

Resummation benchmark: Scale variations in NangaParbat

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TMD evolution equations

🍏 TMD factorisation allows one to obtain the **evolution equations**:

$$\left\{ \begin{array}{l} \frac{d \ln F}{d \ln \mu} = \gamma(\mu, \zeta) \\ \frac{d \ln F}{d \ln \sqrt{\zeta}} = K(\mu) \end{array} \right. , \quad \frac{d^2 \ln F}{d \ln \mu d \ln \sqrt{\zeta}} = \left\{ \begin{array}{l} \frac{d\gamma}{d \ln \sqrt{\zeta}} \\ \frac{dK}{d \ln \mu} \end{array} \right. = \gamma_K(\alpha_s(\mu))$$

🍏 To solve these equations we need to fix **two pairs of (i.e. four) scales**:

🍏 **initial** scales: (μ_0, ζ_0)

🍏 **final** scales: (μ, ζ)

🍏 The solution is **unique** and reads:

$$F(\mu, \zeta) = R[(\mu, \zeta) \leftarrow (\mu_0, \zeta_0)] F(\mu_0, \zeta_0)$$

$$R[(\mu, \zeta) \leftarrow (\mu_0, \zeta_0)] = \exp \left\{ K(\mu_0) \ln \frac{\sqrt{\zeta}}{\sqrt{\zeta_0}} + \int_{\mu_0}^{\mu} \frac{d\mu'}{\mu'} \left[\gamma_F(\alpha_s(\mu')) - \gamma_K(\alpha_s(\mu')) \ln \frac{\sqrt{\zeta}}{\mu'} \right] \right\}$$

🍏 The question is: **how do we choose these four scales?**

Scale variations

- 🍏 A sensible choice of the scales is important to **allow perturbation theory to be reliable**:
 - 🍏 **no large unresummed logarithms** should be introduced,
 - 🍏 each scale has to be set in the **vicinity of its natural (central) value**,
 - 🍏 scale variations (within a reasonable range) give an estimate of HO corrs.

🍏 In TMD factorisation ($q_T \ll Q$) for DY the relevant scales are q_T and Q :

🍏 natural to expect $\mu_0 \sim \sqrt{\zeta_0} \sim q_T \sim b_T^{-1}$ and $\mu \sim \sqrt{\zeta} \sim Q$

🍏 In fact, it turns out that (in the $\overline{\text{MS}}$ scheme) the **central scales** are:

$$\mu_0 = \sqrt{\zeta_0} = \frac{2e^{-\gamma_E}}{b_T} \equiv \mu_b \quad \text{and} \quad \mu = \sqrt{\zeta} = Q$$

🍏 This choice **nullifies** all unresummed logs. One should thus consider:

$$\mu_0 = C_i^{(1)} \mu_b, \quad \sqrt{\zeta_0} = C_i^{(2)} \mu_b, \quad \mu = C_f^{(1)} Q, \quad \sqrt{\zeta} = C_f^{(2)} Q,$$

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Scale variations

🍏 To reason why variations of ζ have **no effect** is that:

$$\frac{d\sigma}{dq_T} \propto H \left(\frac{\mu}{Q} \right) F_1(\mu, \zeta_1) F_2(\mu, \zeta_2) \quad \text{with} \quad \boxed{\zeta_1 \zeta_2 \stackrel{!}{=} Q^4}$$

🍏 It is easy to see that:

$$F_1(\mu, \zeta_1) F_2(\mu, \zeta_2) = \underbrace{R[(\mu, \zeta_1) \leftarrow (\mu_0, \zeta_0)] R[(\mu, \zeta_2) \leftarrow (\mu_0, \zeta_0)]}_{f(\zeta_1 \zeta_2) = f(Q^4)} F_1(\mu_0, \zeta_0) F_2(\mu_0, \zeta_0)$$

🍏 The single dependence on ζ_1 and ζ_2 **drops** in the combination:

🍏 we choose $\zeta_1 = \zeta_2 = Q^2$ but any other choice such that $\zeta_1 \zeta_2 = Q^4$ is **identical**.

🍏 In addition, in NangaParbat we have chosen to set $\mu_0 = \sqrt{\zeta_0}$:

🍏 not strictly necessary but **probably a conservative choice**.

🍏 At the end of the day, we have **two scales** to be varied:

$$\boxed{\mu_0 = \sqrt{\zeta_0} = C_i \mu_b \quad \text{and} \quad \mu = C_f Q}$$

Comparison to q_T resummation

🍏 In q_T resummation, the **resummation scale** M is introduced as:

$$L = \ln \left(\frac{Q}{\mu_b} \right) = \ln \left(\frac{M}{\mu_b} \right) + \ln \left(\frac{Q}{M} \right)$$

🍏 These logs are **exposed** by expressing integral representations of the argument of the Sudakov in terms of the **functions** g_n :

$$\begin{aligned} \int_{\mu_b}^Q \frac{d\mu'}{\mu'} \left[A(\alpha_s(\mu')) \ln \left(\frac{Q}{\mu'} \right) + B(\alpha_s(\mu')) \right] &= Lg_0(\alpha_s L) + \sum_{n=1}^{\infty} \alpha_s^{n-1} g_n(\alpha_s L) \\ &= Lg_0(\alpha_s L) + \sum_{n=1}^k \alpha_s^{n-1} g_n(\alpha_s L) + \mathcal{O}(\alpha_s^{k+n} L^n) \end{aligned}$$

🍏 The series in the r.h.s. is **truncated** according to the log accuracy:

🍏 the truncation is responsible for the **explicit dependence on** M .

🍏 If the l.h.s. integral is computed exactly, no dependence on M appears:

🍏 this is what we do in NangaParbat by computing the integral numerically,

🍏 therefore, we have **no resummation scale dependence**.

Comparison to q_T resummation

- 🍏 The **renormalisation scale** μ_R in q_T resummation is probably to be (partly) identified with the scale μ in the TMD formalism:
 - 🍏 this is the large scale at which the strong coupling α_s is computed.
- 🍏 The **factorisation scale** μ_F present in q_T resummation is absent in the TMD formalism:
 - 🍏 in the TMD approach, PDFs are computed at the low scale μ_0 :
 - 🍏 μ_0 is varied around μ_b ,
 - 🍏 in q_T resummation, PDFs are evolved from *exactly* μ_b up to μ_F :
 - 🍏 μ_F is varied around Q .
 - 🍏 variations of μ_0 are typically much larger than variations of μ_F because at the energies relevant to the benchmark $\alpha_s(\mu_0) \gg \alpha_s(\mu_F)$:
 - 🍏 problems with NangaParbat in using a b_{\max} too large with scale variations.

Problems with b_{max}

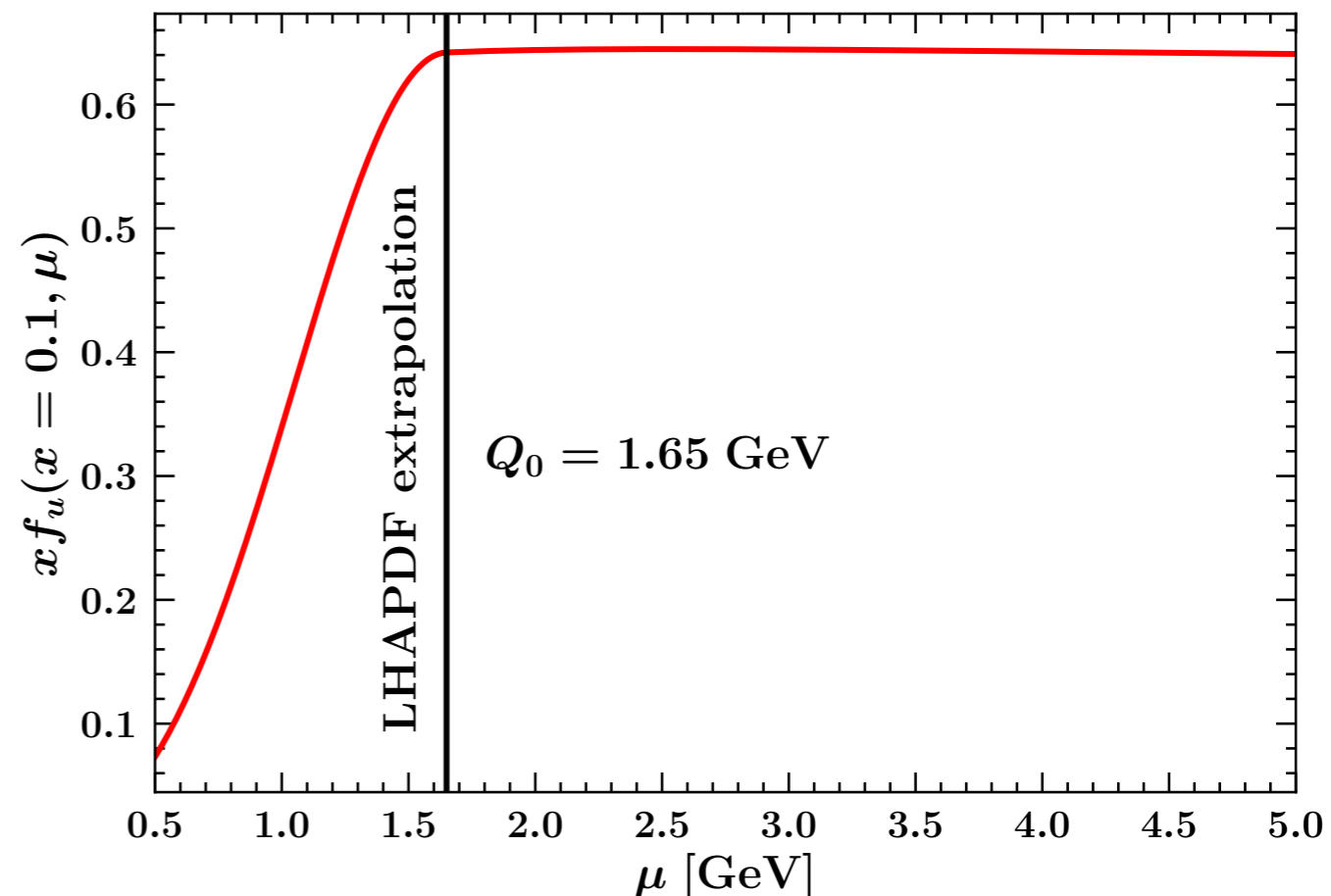
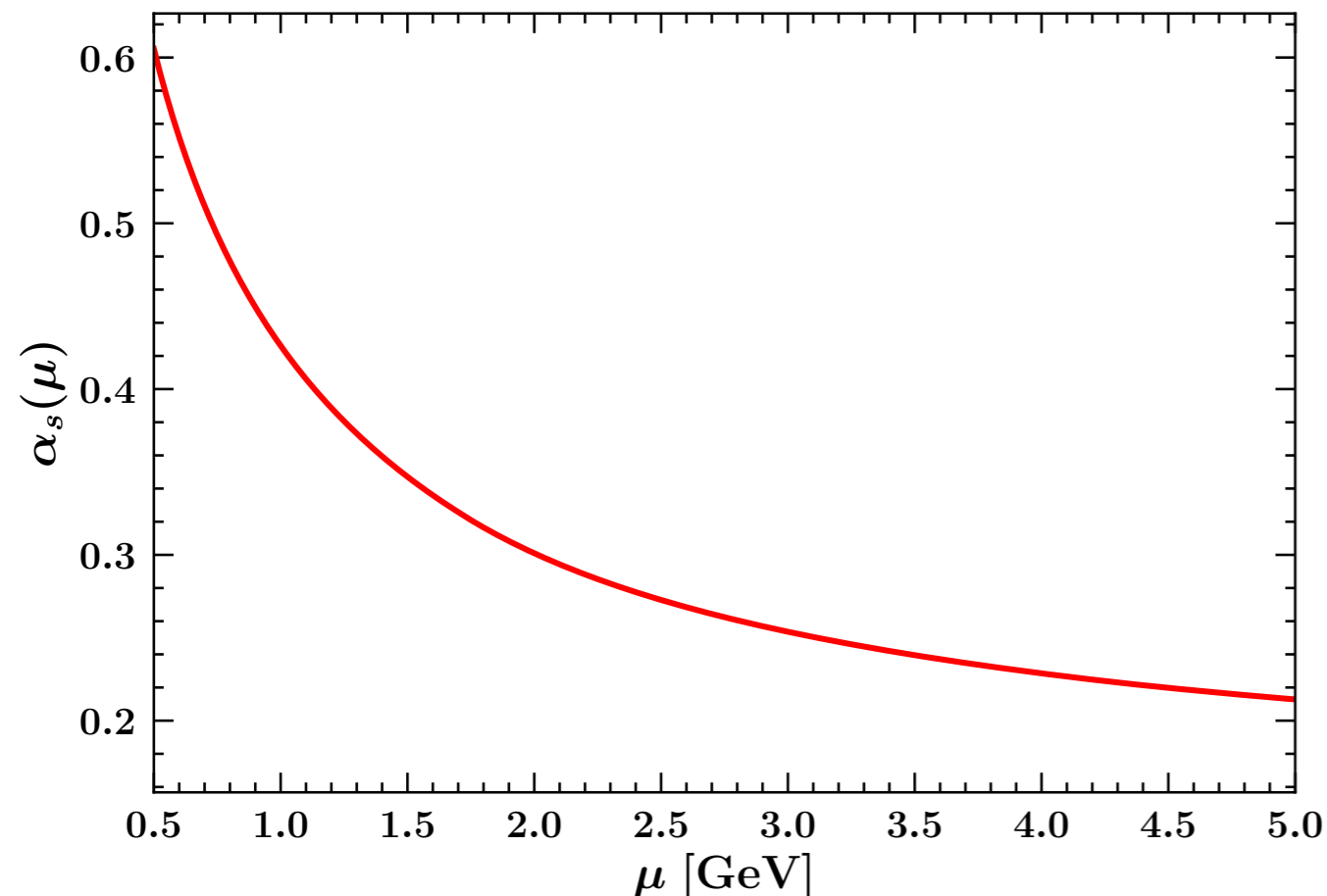
🍏 In the benchmark settings b_{max} is set such that:

$$\mu_{min} = \frac{b_0}{b_{max}} = \frac{2e^{-\gamma_E}}{b_{max}} = 1 \text{ GeV}$$

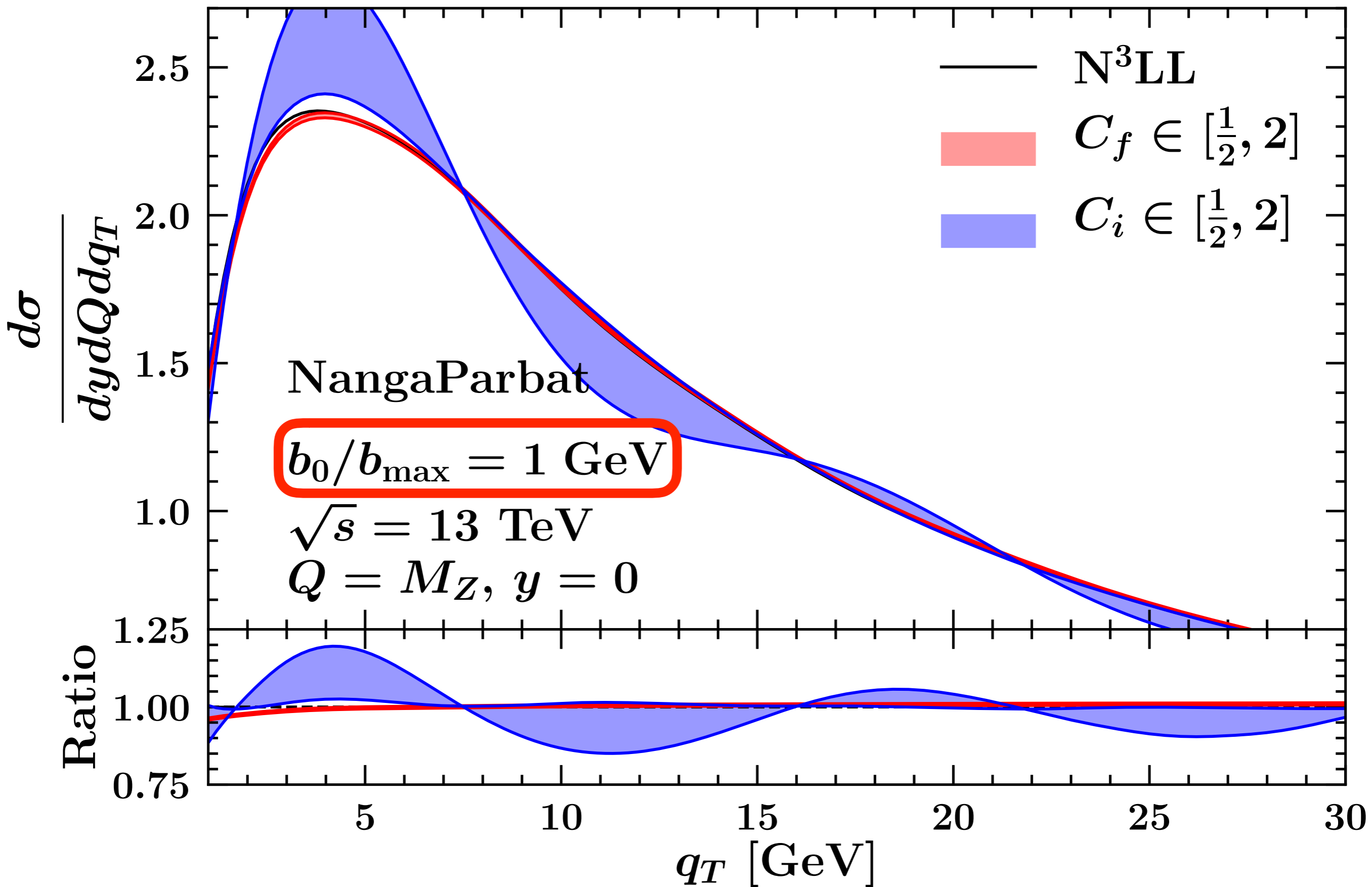
🍏 This is the **minimum** value of the scale at α_s which and PDFs are called.

🍏 Scales variations in NangaParbat cause μ_{min} to be scaled by some factor:

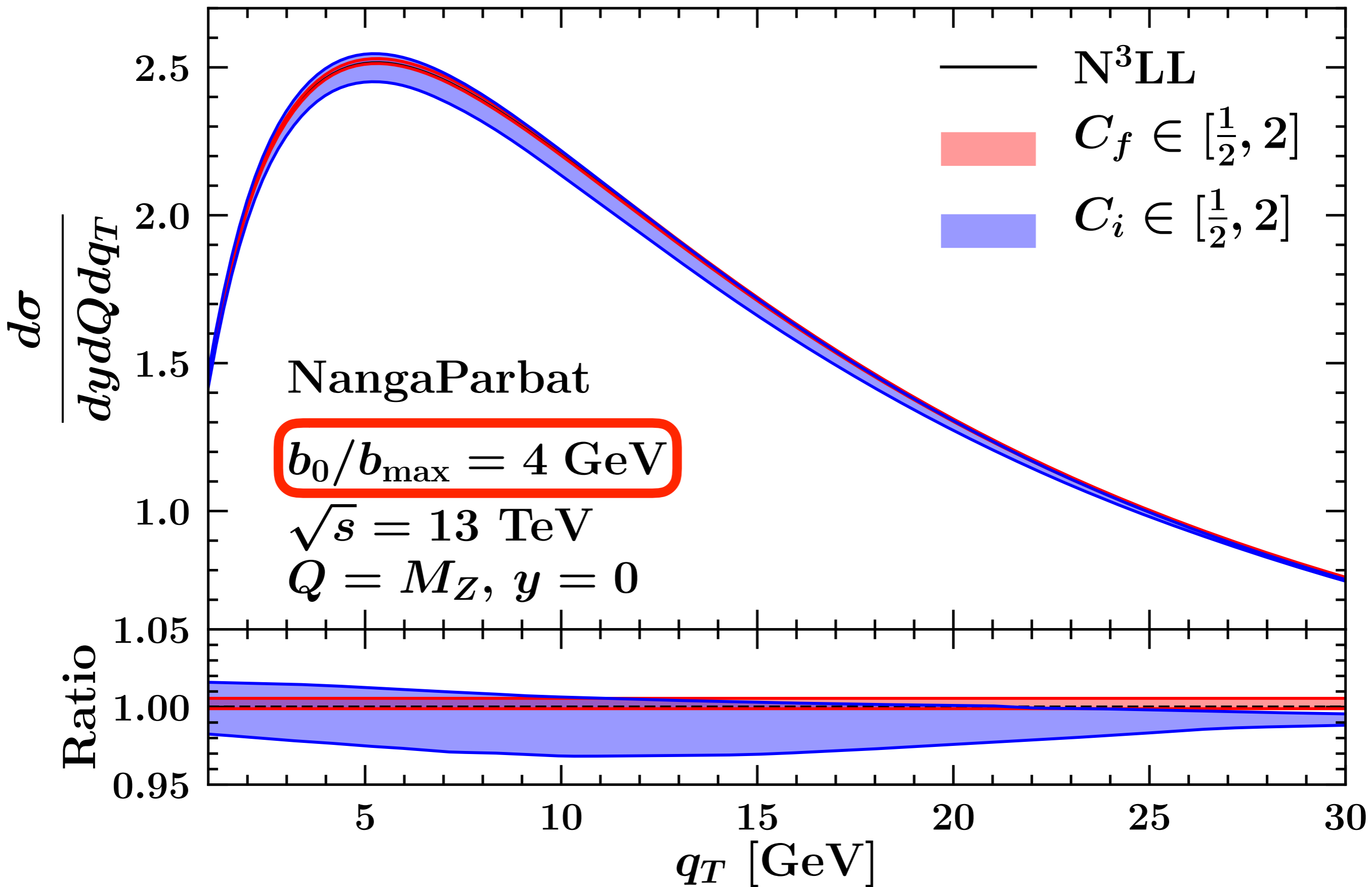
🍏 if the scale variation is a factor 1/2, then $\mu_{min} = 0.5 \text{ GeV}$ α_s where is very large and **PDFs are crazy**.



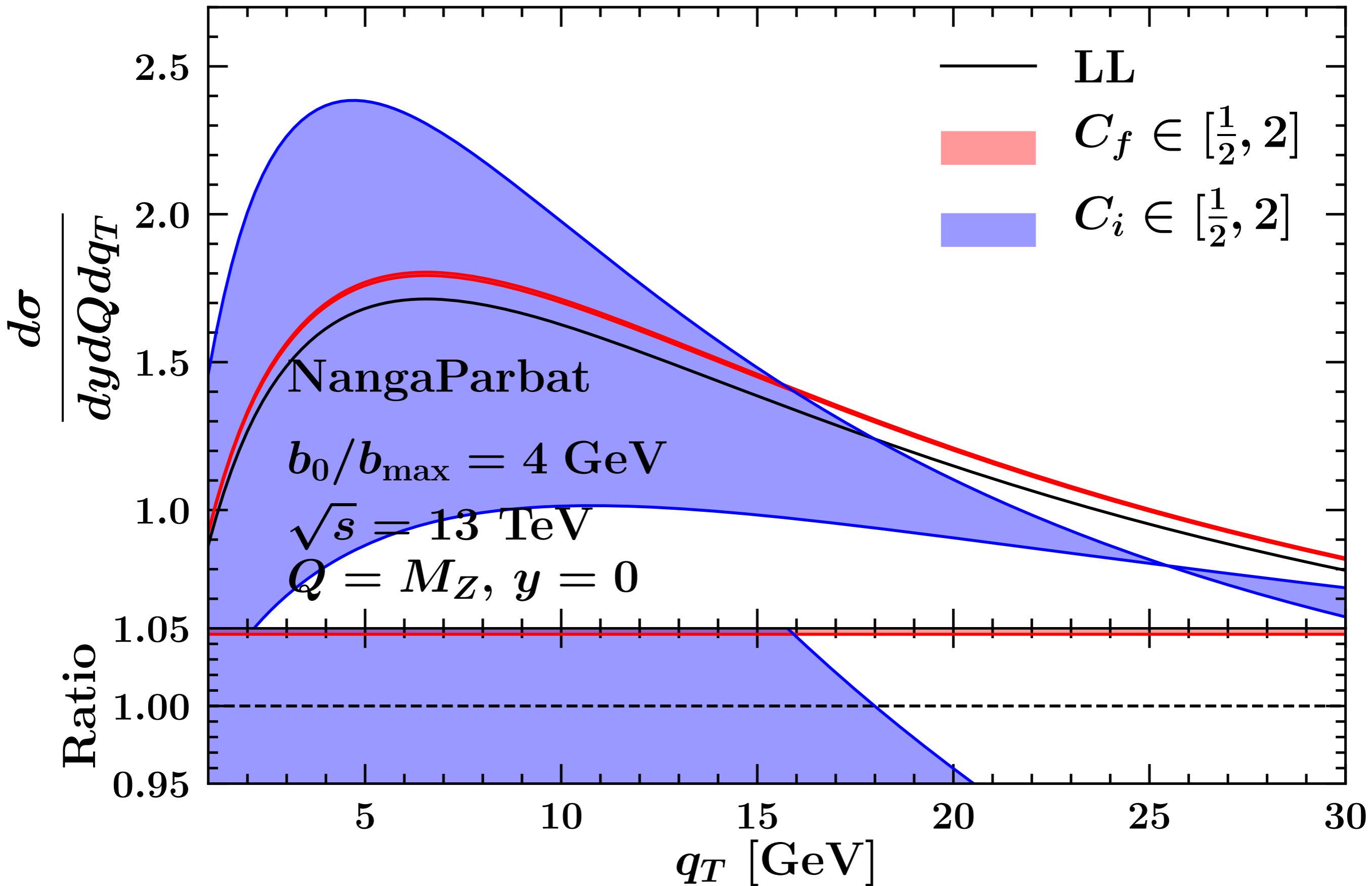
Numerical effect



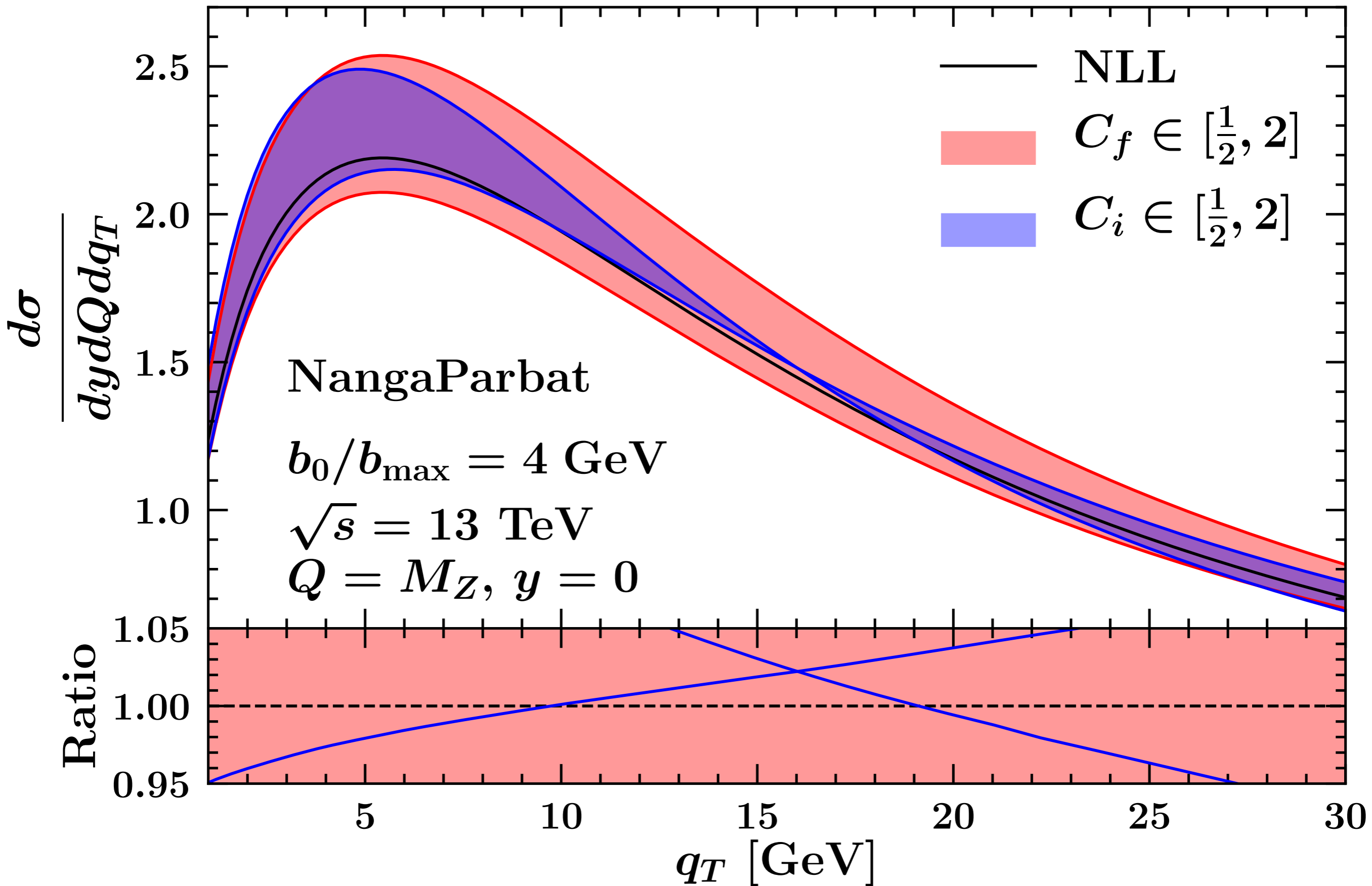
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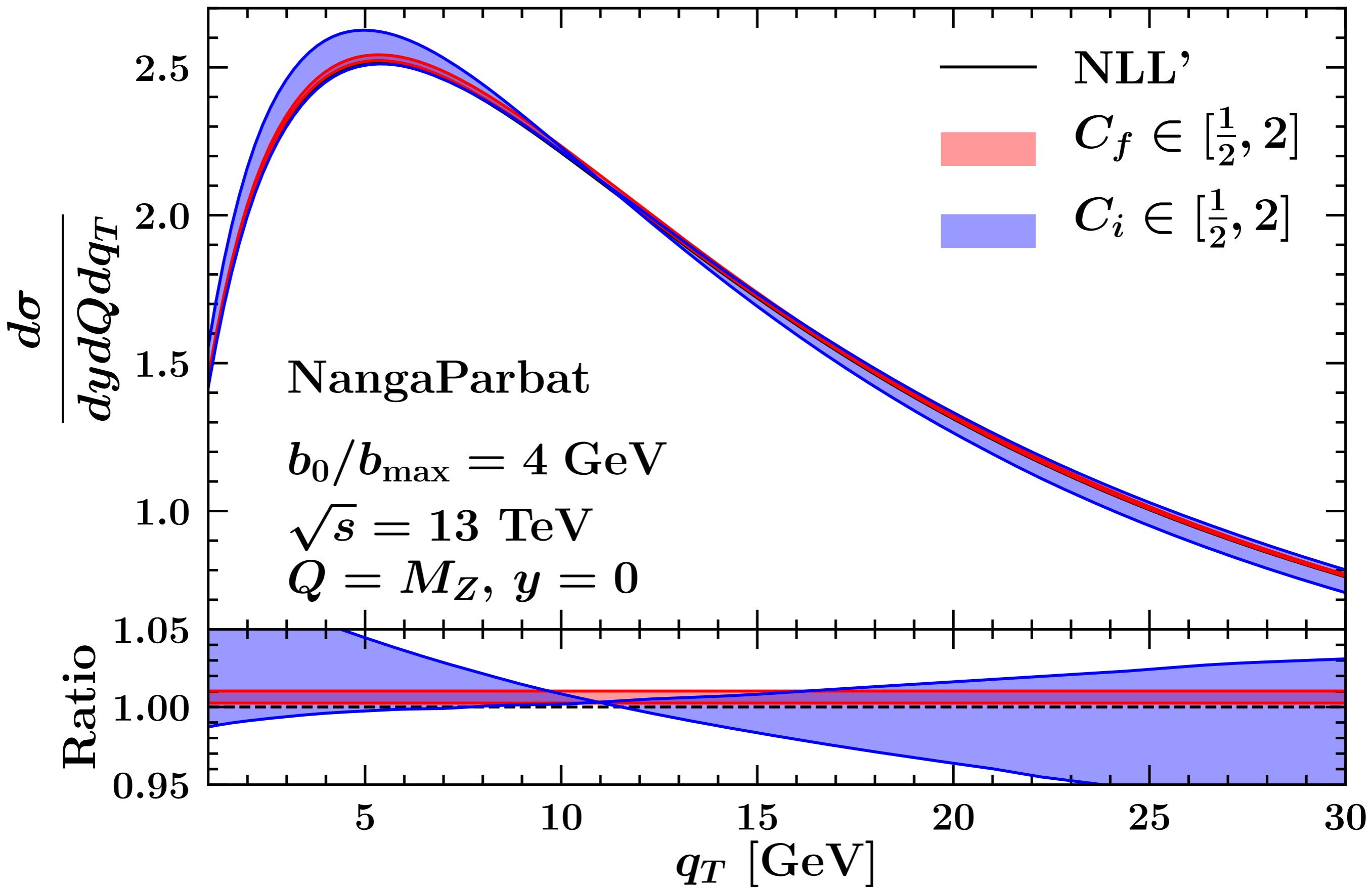
Perturbative convergence



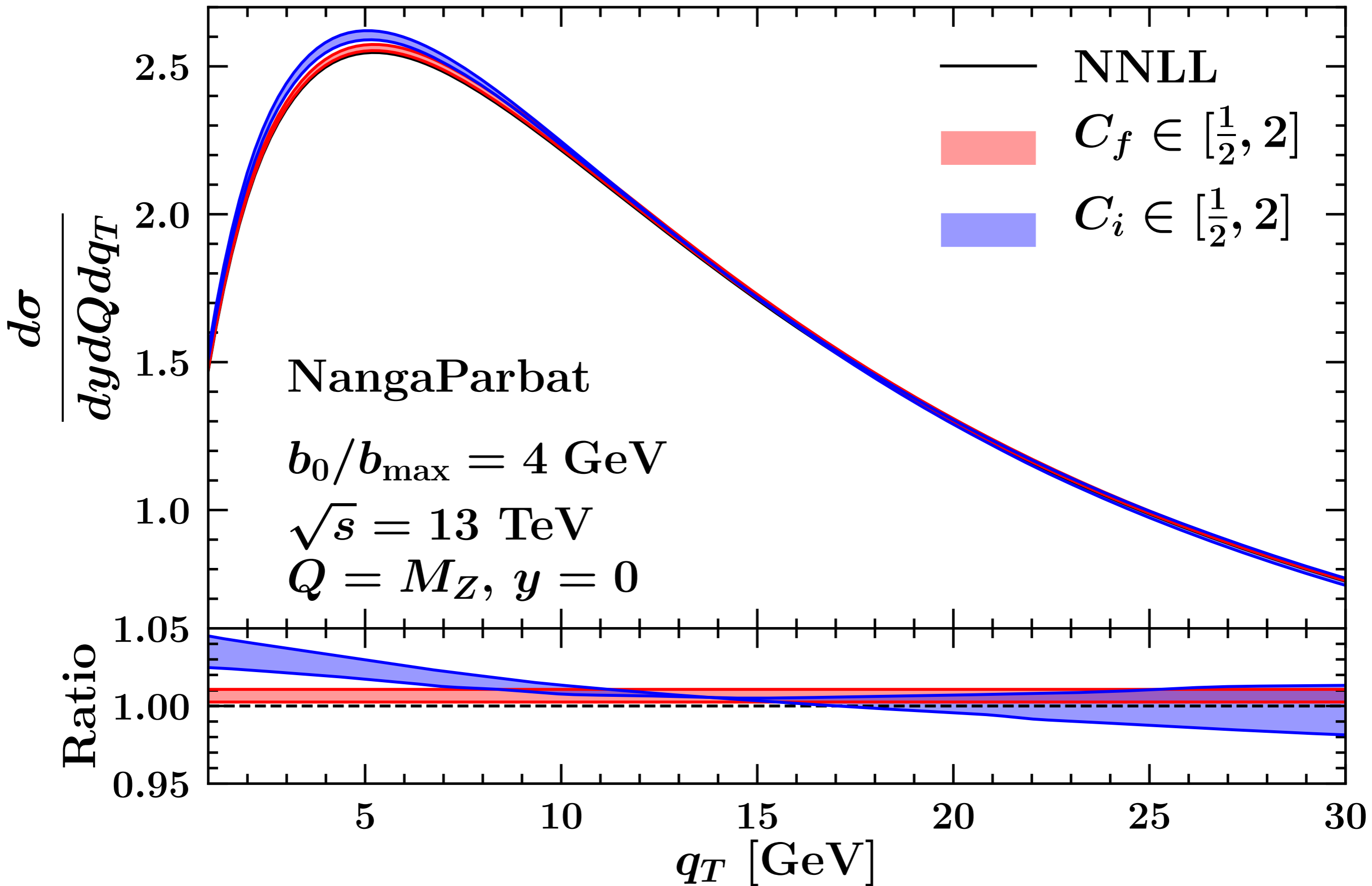
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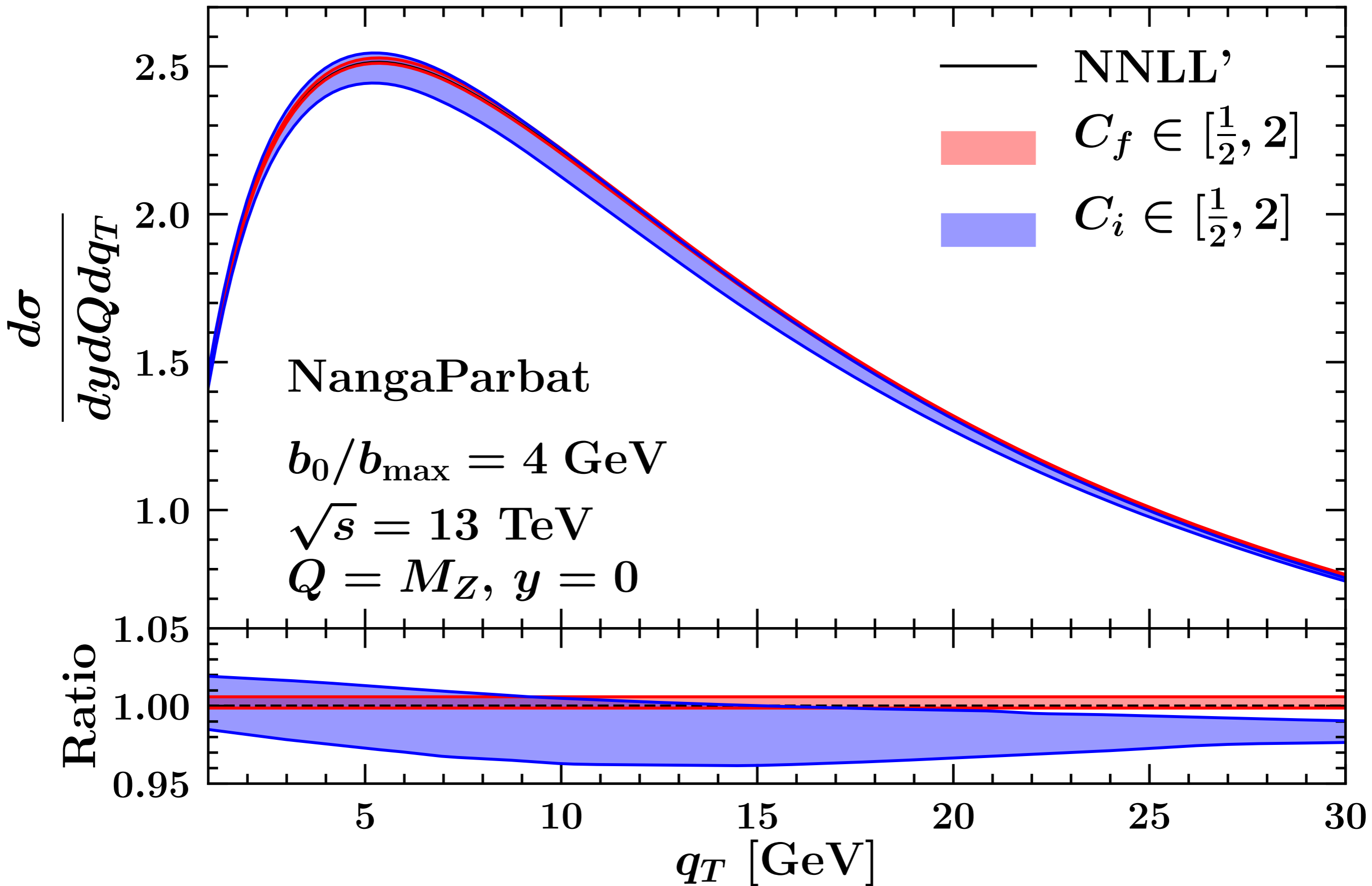
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