

Early Isotropization of the Quark Gluon Plasma

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Viscous Hydrodynamics

I) Macroscopic theory

II) Few parameters: $P_L, P_T, \epsilon, \vec{u}$

III) Need input:

1) Equation of state $f(P_L, P_T) = \epsilon$

2) Small anisotropy

3) Initialization: $\epsilon(\tau_0), P_L(\tau_0)? \dots$

4) viscous coefficients: shear viscosity η, \dots

5) Short isotropization time

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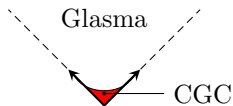
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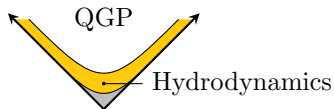
**None of this is easy
to get from QCD**

Early transition: the problem



Huge anisotropy
(negative P_L)

**Isotropization?
Time scale?**



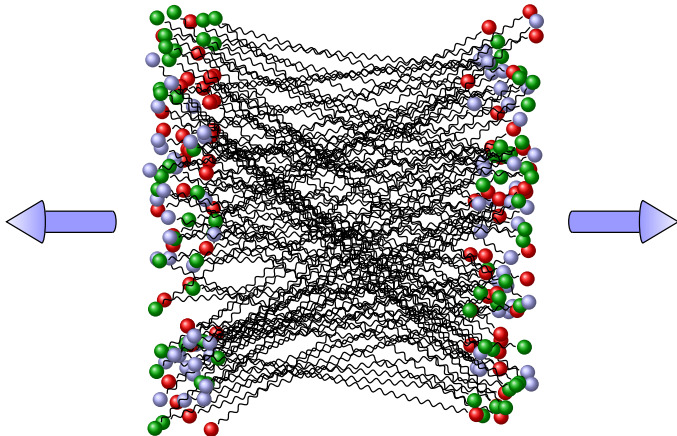
Small anisotropy

Long time puzzle: Does (fast) isotropization occur?

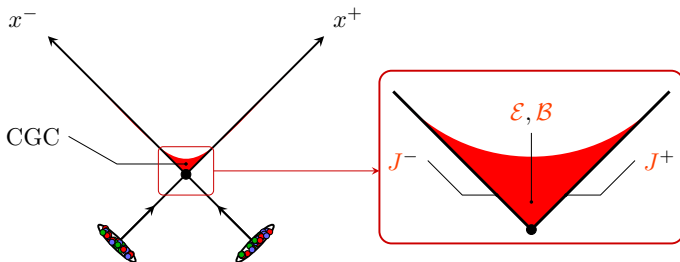
HOW TO STUDY THE TRANSITION?

Weakly coupled method at high density:

$$\alpha_s \ll 1 \text{ but } f_{\text{gluon}} \sim \frac{1}{\alpha_s}$$



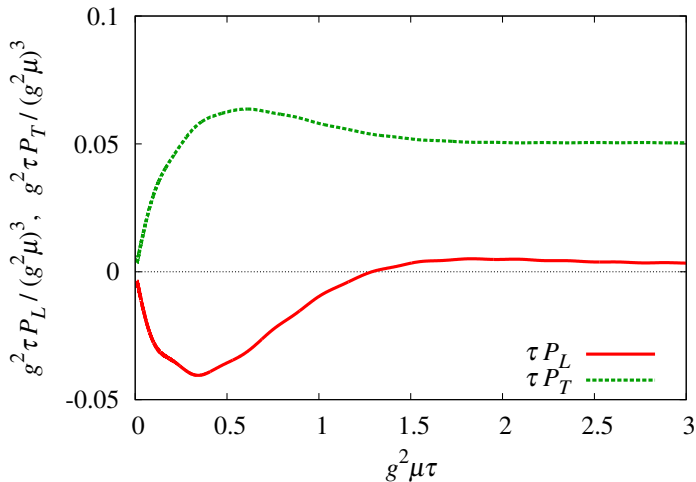
Theoretical framework (Weakly coupled but strongly interacting)



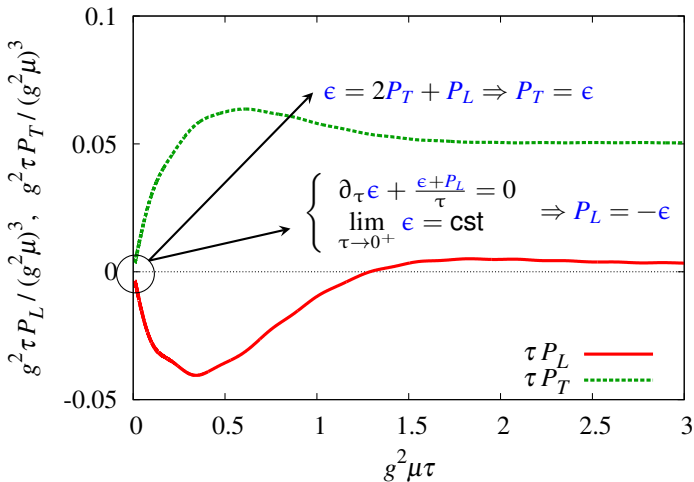
LO:
$$\epsilon = \frac{1}{2} \underbrace{(\vec{\mathcal{E}}^2 + \vec{\mathcal{B}}^2)}_{\text{Classical color fields}}$$

$$\mathcal{D}_\mu \mathcal{F}^{\mu\nu} = \underbrace{J^\nu}_{\text{Color sources on the light cone}}$$

[KRASNITZ, VENUGOPALAN (1998)]

Strong anisotropy at all time

Strong anisotropy at all time



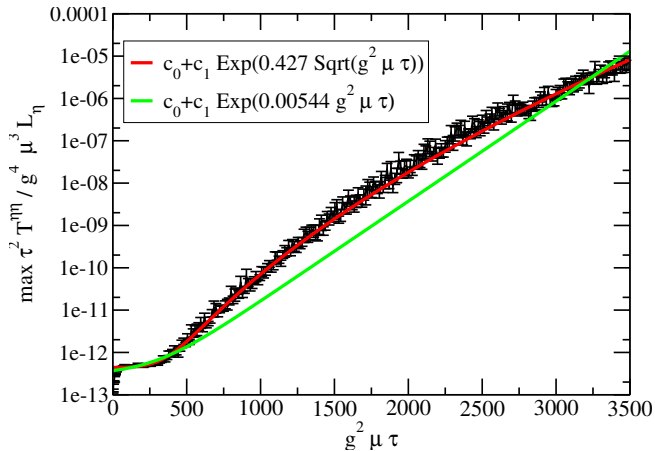
THE COLOR GLASS CONDENSATE AT NLO

$$E^2(x) = \underbrace{\mathcal{E}^2(\mathbf{x}_\perp)}_{\text{LO}} + \underbrace{\frac{1}{2} \int_{\vec{k}} |e_{\vec{k}}(x)|^2}_{\text{NLO}} + \dots$$

$e_{\vec{k}}(x)$ perturbation to $\mathcal{E}(x)$ created by a plane wave of momentum \vec{k} in the remote past.

Obtained by solving the linearized equation of motions.

THE COLOR GLASS CONDENSATE AT NLO

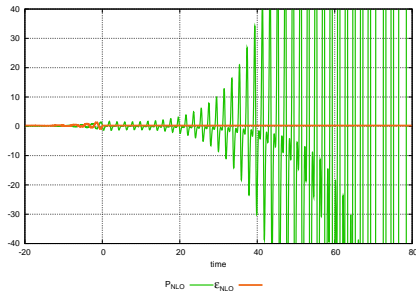
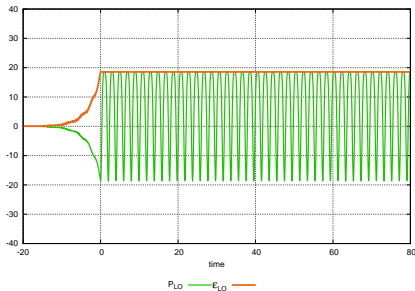


[ROMATSCHKE, VENUGOPALAN (2006)]

Small Fluctuations grow exponentially (Weibel instability)

THE COLOR GLASS CONDENSATE AT NLO

- Because of instabilities, the **NLO** correction eventually becomes as large as the **LO** \Rightarrow Important effect, should be included
- **NLO** alone will grow forever \Rightarrow unphysical effect, should be taken care of



- Such growing contributions are present at all orders of the perturbative expansion

How to deal with them?

- At the initial time $\tau = \tau_0$, take:

$$\vec{E}_0(\tau_0, \vec{x}) = \vec{\mathcal{E}}_0(\tau_0, \vec{x}) + \int_{\vec{k}} c_{\vec{k}} \vec{e}_{\vec{k}}(\tau_0, \vec{x})$$

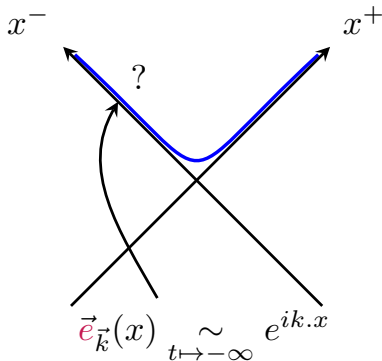
where $c_{\vec{k}}$ are random coefficients: $\langle c_{\vec{k}} c_{\vec{k}'} \rangle \sim \delta_{\vec{k}\vec{k}'}$

- Solve the **Classical** equation of motion $D_{\mu} F^{\mu\nu} = J^{\nu}$
- Compute $\langle \vec{E}^2(\tau, \vec{x}) \rangle$, where $\langle \rangle$ is the average on the $c_{\vec{k}}$ (Monte-Carlo)
- One can show that this resums all the fastest growing terms at each order, leading to a result that remains bounded when $\tau \rightarrow \infty$
[GELIS, LAPPI, VENUGOPALAN (2008)]

This gives: LO+NLO+Subset of higher orders

THE NLO SPECTRUM

- Need to know $\vec{e}_{\vec{k}}(\tau_0, \vec{x})$ at the time τ_0 we start the numerical simulation
- For practical reasons, we must start in the forward light cone ($\tau_0 > 0$)



This can be done analytically [TE,GELIS 1307:1765]

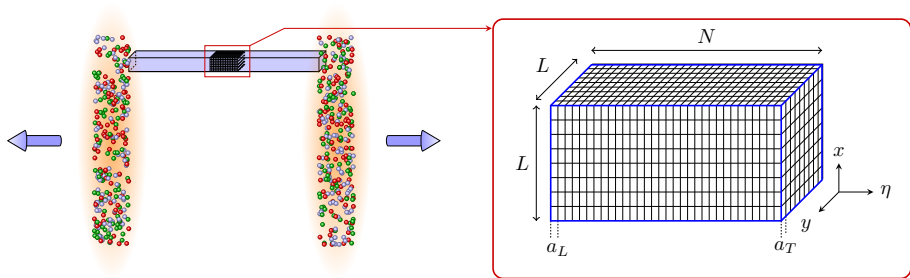
Result

$$e_{\sqrt{k}_\perp}^i(\tau, \mathbf{x}_\perp, \eta) = i\nu e^{i\nu\eta} \left[F_{\sqrt{k}_\perp}^{i,2}(\mathcal{U}_2, \tau, \mathbf{x}_\perp) - F_{\sqrt{k}_\perp}^{i,1}(\mathcal{U}_1, \tau, \mathbf{x}_\perp) \right]$$

$$e_{\sqrt{k}_\perp}^\eta(\tau, \mathbf{x}_\perp, \eta) = e^{i\nu\eta} \mathcal{D}^i \left[F_{\sqrt{k}_\perp}^{i,2}(\mathcal{U}_2, \tau, \mathbf{x}_\perp) - F_{\sqrt{k}_\perp}^{i,1}(\mathcal{U}_1, \tau, \mathbf{x}_\perp) \right]$$

- \mathcal{U}_1 depends on the color source J^+ of the first nucleus
- \mathcal{U}_2 depends on the color source J^- of the second nucleus
- Analytical checks performed on the solution
 - Gauss's law
 - linearized Yang-Mills EOM
 - Orthonormality of the mode functions

Gauge potential $A^\mu \rightarrow$ link variables (exact gauge invariance on the lattice)

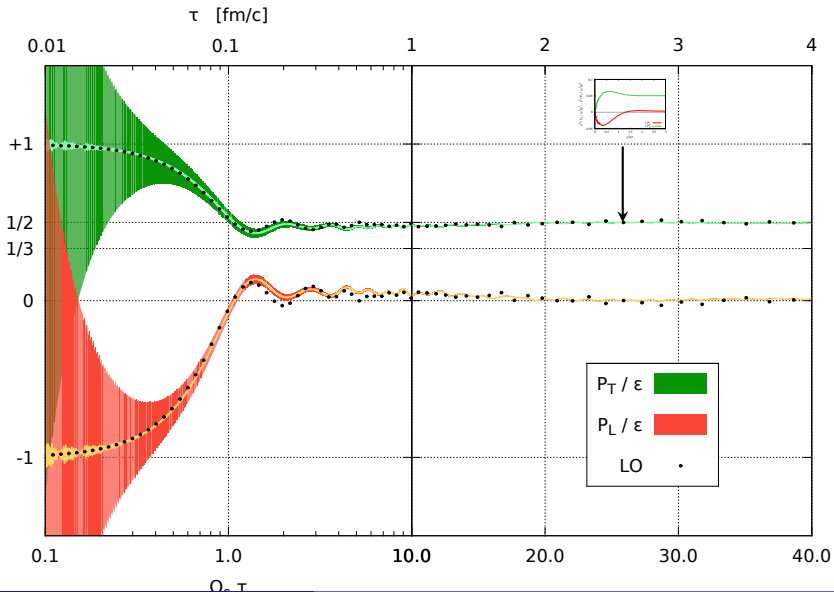


Numerical parameters

- Transverse lattice size $L = 64$, transverse lattice spacing $Q_s a_T = 1$
- Longitudinal lattice size $N = 128$, longitudinal lattice spacing $a_L = 0.016$
- Number of configurations for the Monte-Carlo $N_{\text{conf}} = 200$ to 2000
- Initial time $Q_s \tau_0 = 0.01$

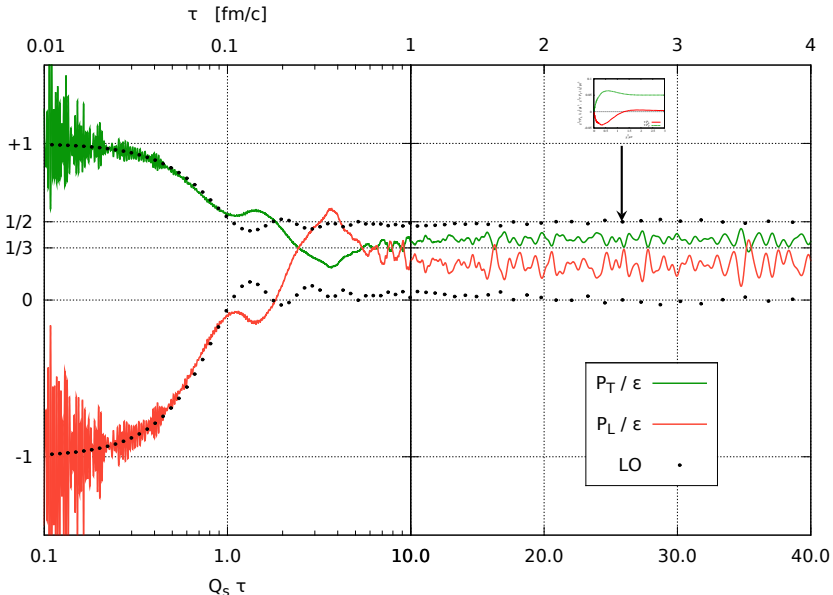
NUMERICAL RESULTS [TE,GELIS (2013)]

$$\alpha_s = 8 \cdot 10^{-4} \quad (g = 0.1)$$



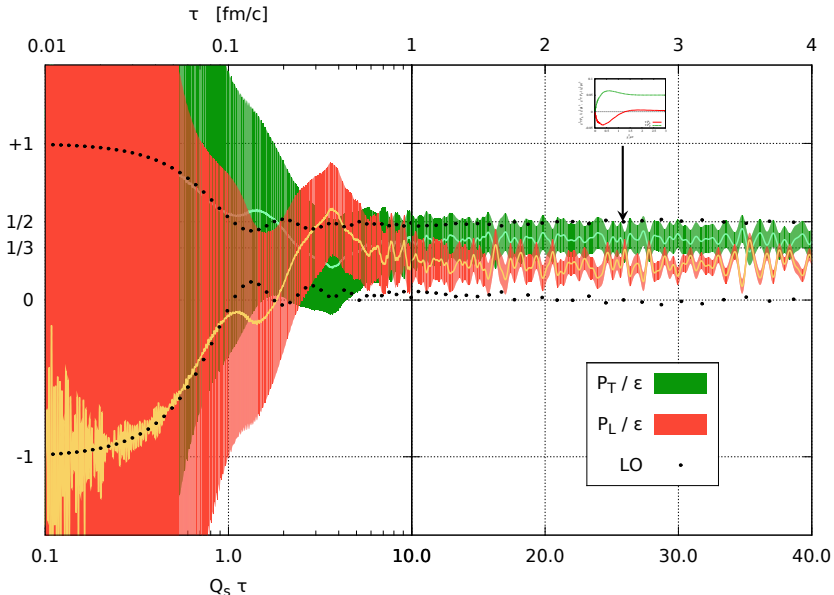
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ANOMALOUSLY SMALL VISCOSITY

Assuming simple first order viscous hydrodynamics

$$\epsilon \approx \underbrace{\epsilon_0 \tau^{-\frac{4}{3}}}_{\text{Ideal hydro}} - \underbrace{2\eta_0 \tau^{-2}}_{\text{first order correction}}$$

we can compute the dimensionless ratio ($\eta = \eta_0 \tau^{-1}$)

$$\eta \epsilon^{-\frac{3}{4}} \lesssim 1$$

In contrast, perturbation theory at LO gives $\eta \epsilon^{-\frac{3}{4}} \sim 300$.

If the system is nearly thermal

$$\epsilon^{\frac{3}{4}} \sim s \implies \frac{\eta}{s} \text{ close to } \frac{1}{4\pi}$$

CONCLUSION

- Correct NLO spectrum from first principles
- Fixed anisotropy for $g = 0.5$ at $\tau \sim 1fm/c$
- Assuming first order viscous hydrodynamics: $\eta\epsilon^{-\frac{3}{4}} \lesssim 1$
- Compatible with viscous hydrodynamical expansion
- No need for strong coupling to get hydrodynamization

BACKUP: COMPLETELY INCOHERENT INITIAL FIELDS VS CGC

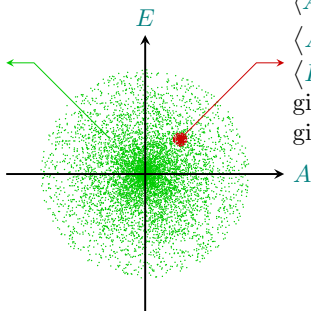
$$\langle A \rangle \sim 0, \langle E \rangle \sim 0$$

$$\langle A^2 \rangle - \langle A \rangle^2 \sim \frac{Q_s^2}{g^2}$$

$$\langle E^2 \rangle - \langle E \rangle^2 \sim \frac{Q_s^4}{g^2}$$

May give correct answer at LO

Not correct at NLO



$$\langle A \rangle \sim \frac{Q_s}{g}, \langle E \rangle \sim \frac{Q_s^2}{g}$$

$$\langle A^2 \rangle - \langle A \rangle^2 \sim Q_s^2$$

$$\langle E^2 \rangle - \langle E \rangle^2 \sim Q_s^4$$

give correct answer at LO

give correct answer at NLO

EOM ON A LATTICE

Writing

$$E^\mu(x) = \overset{x}{\underset{\mu}{\bullet}} \quad U_\mu(x) = \overset{x}{\bullet} \xrightarrow{\hat{\mu}} \quad U_\mu^\dagger(x) = \overset{\hat{\mu}}{\bullet} \xleftarrow{x} + \hat{\mu}$$

and

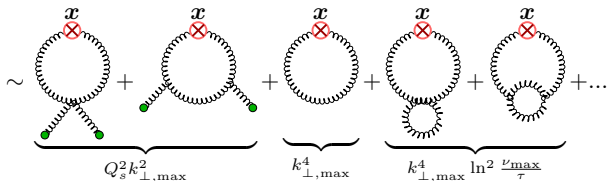
$$U_{\mu\nu}(x) = \begin{array}{c} \text{---} \\ \downarrow \\ \text{---} \\ \uparrow \\ \text{---} \end{array} \begin{array}{c} \hat{\nu} \\ \hat{\mu} \end{array} \quad U_{\mu\nu}^\dagger(x) = \begin{array}{c} \hat{\mu} \\ \hat{\nu} \end{array} \begin{array}{c} \text{---} \\ \downarrow \\ \text{---} \\ \uparrow \\ \text{---} \end{array} \\ U_{\mu-\nu}(x) = \begin{array}{c} \text{---} \\ \uparrow \\ \text{---} \\ \downarrow \\ \text{---} \end{array} \begin{array}{c} \hat{\mu} \\ \hat{\nu} \end{array} \quad U_{\mu-\nu}^\dagger(x) = \begin{array}{c} \hat{\nu} \\ \hat{\mu} \end{array} \begin{array}{c} \text{---} \\ \uparrow \\ \text{---} \\ \downarrow \\ \text{---} \end{array} .$$

We can therefore rewrite the EOM as

$$\partial_\tau \overset{x}{\underset{i}{\bullet}} = \frac{-i}{2ga_I a_J^2} \sum_J \left[\begin{array}{c} \text{---} \\ \downarrow \\ \text{---} \\ \uparrow \\ \text{---} \end{array} \begin{array}{c} \hat{i} \\ \hat{j} \end{array} - \begin{array}{c} \hat{j} \\ \hat{i} \end{array} \begin{array}{c} \text{---} \\ \downarrow \\ \text{---} \\ \uparrow \\ \text{---} \end{array} + \begin{array}{c} \text{---} \\ \downarrow \\ \text{---} \\ \uparrow \\ \text{---} \end{array} \begin{array}{c} \hat{j} \\ \hat{i} \end{array} - \begin{array}{c} \hat{i} \\ \hat{j} \end{array} \begin{array}{c} \text{---} \\ \downarrow \\ \text{---} \\ \uparrow \\ \text{---} \end{array} \right] \\ \partial_\tau \overset{x}{\bullet} \xrightarrow{i} = -i g a_I \overset{x}{\underset{i}{\bullet}} \xrightarrow{i}$$

RENORMALIZATION PROCEDURE

$$\langle E_{L,\text{div}}^2 \rangle \sim Q_s^2 k_{\perp,\text{max}}^2 + k_{\perp,\text{max}}^4 + k_{\perp,\text{max}}^4 \ln^2 \frac{\nu_{\text{max}}}{\tau} + \dots$$



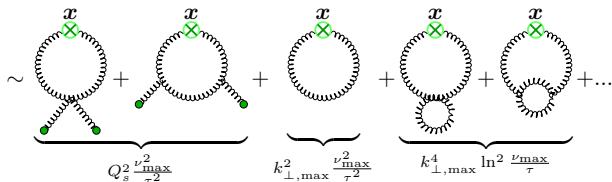
3 last diagrams can be subtracted with a simulation where

$$A_{\mu}^a(x) = 0 + a_{\mu}^a(x)$$

$$E_L^2 \text{ "fine" } (B_L^2 \text{ too})$$

RENORMALIZATION PROCEDURE

$$\langle E_{T,\text{div}}^2 \rangle \sim Q_s^2 \frac{\nu_{\text{max}}^2}{\tau^2} + k_{\perp,\text{max}}^2 \frac{\nu_{\text{max}}^2}{\tau^2} + k_{\perp,\text{max}}^4 \ln^2 \frac{\nu_{\text{max}}}{\tau} + \dots$$



3 last diagrams can be subtracted with a simulation where

$$A_{\mu}^a(x) = 0 + a_{\mu}^a(x)$$

How to deal with the first 2? \rightarrow fit for the time being.

Otherwise E_T^2 and B_T^2 behaves as τ^{-2} at early time.

RENORMALIZATION PROCEDURE

$$\epsilon = E_T^2 + B_T^2 + \underbrace{E_L^2}_{\text{fine}} + \underbrace{B_L^2}_{\text{fine}}$$

$$P_T = \underbrace{E_L^2}_{\text{fine}} + \underbrace{B_L^2}_{\text{fine}}$$

$$P_L = E_T^2 + B_T^2 - \underbrace{E_L^2}_{\text{fine}} - \underbrace{B_L^2}_{\text{fine}}$$

RENORMALIZATION PROCEDURE

$$\begin{aligned}
 \langle P_T \rangle_{\text{phys.}} &= \langle P_T \rangle_{\substack{\text{backgd.} \\ + \text{fluct.}}} - \langle P_T \rangle_{\substack{\text{fluct.} \\ \text{only}}} \\
 \langle \epsilon, P_L \rangle_{\text{phys.}} &= \underbrace{\langle \epsilon, P_L \rangle_{\substack{\text{backgd.} \\ + \text{fluct.}}}}_{\text{computed}} - \underbrace{\langle \epsilon, P_L \rangle_{\substack{\text{fluct.} \\ \text{only}}}}_{\text{computed}} + \underbrace{A \tau^{-2}}_{\text{fitted}} .
 \end{aligned}$$

τ^- term only one to satisfy Bjorken law and EOS:

$$\partial_\tau \tau^{-\alpha} + 2\tau^{-\alpha-1} = 0$$

RENORMALIZATION PROCEDURE

How come that problematic divergent diagrams behaves as mass terms?

In the continuum limit, they don't exist local gauge invariant operators of dimension two.

On the lattice though, they could be terms like

$$g^2 \frac{v_{\max}^2}{k_{\perp, \max}^2 \tau^2} \text{Tr} F^2,$$

where

$$F_{\mu\nu}(x) \sim \begin{array}{c} \left[\begin{array}{c} \leftarrow \\ \downarrow \\ \rightarrow \\ \uparrow \end{array} \right]_{\hat{\mu}}^{\hat{\nu}} - \begin{array}{c} \left[\begin{array}{c} \rightarrow \\ \downarrow \\ \leftarrow \\ \uparrow \end{array} \right]_{\hat{\nu}}^{\hat{\mu}} \end{array}$$