i = 0$. The IR subtraction terms for the massive case $m_i \neq 0$ have also been derived and are only slightly more involved.

Subtraction terms - UV subtraction I

Consider again, for $m_i = 0$, our one-loop integrand $G(k, n) \equiv P_a(k) \prod_{j=1}^n \frac{1}{k_j^2}$, with $k_j = k - \sum_i p_i$ and $P_a(k)$ is again a polynomial of degree a in k .

It can be shown that the **UV divergent diagrams** are only those which contribute to **propagator or vertex corrections**.

So, G contains all one-loop diagrams that contribute to a given correction. Still, refer to it representatively as "Graph".

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We want a term that mimics the [exact] local UV behavior

- **Expand around** the inverse loop propagator in the UV limit, more precisely around $(\bar{k}^2 - \mu_{UV}^2)^{-1}$, where $\bar{k} \equiv k - Q$.
- $G(k, n) \approx \frac{P_a(\bar{k})}{(\bar{k}^2 - \mu_{UV}^2)^n} \left(1 + \sum_{m=1}^{\ell} \frac{X_m(\bar{k})}{(\bar{k}^2 - \mu_{UV}^2)^m} \right) \equiv G(\bar{k}, n, \ell)$, with $X_m(\bar{k})$ polynomial of order m in \bar{k} μ_{UV} and Q see below
- The **cut on ℓ** depends on the degree of divergence of the graph G : $\ell = 0$ logarithmic; $\ell = 1$ linear; $\ell = 2$ quadratic. E.g. for the gluon self energy use $\ell = 2$, since quadratically divergent. Ultimately: Count only UV divergent powers of \bar{k} .
- Small example: $\frac{1}{(k-p)^2 - m^2} = \frac{1}{\bar{k}^2 - \mu_{UV}^2} \left\{ 1 + \frac{2\bar{k} \cdot (p-Q)}{\bar{k}^2 - \mu_{UV}^2} - \frac{(p-Q)^2 - m^2 + \mu_{UV}^2}{\bar{k}^2 - \mu_{UV}^2} + \frac{(2\bar{k} \cdot (p-Q))^2}{(\bar{k}^2 - \mu_{UV}^2)^2} \right\} + O(1/|\bar{k}|^5)$

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Upon integration

... omit a factor $\mu_S^{2\epsilon} (2\pi)^{-D}$ in the integration measure

- $\int d^D k G(\bar{k}, n, \ell) = C \left(\frac{1}{\epsilon} - \log\left(\frac{\mu_{UV}^2}{\mu_S^2}\right)\right) A_n^{(0)} + R$, with C a constant factor of proportionality and R a finite [rational] term.
- Re-define $G_{UV}(\bar{k}, n, \ell) = G(\bar{k}, n, \ell) - \frac{-2\mu_{UV}^2}{(\bar{k}^2 - \mu_{UV}^2)^3} R$, to absorb the finite term.
- Then: $\int d^D k G_{UV}(\bar{k}, n, \ell) = C \left(\frac{1}{\epsilon} - \log\left(\frac{\mu_{UV}^2}{\mu_S^2}\right)\right) A_n^{(0)} \propto$ "common pole part" \times "Born amplitude", exactly as we want it!

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Use μ_{UV}^2 and Q to control the quality of the contour deformation in the UV region

- **Re** $(\mu_{UV}^2) = 0$ and **Im** $(\mu_{UV}^2) < 0$ ensures that the integration contour in the UV region never approaches the singular surface defined by $(\bar{k}^2 - \mu_{UV}^2) = 0$.
- The **integrated UV subtraction terms** are **independent of the four-vector Q** . We can choose Q to our will in order to enhance the numerical stability of the loop integrand upon contour deformation.
- The results of our calculation [of the one-loop amplitude] are in the end independent of μ_{UV}^2 and Q !

Subtraction terms - UV subtraction II

Total [local] UV subtraction terms

New!

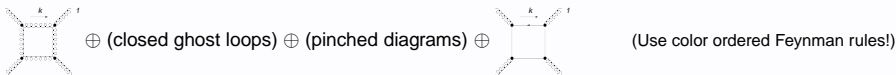
- Only propagator and vertex corrections are UV divergent, quadratically at maximum!
- We can derive a relatively small set of simple local subtraction terms!
- Simple analytic result upon integration, proportional to a common pole part and the respective Born level term!
- These subtraction terms can be used as [local] "counter" terms to recursively construct the total UV subtraction term! The total UV subtraction term will also be proportional to the common pole part and the Born level amplitude!

[Assadsolimani, Becker, Reuschle, Weinzierl - Nucl.Phys.Proc.Suppl.205-206:224-229,2010]

[Becker, Reuschle, Weinzierl - JHEP 1012:013,2010]

Example: 4-gluon vertex

... omit a factor $\mu_s^{2\epsilon} (2\pi)^{-D}$ in the integration measure



$$G_{uv,gggg,lc}^{(1)\mu\nu\lambda\kappa} = \left[\frac{32(1-\epsilon)\bar{k}^\mu\bar{k}^\nu\bar{k}^\lambda\bar{k}^\kappa}{(\bar{k}^2 - \mu_{UV}^2)^4} + \frac{-8(1-\epsilon)W(\bar{k})^{\mu\nu\lambda\kappa} - \frac{4}{3}\mu_{UV}^2 V_4^{\mu\nu\lambda\kappa}}{(\bar{k}^2 - \mu_{UV}^2)^3} + \frac{2(1-\epsilon)(\eta^{\mu\nu}\eta^{\lambda\kappa} + \eta^{\mu\kappa}\eta^{\nu\lambda})}{(\bar{k}^2 - \mu_{UV}^2)^2} \right],$$

$$G_{uv,gggg,nf}^{(1)\mu\nu\lambda\kappa} = \left[\frac{-32\bar{k}^\mu\bar{k}^\nu\bar{k}^\lambda\bar{k}^\kappa}{(\bar{k}^2 - \mu_{UV}^2)^4} + \frac{8W(\bar{k})^{\mu\nu\lambda\kappa}}{(\bar{k}^2 - \mu_{UV}^2)^3} + \frac{4(V_4^{\mu\nu\lambda\kappa} - \eta^{\mu\lambda}\eta^{\nu\kappa})}{(\bar{k}^2 - \mu_{UV}^2)^2} \right], \quad \text{where } W(\bar{k})^{\mu\nu\lambda\kappa} \equiv \eta^{\mu\nu}\bar{k}^\lambda\bar{k}^\kappa + \eta^{\mu\kappa}\bar{k}^\nu\bar{k}^\lambda + \eta^{\nu\lambda}\bar{k}^\mu\bar{k}^\kappa + \eta^{\lambda\kappa}\bar{k}^\mu\bar{k}^\nu$$

$$\int d^D k G_{uv,gggg,lc}^{(1)\mu\nu\lambda\kappa} = \frac{i}{(4\pi)^2} \left(\frac{2}{3} V_4^{\mu\nu\lambda\kappa} \left(\frac{1}{\epsilon} - \log\left(\frac{\mu_{UV}^2}{\mu_s^2}\right) \right) \right) \quad \text{and} \quad \int d^D k G_{uv,gggg,nf}^{(1)\mu\nu\lambda\kappa} = \frac{i}{(4\pi)^2} \left(\frac{4}{3} V_4^{\mu\nu\lambda\kappa} \left(\frac{1}{\epsilon} - \log\left(\frac{\mu_{UV}^2}{\mu_s^2}\right) \right) \right)$$

Check against renormalized 4-gluon vertex $\propto 1 - \frac{\alpha_s}{4\pi} \frac{N_C}{2} \left(\frac{2}{3} + \frac{4}{3} \frac{n_f}{N_C} \right) \frac{1}{\epsilon_{UV}}$

$V_4^{\mu\nu\lambda\kappa}$ = color ordered 4-gluon-vertex

Local counter terms (lc ⊕ nf):  $\hat{=} G_{uv,gggg,lc}^{(1)} \oplus G_{uv,gggg,nf}^{(1)}$

- Utilizing Berends-Giele type **recursion relations**, based on color ordered **[one-loop] off-shell currents** [Berends & Giele, v. Hameren, ...], to construct the [total] bare one-loop integrand $G_{bare}^{(1)}$ and the [total] UV subtraction term $G_{UV}^{(1)}$!

Intermezzo: Recursive methods

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Example: One-loop off-shell gluon current in a 3-valent toy-model

$$\begin{aligned}
 & \text{Diagram 1} = \sum_{i=m}^{n-1} \text{Diagram 2} + \sum_{i=m}^{n-1} \text{Diagram 3} + \text{Diagram 4} \\
 & \text{Diagram 5} = \sum_{i=m}^{n-1} \text{Diagram 6} \\
 & \text{Diagram 7} = \sum_{i=m-1}^{n-1} \text{Diagram 8}
 \end{aligned}$$

- The **recursive construction of $G_{bare}^{(1)}$** ensures the correct incorporation of all necessary one-loop diagrams to a given partial/primitive amplitude.

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Example: One-loop off-shell gluon current in a 3-valent toy-model

$$\begin{aligned}
 & \text{Diagram 1: } n+1 \text{ line entering a circle with } m \text{ lines exiting, } n \text{ lines on the right.} \\
 & = \sum_{i=m}^{n-1} \text{Diagram 2: } n+1 \text{ line entering a circle with } m \text{ lines exiting, } i \text{ lines on the left, } i+1 \text{ lines on the right, } n \text{ lines on the right.} \\
 & + \sum_{i=m}^{n-1} \text{Diagram 3: } n+1 \text{ line entering a circle with } m \text{ lines exiting, } i \text{ lines on the left, } i+1 \text{ lines on the right, } n \text{ lines on the right.} \\
 & + \text{Diagram 4: } n+1 \text{ line entering a circle with } m \text{ lines exiting, } k_{m-1} \text{ arrow pointing to the circle, } n \text{ lines on the right.} \\
 \\
 & \text{Diagram 5: } n+1 \text{ line entering a circle with } m \text{ lines exiting, } n \text{ lines on the right.} \\
 & = \sum_{i=m}^{n-1} \text{Diagram 6: } n+1 \text{ line entering a circle with } m \text{ lines exiting, } i \text{ lines on the left, } i+1 \text{ lines on the right, } n \text{ lines on the right.} \\
 \\
 & \text{Diagram 7: } n+2 \text{ lines entering a circle with } m \text{ lines exiting, } n \text{ lines on the right.} \\
 & = \sum_{i=m-1}^{n-1} \text{Diagram 8: } n+2 \text{ lines entering a circle with } m \text{ lines exiting, } n+1 \text{ line entering a circle with } m \text{ lines exiting, } i \text{ lines on the left, } i+1 \text{ lines on the right, } n \text{ lines on the right.}
 \end{aligned}$$

- The **recursive construction of $G_{bare}^{(1)}$** ensures the correct incorporation of all necessary one-loop diagrams to a given partial/primitive amplitude.

Example: Total UV subtraction term to the one-loop off-shell gluon current in this toy-model

$$\text{Diagram 9: } n+1 \text{ line entering a circle with } m \text{ lines exiting, } n \text{ lines on the right, marked with an } \otimes. \\
 = \sum_{i=m}^{n-1} \left(\text{Diagram 10: } n+1 \text{ line entering a circle with } m \text{ lines exiting, } i \text{ lines on the left, } i+1 \text{ lines on the right, } n \text{ lines on the right, marked with an } \otimes. \right. \\
 + \text{Diagram 11: } n+1 \text{ line entering a circle with } m \text{ lines exiting, } i \text{ lines on the left, } i+1 \text{ lines on the right, } n \text{ lines on the right.} \\
 + \text{Diagram 12: } n+1 \text{ line entering a circle with } m \text{ lines exiting, } i \text{ lines on the left, } i+1 \text{ lines on the right, } n \text{ lines on the right, marked with an } \otimes. \\
 \left. + \text{Diagram 13: } n+1 \text{ line entering a circle with } m \text{ lines exiting, } i \text{ lines on the left, } i+1 \text{ lines on the right, } n \text{ lines on the right, marked with an } \otimes. \right)$$

- The **recursive construction of $G_{UV}^{(1)}$** ensures the correct incorporation of all necessary UV counterterms to a given partial/primitive amplitude.

Implemented the recursive algorithms for the full theory [so far at leading color] in several C++ libraries! In spinor formalism!

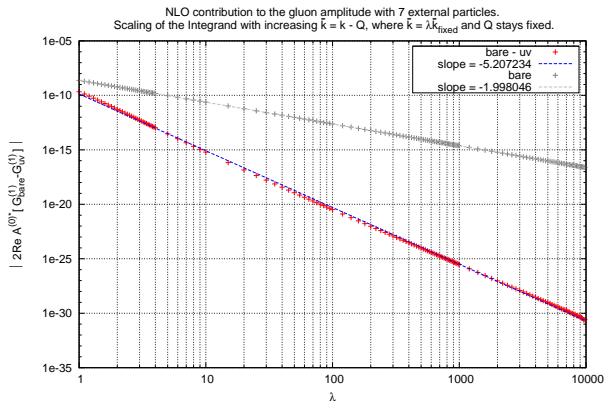
UV subtraction - Consistency check I

Check whether the implemented recursive constructions of the one-loop integrand and the UV subtraction term play along well for large values of the loop momentum.

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Check whether the implemented recursive constructions of the one-loop integrand and the UV subtraction term play along well for large values of the loop momentum.

In the plot we show $|2\text{Re}(A^{(0)*} G^{(1)})|$ vs. a UV scaling parameter λ for an n -gluon amplitude with $n = 7$:
[preliminary, S. Becker, CR]



Summing over all helicities!

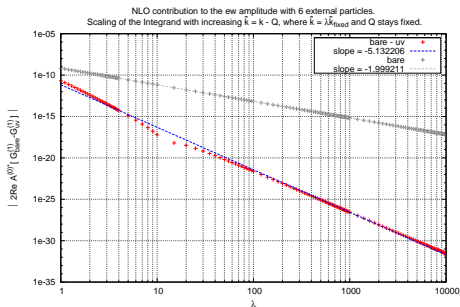
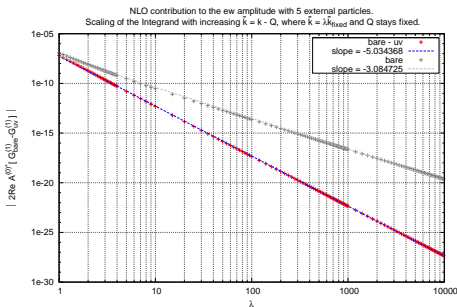
The unsubtracted integrand grows like $\int \frac{d^4 k}{k^2}$: Quadr. UV divergent!

The UV subtracted integrand grows like $\int \frac{d^4 k}{k^5}$: UV finite!

In gray we see the unsubtracted [total] integrand, which is obviously quadratically UV divergent. In red we see the fully UV subtracted [total] integrand, which shows clearly a finite behavior!

UV subtraction - Consistency check II

In these plots we show $|2\text{Re}(A^{(0)*} G^{(1)})|$ vs. a UV scaling parameter λ for $e^+e^- \rightarrow 3/4$ jets [i.e. $q\bar{q} + 1/2 g$'s]:
[preliminary, S. Becker, CR]



In gray we see again the unsubtracted [total] integrand, which is obviously UV divergent. In red we see the fully UV subtracted [total] integrand, which shows again clearly a finite behavior!

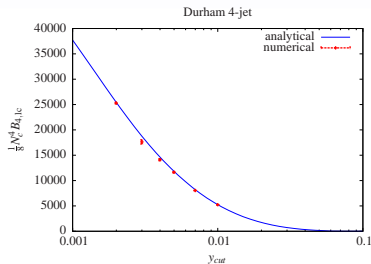
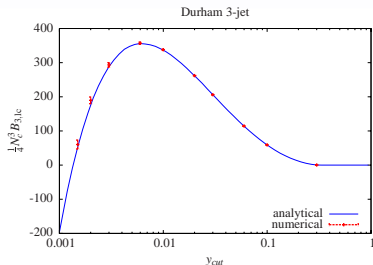
Something to learn: We notice that in the unsubtracted 3 jet event there is only a total linear divergence $\sim d^4k/k^3$ in contrast to the total quadratic divergence $\sim d^4k/k^2$ in the unsubtracted 4 jet event.

Due to the fact that in the 4 jet event the gluon propagator, and hence the gluon self energy, appears off-shell for the first time.

Proof of principle - Computing jet rates in $e^+e^- \rightarrow jets$

Use the whole method in a calculation of jet rates in e^+e^- , together with a simple phase space generating algorithm [RAMBO] and the Durham jet algorithm.

The plots show $e^+e^- \rightarrow 3$ and 4 jets at leading color: $1/8N_c^{3/4} B_{3/4,lc}$ vs. y_{cut} , where $B_{3/4,lc}$ is the NLO coefficient in the perturbative expansion for the 3/4-jet rate and y_{cut} the jet resolution parameter. [S. Weinzierl, preliminary]



Blue shows the results from previous analytic calculations. Red shows the numerical results of our method. Good agreement!

Brief reminder on jet rates

- The production rate for n -jet events, or short the ' n -jet rate', is given by the ratio of the cross section for n -jet events divided by the total hadronic cross section.
- In e^+e^- annihilation:

$$R_n = \frac{\sigma_{n-jet}}{\sigma_{tot}} = \left(\frac{\alpha_s}{2\pi}\right)^{n-2} \bar{A}_n + \left(\frac{\alpha_s}{2\pi}\right)^{n-1} \bar{B}_n + O(\alpha_s^n)$$

- In practice we calculate:

$$\frac{\sigma_{n-jet}}{\sigma_0} = \left(\frac{\alpha_s}{2\pi}\right)^{n-2} A_n + \left(\frac{\alpha_s}{2\pi}\right)^{n-1} B_n + O(\alpha_s^n)$$

with σ_0 the leading order cross section for $e^+e^- \rightarrow hadrons$.

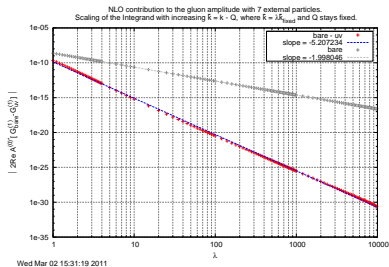
The relation between \bar{A}_n (\bar{B}_n) and A_n (B_n) can be determined from the perturbative expansion of the total hadronic cross section.

There is need for fast high accuracy tools in LHC physics!

- We present a **fully numerical algorithm to compute NLO QCD amplitudes with many legs in the final state** for fixed order of α_s in perturbation theory! [Becker, Reuschle, Weinzierl - JHEP 1012:013,2010]

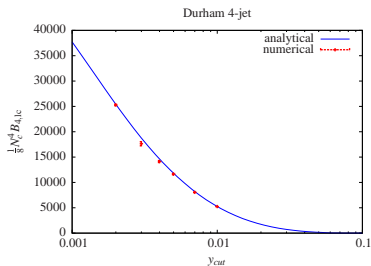
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- We can **reproduce** known analytic results for **$e^+e^- \rightarrow$ up to 4 jets** within high accuracy!



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- We present a **fully numerical algorithm to compute NLO QCD amplitudes with many legs in the final state** for fixed order of α_s in perturbation theory! [Becker, Reuschle, Weinzierl - JHEP 1012:013,2010]
- The **local virtual subtraction method** on the amplitude level is an ideal candidate for a fast **numerical [MC] evaluation of the one-loop integration** and **works very well** in the presented form!
- We can **reproduce** known analytic results for **$e^+e^- \rightarrow$ up to 4 jets** within high accuracy!

In essence

- Our subtraction terms are simple and fast!
- They are tailored such that they go hand in hand with the contour deformation and the real emission!

- Proof of principle: Possible to reproduce the results for $e^+e^- \rightarrow$ up to 4 jets with a purely numerical approach! 5 and 6 jets is in preparation.
- So far we computed our results for leading color! Extension to full color is work in progress.
- The next bigger step would be to apply our method to Z production plus jets at the LHC.

Thank you for your attention!

Backup: Pole structure / Insertion operators L and I

After integration, the soft and collinear poles of a primitive one-loop amplitude with massless partons are given by [Becher & Neubert, Magnea]:

$$S_\varepsilon^{-1} \mu_s^{2\varepsilon} A_{bare}^{(1)} = \frac{\alpha_s}{4\pi} \frac{\exp(\varepsilon\gamma_E)}{\Gamma(1-\varepsilon)} \sum_{i \in I_g} \left[\frac{2}{\varepsilon^2} \left(\frac{-2p_i \cdot p_{i+1}}{\mu_s^2} \right)^{-\varepsilon} + \frac{2}{\varepsilon} (S_i + S_{i+1}) \right] A_i^{(0)},$$

where $S_\varepsilon = (4\pi)^\varepsilon \exp(-\varepsilon\gamma_E)$ the typical volume factor in dimensional regularization. The index i in S_i refers to the external particle: $S_q = S_{\bar{q}} = 1$, $S_g = 1/2$. This is exactly the pole structure also reproduced by our IR subtraction terms. A more familiar [and also involving the UV poles] formulation of this may be:

$$S_\varepsilon^{-1} \mu_s^{2\varepsilon} A_{bare}^{(1)} = \frac{\alpha_s}{4\pi} \frac{\exp(\varepsilon\gamma_E)}{\Gamma(1-\varepsilon)} \left[\frac{(n-2)}{2} \frac{\beta_0}{\varepsilon} + \sum_i \sum_{j \neq i} \mathbf{T}_i \mathbf{T}_j \left(\frac{1}{\varepsilon^2} + \frac{\gamma_i}{\mathbf{T}_i^2} \frac{1}{\varepsilon} \right) \left(\frac{-2p_i \cdot p_j}{\mu_s^2} \right)^{-\varepsilon} \right] A^{(0)},$$

where the first part in the squared brackets is exactly the negative of the counterterm contribution $S_\varepsilon^{-1} \mu_s^{2\varepsilon} A_{CT}^{(1)}$ from UV renormalization.

The sum of the collinear subtraction part for the initial state plus the one-particle phase-space integration over the real subtraction part can be written as:

$$d\sigma^C + \int_1 d\sigma^A = \mathbf{I} \otimes d\sigma^B + \mathbf{K} \otimes d\sigma^B + \mathbf{P} \otimes d\sigma^B,$$

where color correlations still remain. The insertion operators \mathbf{K} and \mathbf{P} pose no problem for the numerical evaluation. The term $\mathbf{I} \otimes d\sigma^B$ has the appropriate pole structure to cancel the IR divergences coming from the loop. Hence, $d\sigma^V + \mathbf{I} \otimes d\sigma^B$ is IR finite.

Remember now that $d\sigma^V \propto 2\text{Re}(\mathcal{A}_n^{(0)*} \mathcal{A}_n^{(1)}) d\phi_n$, where further $\mathcal{A}_n^{(1)} = \mathcal{A}_{bare}^{(1)} + \mathcal{A}_{CT}^{(1)}$. We make $\mathcal{A}_{bare}^{(1)}$ finite by introducing our subtraction terms locally in the "first bracket" and are in the "second bracket" left with an analytical structure of the form

$\mathcal{A}_{CT}^{(1)} + \mathcal{A}_{soft}^{(1)} + \mathcal{A}_{coll}^{(1)} + \mathcal{A}_{UV}^{(1)}$. This structure defines us a new insertion operator \mathbf{L} via:

$$2\text{Re}(\mathcal{A}_n^{(0)*} \mathcal{A}_{CT}^{(1)} + \mathcal{A}_{soft}^{(1)} + \mathcal{A}_{coll}^{(1)} + \mathcal{A}_{UV}^{(1)}) d\phi_n = \mathbf{L} \otimes d\sigma^B.$$

The insertion operator \mathbf{L} contains the explicit poles in the dimensional regularization parameter related to the IR singularities of the one-loop amplitude. These poles cancel when combined with the insertion operator \mathbf{I} :

$$(\mathbf{I} + \mathbf{L}) \otimes d\sigma^B = \text{finite.}$$

Backup: Derivation of the IR subtraction terms - More rigorously I

Soft subtraction term

When gluon i is soft the corresponding propagator goes on-shell and we may replace:

$$\frac{-ig^{\mu\nu}}{k_j^2} \rightarrow \frac{i}{k_j^2} \left(d^{\mu\nu}(k_j^b, n) - 2 \frac{k_j^{b\mu} n^\nu + n^\mu k_j^{b\nu}}{2k_j^b \cdot n} \right),$$

with k_j^b the on-shell limit of k_j , n a light-like reference vector and $d^{\mu\nu}$ the sum over physical polarizations. Adding self-energy diagrams will not change the soft limit. With this inclusion and a similar replacement as above the contribution from the polarization sum makes a partial tree-level amplitude, where two gluons with momenta k_j^b and $-k_j^b$ have been inserted between the legs j and $j+1$:

$$\lim_{k_j \rightarrow 0} \left[\text{Diagram: Hard process} \text{---} \text{Gluon loop } k_j \right] = \lim_{k_j \rightarrow 0} \left[\text{Diagram: Hard process} \text{---} \text{Gluon } (k_j, \lambda) \text{---} \text{Gluon } (-k_j, -\lambda) \right]$$

In the soft limit this tree-level partial amplitude is given by two eikonal factors times the tree-level amplitude without these two additional gluons:

$$\left(\frac{p_j^\mu}{p_j \cdot k_j^b} \right) g^{\mu\nu} \left(\frac{p_{j+1}^\nu}{p_{j+1} \cdot (-k_j^b)} \right) A_j^{(0)}.$$

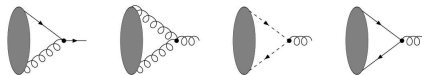
In the soft limit we may replace $2p_j \cdot k_j^b$ by k_{j-1}^2 and $2p_{j+1} \cdot (-k_j^b)$ by k_{j+1}^2 , which then leads to the form of our local soft subtraction terms.

The terms with $k_j^{b\mu} n^\nu + n^\mu k_j^{b\nu}$ in our replacement at the beginning vanish for the sum of all diagrams due to gauge invariance.

Backup: Derivation of the IR subtraction terms - More rigorously II

Collinear subtraction term

We have to consider configurations where two adjacent propagators go on-shell with a massless leg in between:



The diagrams where an external gluon splits into two ghosts or a $q\bar{q}$ -pair are in the collinear limit not singular enough to yield a divergence after integration. We are left with the $q \rightarrow qg$ - and the $g \rightarrow gg$ -splittings.

In $q \rightarrow qg$ one can show that only the longitudinal polarization of the gluon contributes to the collinear limit. The same holds for $g \rightarrow gg$, here the collinear limit receives contributions when one of the two gluons in the loop carries a longitudinal polarization (not both). The external gluon has of course physical transverse polarization.

We use the fact that contraction of a longitudinal polarization into a gauge invariant set of diagrams yields zero. Now the "blobs" of the two cases we just discussed consist almost of a gauge invariant set of diagrams. There is only one missing, where the longitudinal polarized gluon couples directly to the other parton connected to the "blob". This is a self-energy insertion on an external line, by definition absent from the amputated one-loop amplitude.

We turn the argument around and replace the sum of collinear singular diagrams by the negative of the respective self-energy insertions on the external line:

$$\begin{aligned}
 \lim_{k_{j-1} \parallel k_j} \text{Diagram 1} &= - \lim_{k_{j-1} \parallel k_j} \text{Diagram 2} \\
 \lim_{k_{j-1} \parallel k_j} \text{Diagram 3} &= - \lim_{k_{j-1} \parallel k_j} (\text{Diagram 4} + \text{Diagram 5})
 \end{aligned}$$

Collinear subtraction term ctd.

As parametrization for the collinear limit we use the same as is usually used in the real emission case. The singular part of the self-energies is then proportional to:

$$P_{q \rightarrow qg}^{long} = -\frac{2}{2k_{j-1} \cdot k_j} \left(-\frac{2}{1-x} + 2 \right) p',$$

$$P_{g \rightarrow gg}^{long} = -\frac{2}{2k_{j-1} \cdot k_j} \left(-\frac{2}{x} - \frac{2}{1-x} + 2 \right) \left(-g^{\mu\nu} + 2 \frac{p^\mu n^\nu + n^\mu p^\nu}{2p \cdot n} \right).$$

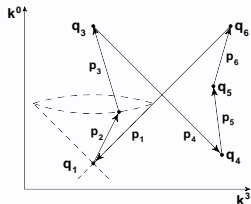
The soft singularities for $x = 0, 1$ must not be double counted. We therefore just have to consider the terms which are non-singular in the soft limit. These terms are independent of x and lead to the form of our local collinear subtraction terms.

Remark: The self-energy insertions introduce a spurious $1/p_j^2$ -singularity. In order to calculate the singular part of the self-energies we regulate this singularity by putting p_j^2 slightly off-shell, but keeping k_{j-1} and k_j on-shell and imposing momentum conservation.

Backup: Integration contour

Consider $m_i = 0$. The denominator in the loop integral becomes singular for $k_i \equiv (k - q_i)^2 = 0$.

- In a diagram where we plot the 0- and 3-components of the loop momentum, $(k - q_i)^2 = 0$ describes a light cone centered on the point q_i .
- Whenever $(k - q_i)^2 = 0$ holds for k , we need to deform the integration contour away from the light cone and into the complex plane.
- We choose for example $-p_1$ and $-p_4$ as our incoming momenta and in such a way that they have components only in the 0-3-plane.
- All other lines p_i are projections onto this plane. They connect the events at q_i , where $p_{i+1} = q_{i+1} - q_i$.



We deform the integration contour into the complex plane, without changing the value of the integral, by deforming the loop momentum into the complex plane:

$$\frac{1}{(k - q_i)^2} \xrightarrow{l(k)=k+i\kappa(k)} \frac{1}{(l(k) - q_i)^2} \xrightarrow{\kappa \text{ "small" }} \frac{1}{(k - q_i)^2 + 2i\kappa \cdot (k - q_i)}$$

We have to choose κ such that whenever $(k - q_i)^2 = 0$ we will get $\kappa \cdot (k - q_i) > 0$, where the numerical stability depends strongly on κ .

The deformation is not possible whenever $k = q_i$ or $k = q_i - xp_i$. These "pinch" singularities are taken care of by the IR subtraction terms.

In practice

First subtract all IR and UV subtraction terms, which yields a UV- and IR-finite integrand. With $P(k)$ and $P_{uv}(k)$ Polynomials in k we have generically:

$$G_{\text{bare}}^{(1)} - G_{\text{soft}}^{(1)} - G_{\text{coll}}^{(1)} - G_{\text{uv}}^{(1)} = \frac{P(k)}{\prod_{j=1}^n (k_j^2 - m_j^2)} - \frac{P_{uv}(k)}{(\bar{k}^2 - \mu_{uv}^2)^{n_{uv}}}$$

We deform the integration contour whenever one of the propagators $1/(k_j^2 - m_j^2)$ or $1/(\bar{k}^2 - \mu_{uv}^2)$ goes on-shell. At the moment we use an algorithm following the idea of Gong, Nagy and Soper [GNS, 2009].