

Pre-equilibrium dynamics in heavy ion collisions

I – Parton evolution at small x , Saturation



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Phenomenology of saturation

General introduction



Stages of a nucleus-nucleus collision

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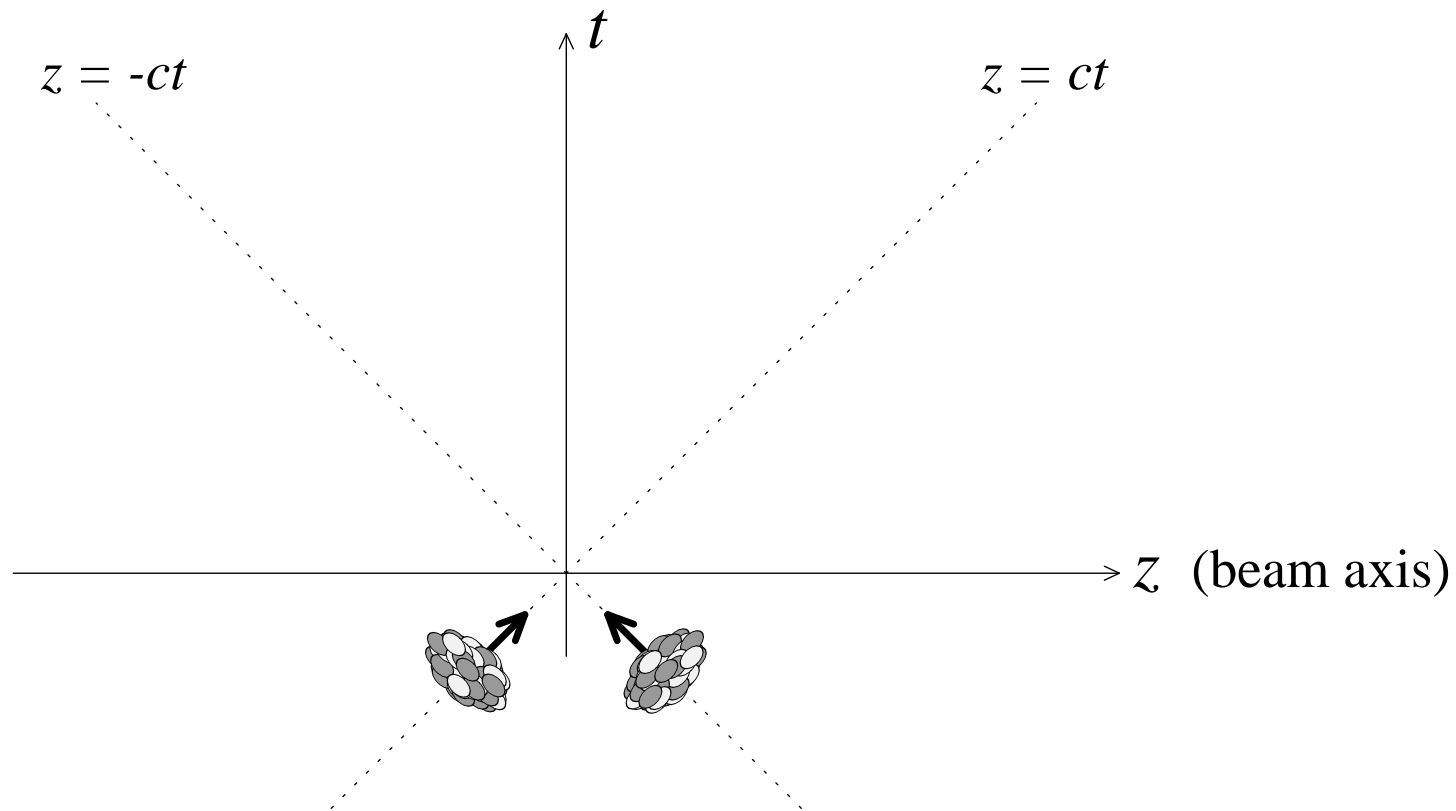
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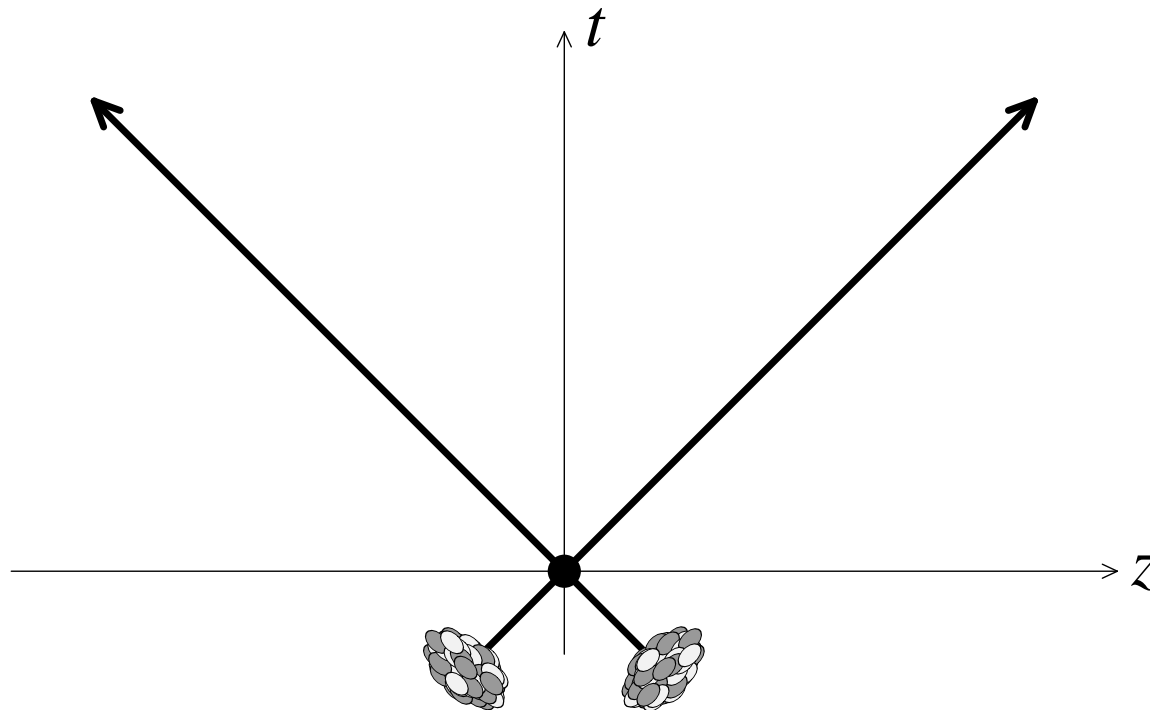
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- $\tau \sim 0 \text{ fm}/c$
- Production of hard particles :
 - ◆ jets, direct photons
 - ◆ heavy quarks
- calculable with perturbative QCD (leading twist)

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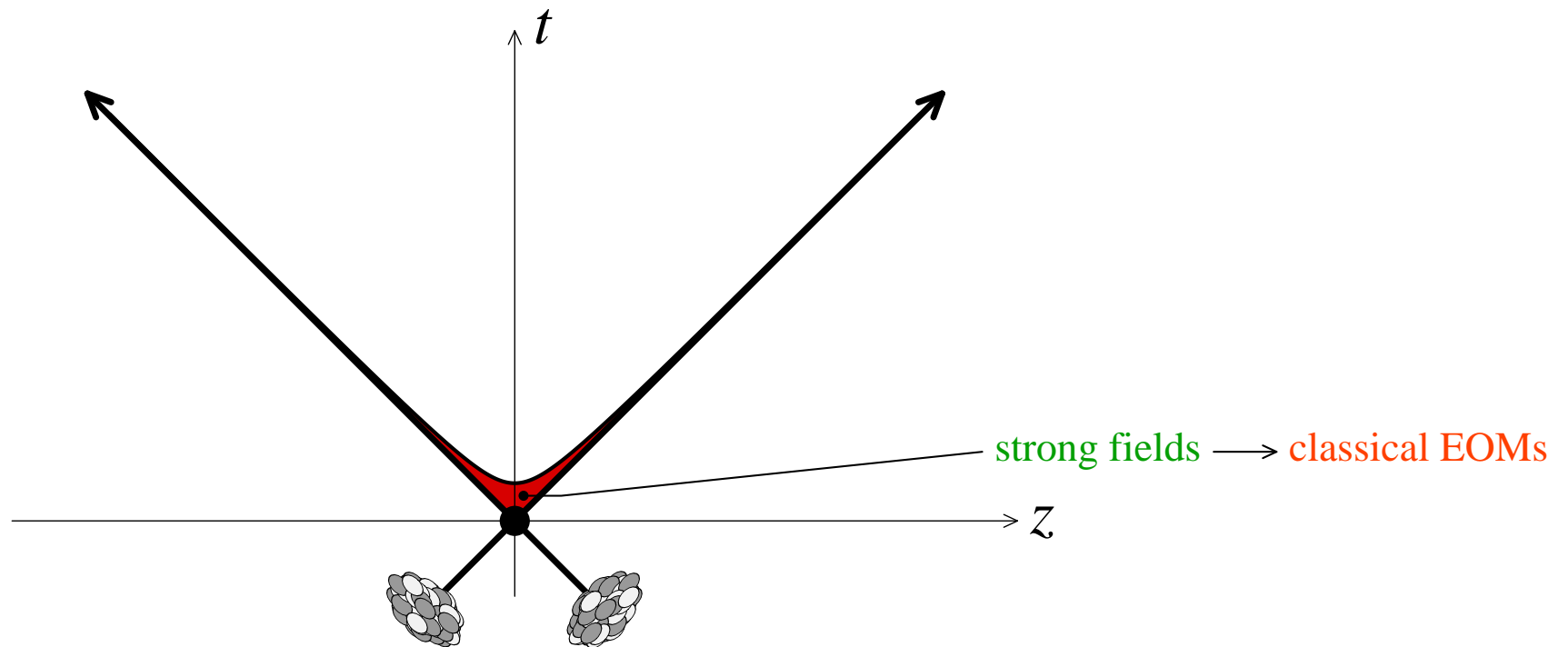
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- $\tau \sim 0.2 \text{ fm}/c$
- Production of semi-hard particles : gluons, light quarks
- relatively small momentum : $p_{\perp} \lesssim 2\text{--}3 \text{ GeV}$
- make up for most of the multiplicity
- sensitive to the physics of saturation (higher twist)

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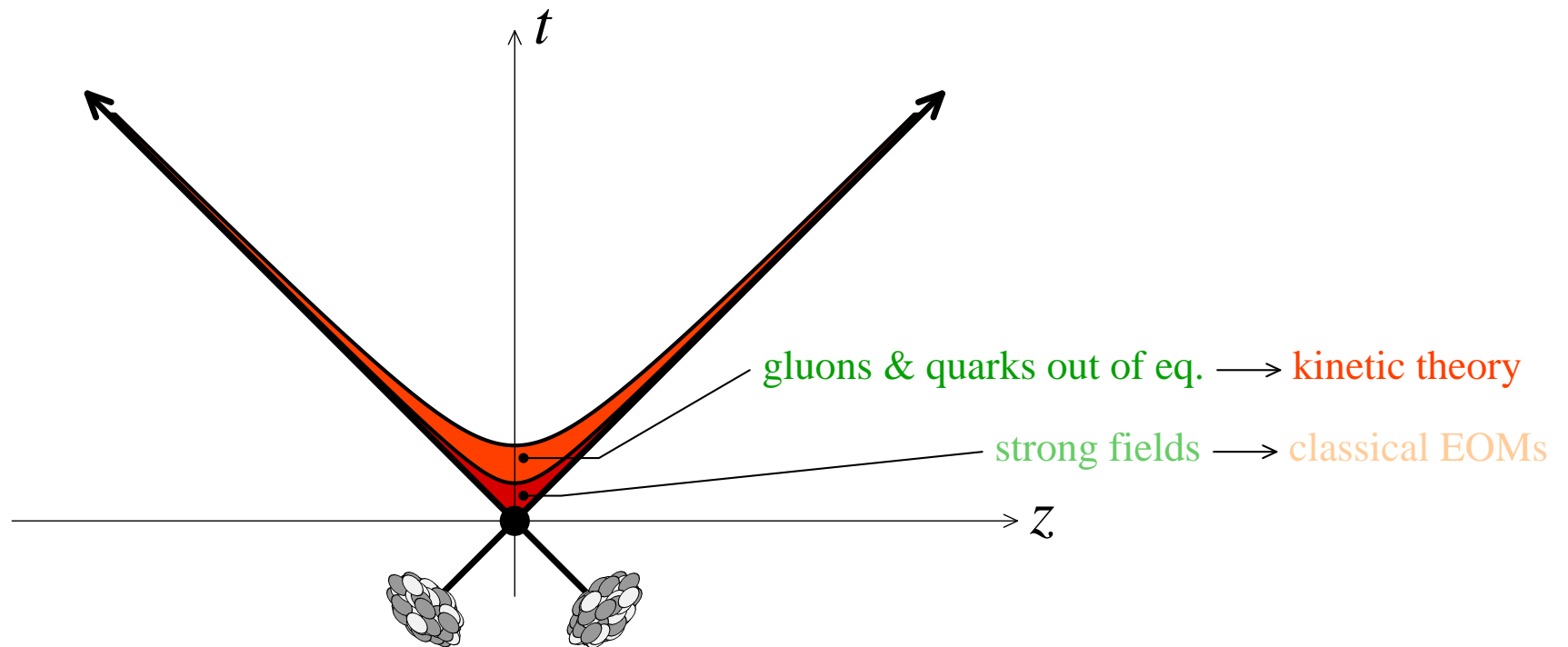
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- $\tau \sim 1-2 \text{ fm}/c$
- **Thermalization**
 - ◆ experiments suggest a fast thermalization
 - ◆ but this is still not understood from QCD

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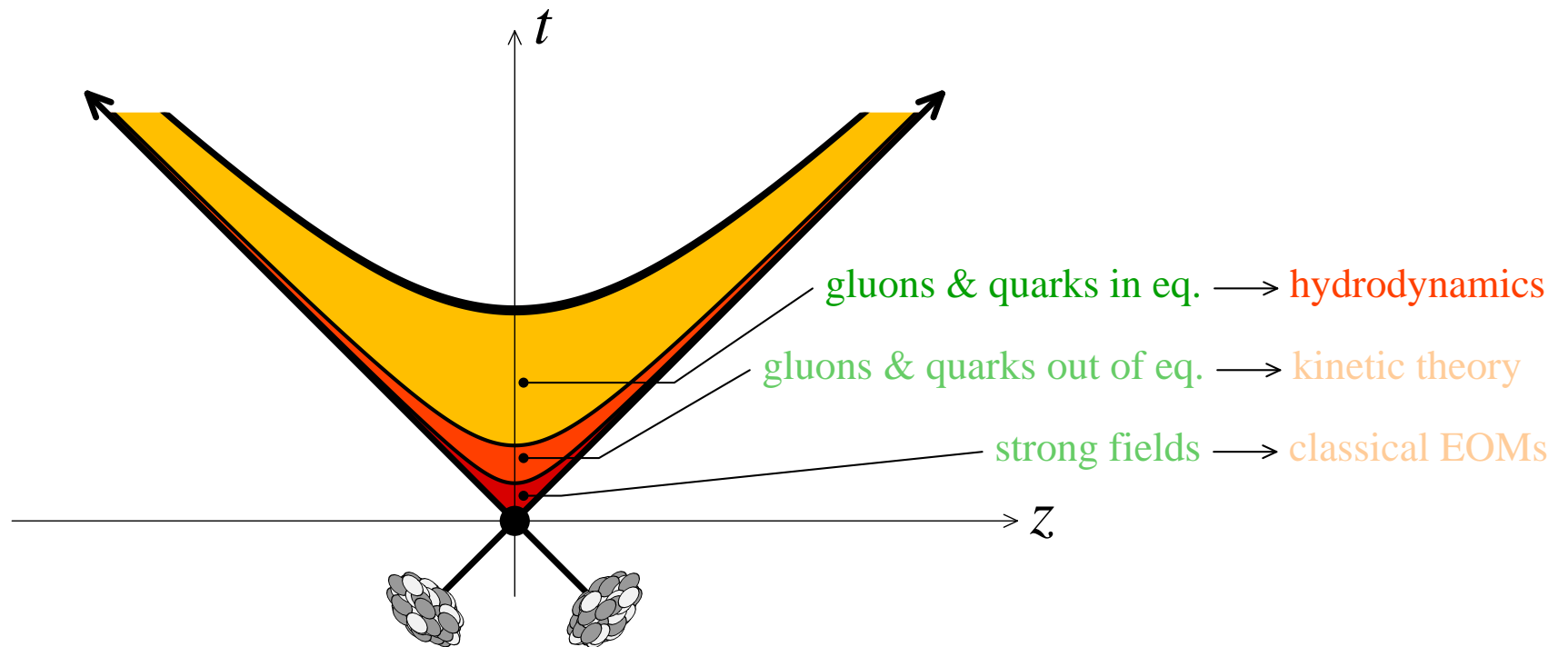
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- $2 \leq \tau \lesssim 10 \text{ fm}/c$
- Quark gluon plasma

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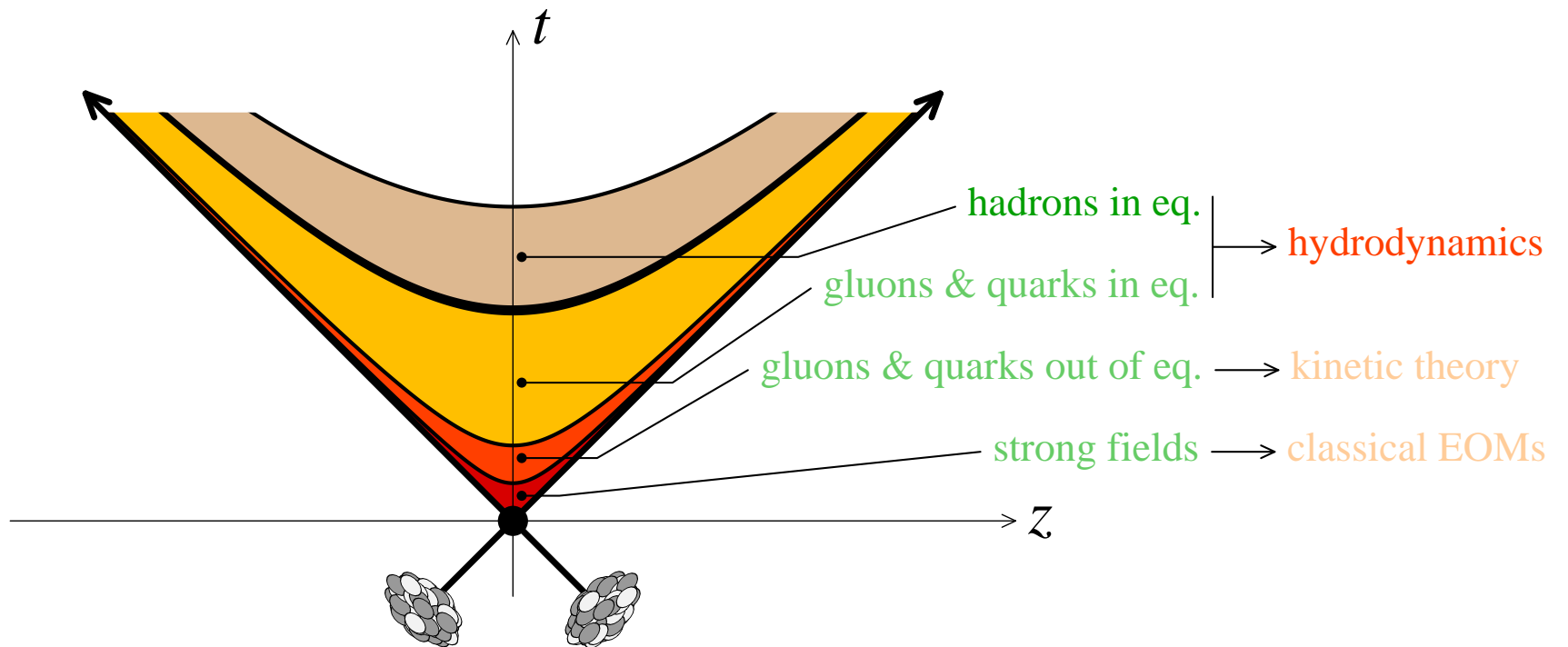
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- $10 \lesssim \tau \lesssim 20 \text{ fm}/c$
- Hot hadron gas

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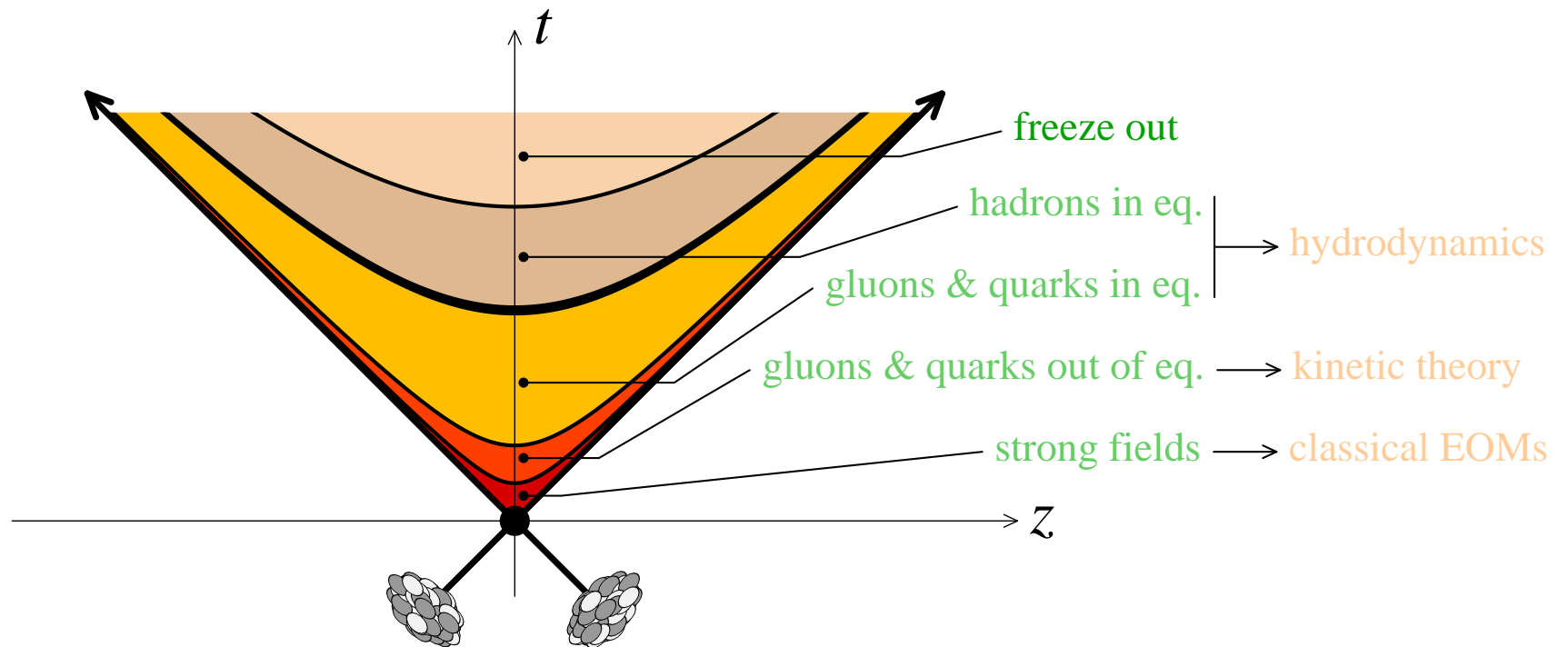
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- $\tau \rightarrow +\infty$
- **Chemical freeze-out :**
density too small to have inelastic interactions
- **Kinetic freeze-out :**
no more elastic interactions

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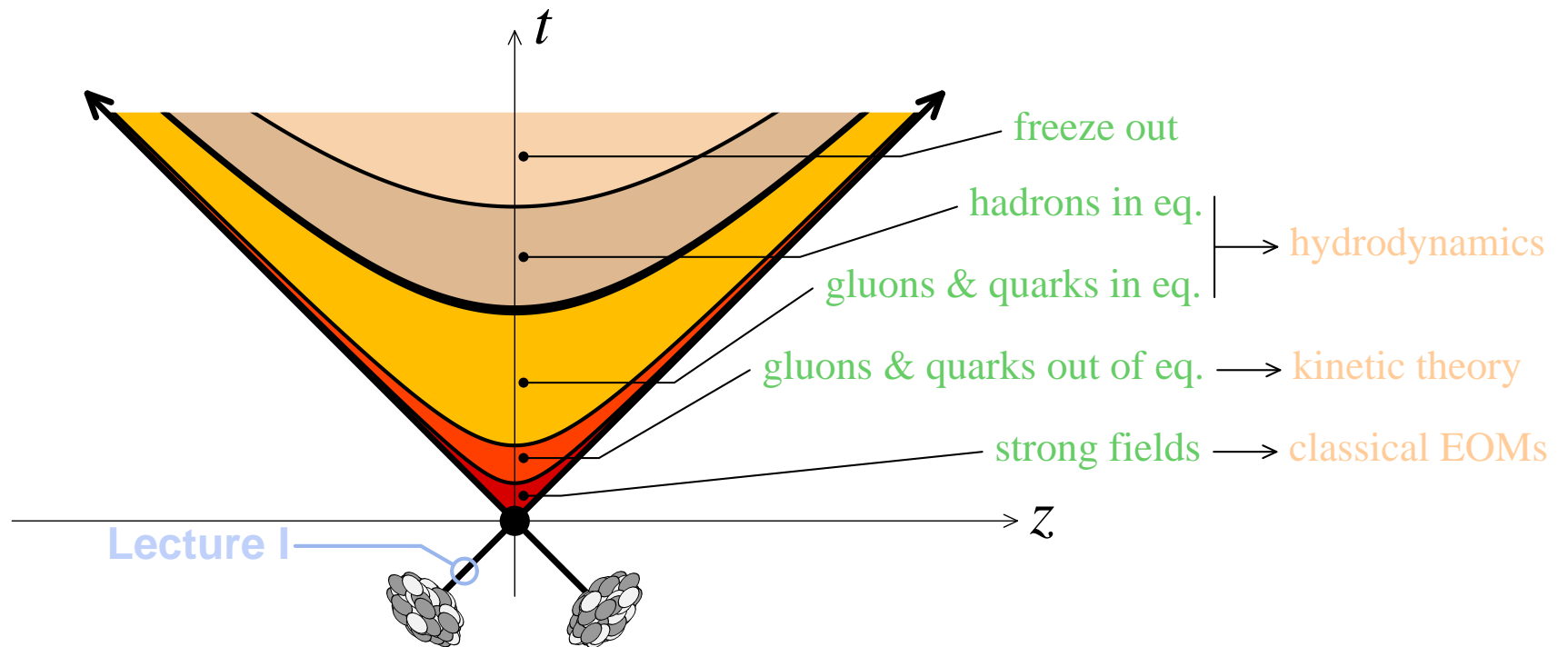
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- Lecture I : Parton evolution at small x , Saturation

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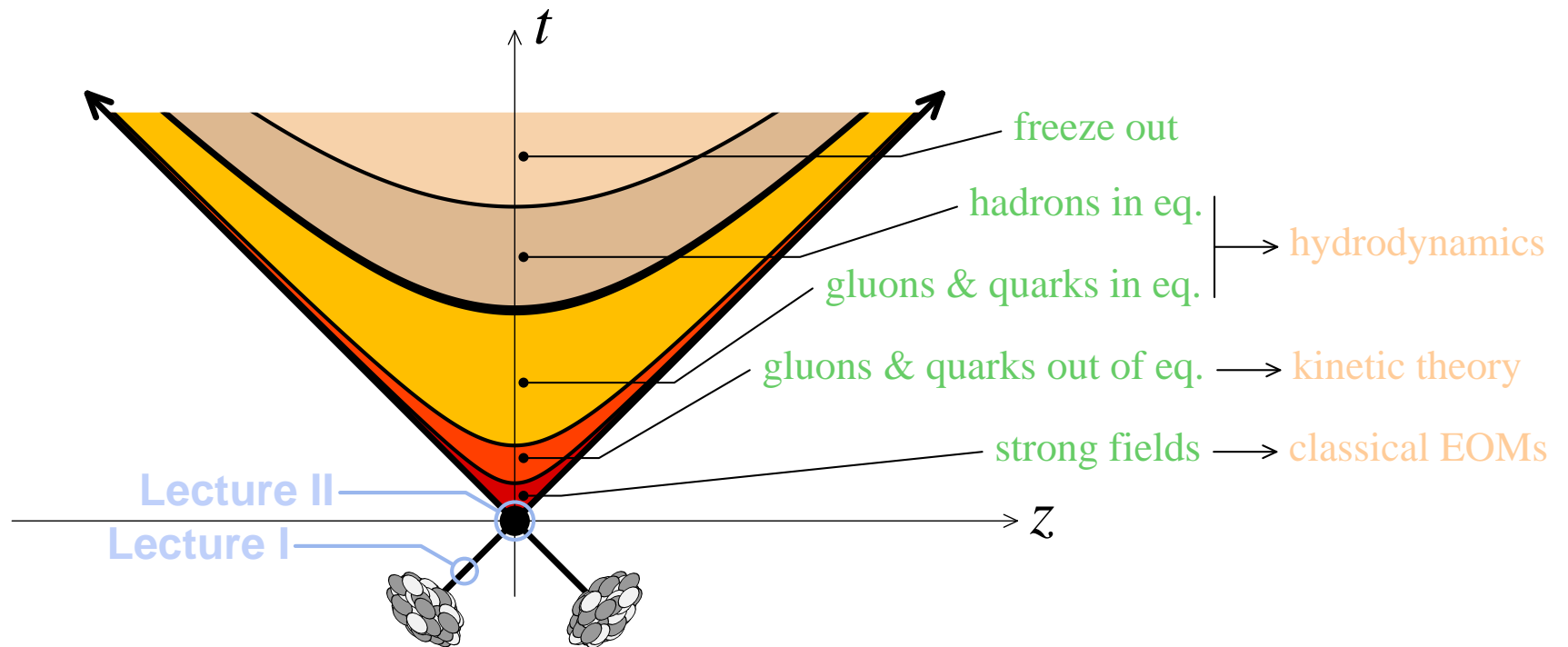
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- **Lecture I** : Parton evolution at small x , Saturation
- **Lecture II** : Initial particle production

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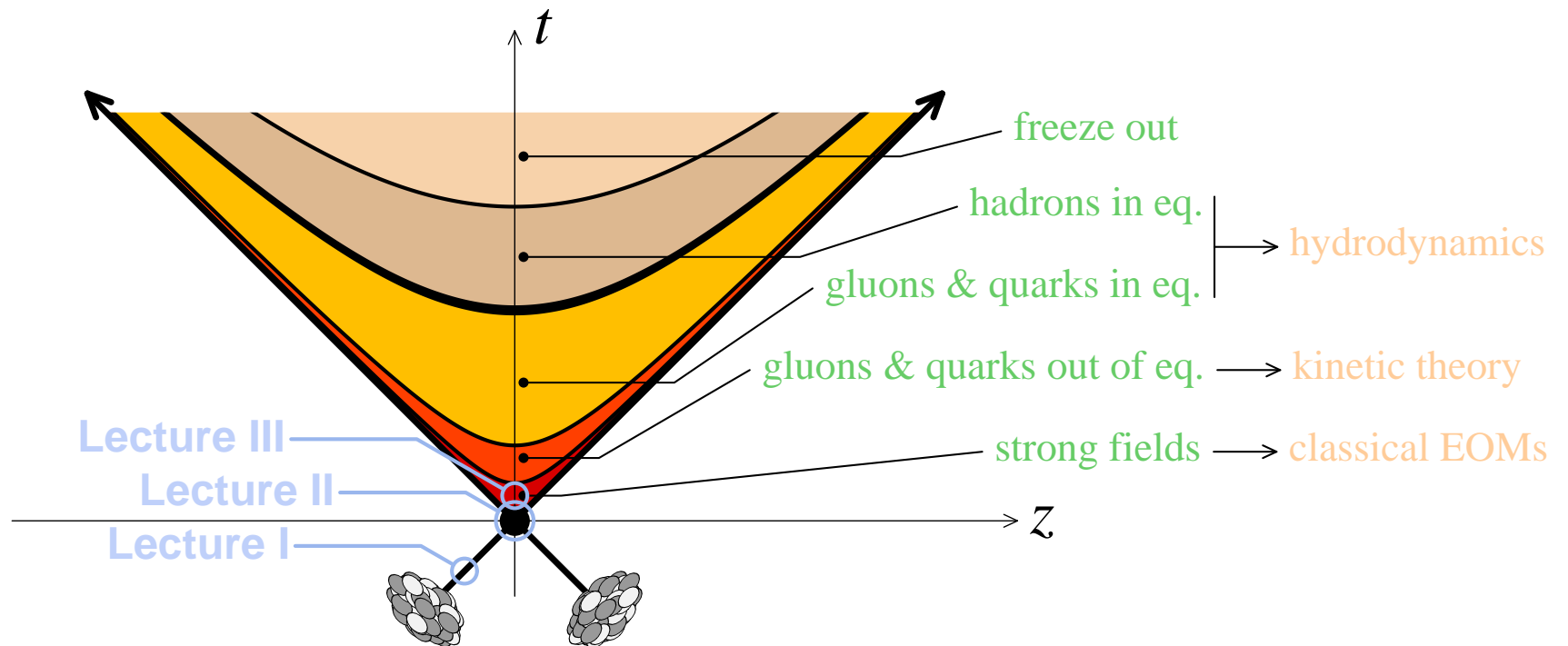
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- **Lecture I** : Parton evolution at small x , Saturation
- **Lecture II** : Initial particle production
- **Lecture III** : Instabilities and thermalization

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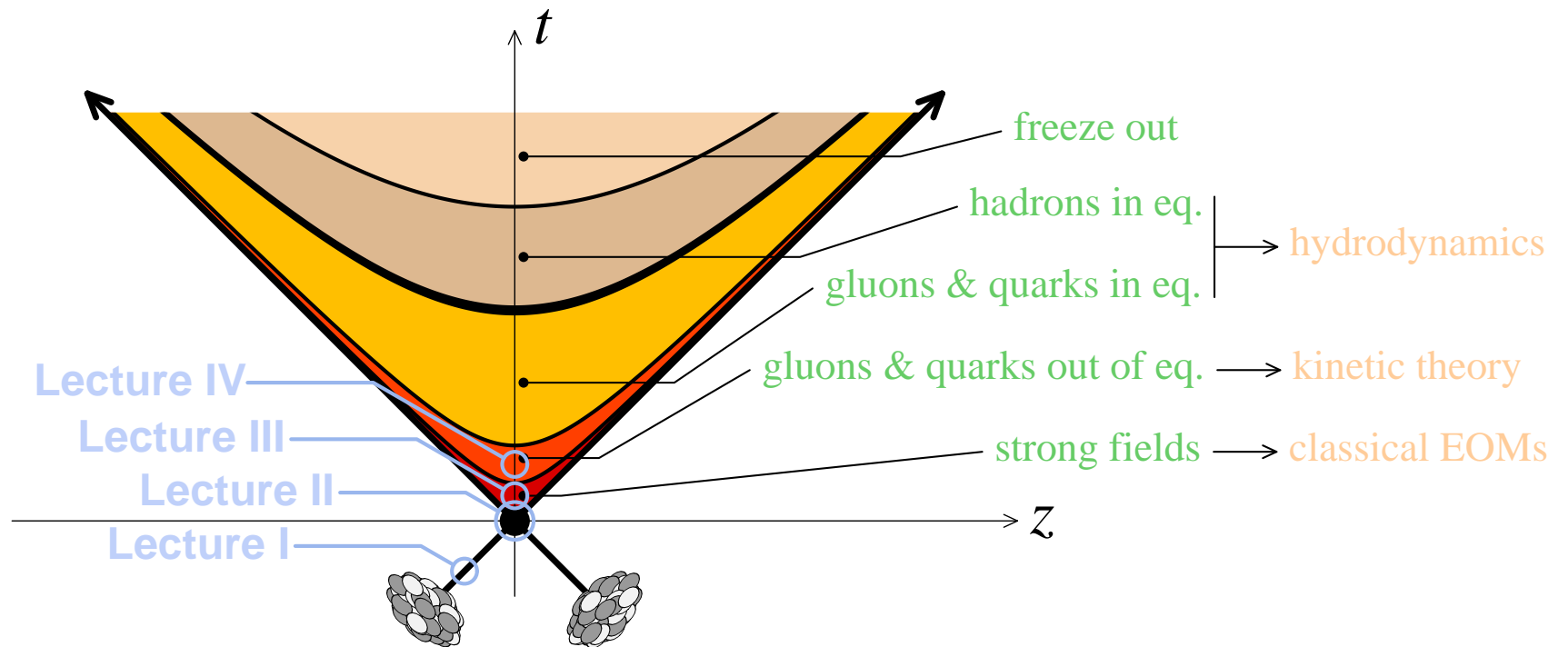
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- **Lecture I** : Parton evolution at small x , Saturation
- **Lecture II** : Initial particle production
- **Lecture III** : Instabilities and thermalization
- **Lecture IV** : Kinetic theory, Near-Equilibrium dynamics



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Lecture I : Parton saturation

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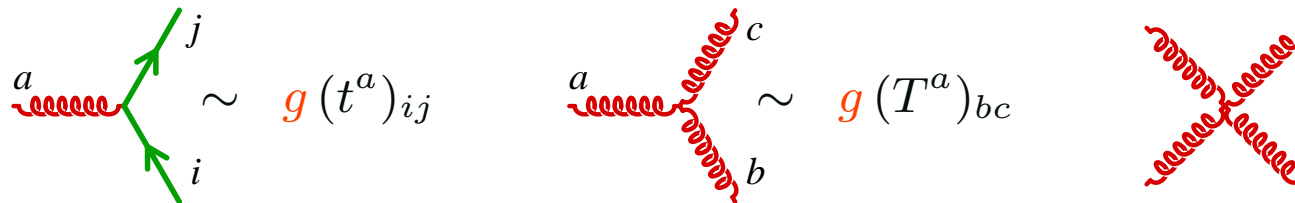
■ Electromagnetic interaction : Quantum electrodynamics

- ◆ Matter : **electron** , interaction carrier : **photon**
- ◆ Interaction :



■ Strong interaction : Quantum chromo-dynamics

- ◆ Matter : **quarks** , interaction carriers : **gluons**
- ◆ Interactions :



- ◆ i, j : colors of the quarks (3 possible values)
- ◆ a, b, c : colors of the gluons (8 possible values)
- ◆ $(t^a)_{ij}$: 3×3 matrix , $(T^a)_{bc}$: 8×8 matrix



QCD Lagrangian

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■ QCD Lagrangian :

$$\mathcal{L} = -\frac{1}{2} \text{tr} (F_{\mu\nu} F^{\mu\nu}) + \bar{\psi} (i\not{D} - m) \psi$$

- ◆ the gauge field A^μ belongs to $SU(3)$
- ◆ $D^\mu \equiv \partial^\mu - igA^\mu$ is the covariant derivative
- ◆ $F^{\mu\nu} \equiv i[D^\mu, D^\nu]/g = \partial^\mu A^\nu - \partial^\nu A^\mu - ig[A^\mu, A^\nu]$

■ The Lagrangian is invariant under gauge transformations :

$$A^\mu(x) \rightarrow \Omega(x) A^\mu(x) \Omega^{-1}(x) + \frac{i}{g} \Omega(x) \partial^\mu \Omega^{-1}(x)$$
$$\psi(x) \rightarrow \Omega(x) \psi(x)$$

where $\Omega(x) \in SU(3)$

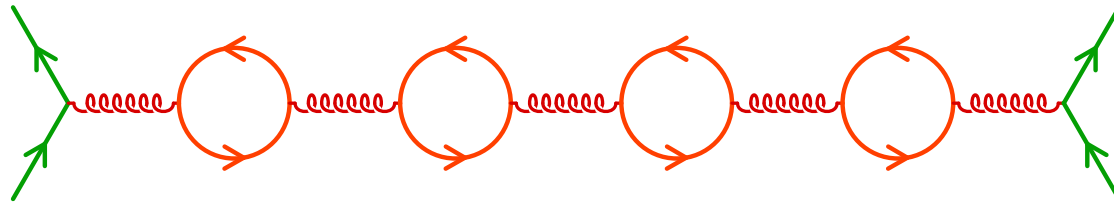
- ◆ Note: the field strength is not invariant but transforms as :

$$F^{\mu\nu}(x) \rightarrow \Omega(x) F^{\mu\nu}(x) \Omega^{-1}(x)$$

Asymptotic freedom

- Running coupling : $\alpha_s = g^2/4\pi$

$$\alpha_s(r) = \frac{2\pi N_c}{(11N_c - 2N_f) \log(1/r\Lambda_{QCD})}$$



- The effective charge seen at large distance is screened by fermionic fluctuations (as in QED)

Asymptotic freedom

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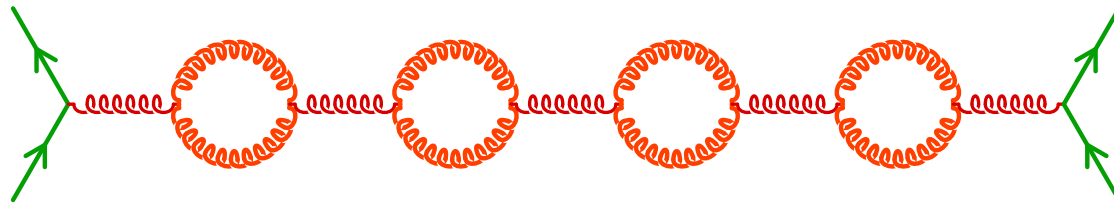
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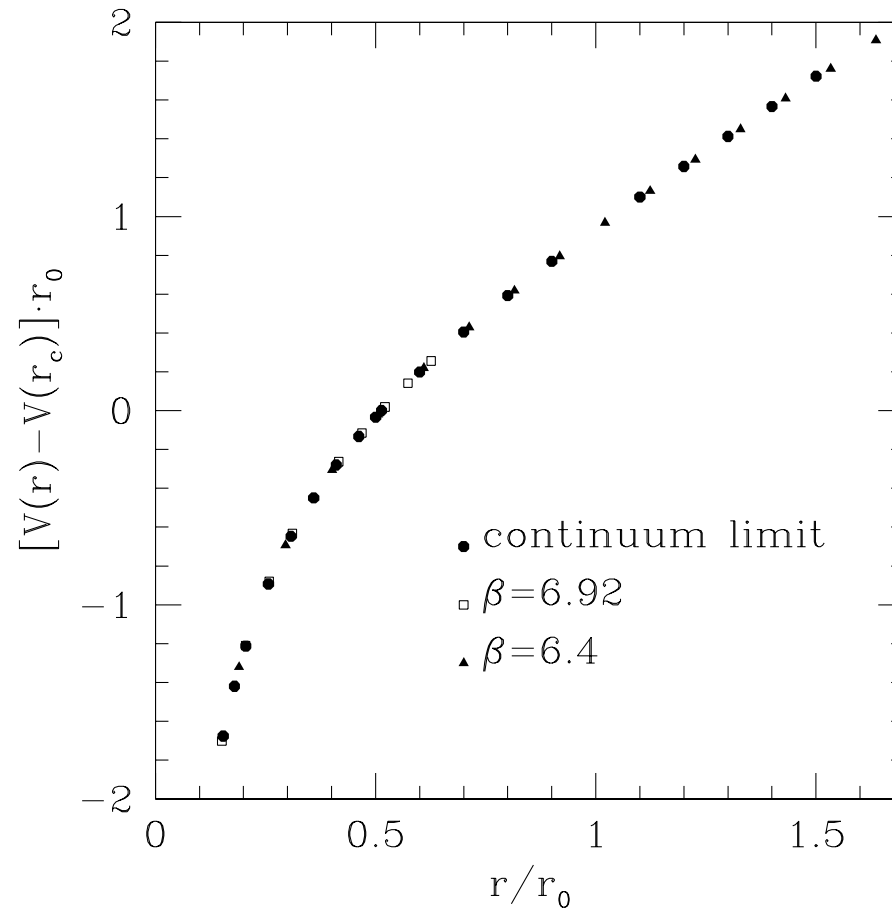
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- Running coupling : $\alpha_s = g^2/4\pi$

$$\alpha_s(r) = \frac{2\pi N_c}{(11N_c - 2N_f) \log(1/r\Lambda_{QCD})}$$



- The effective charge seen at large distance is screened by fermionic fluctuations (as in QED)
- But gluonic vacuum fluctuations produce an anti-screening (because of the non-abelian nature of their interactions)
- As long as $N_f < 11N_c/2 = 16.5$, the gluons win...



- The quark potential increases linearly with distance
- Color singlet hadrons : no free quarks and gluons in nature



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- QCD is the fundamental theory of strong interactions among quarks and gluons
- Experiments involve hadrons in their initial and final states, not quarks and gluons
- Hadrons cannot be described perturbatively in QCD
- Scattering amplitudes with time-like on-shell momenta cannot be computed on the lattice
 - ▷ How can we compare theory and experiments?
 - ▷ **Factorization** : separation of short distances (perturbative) and long distance (non perturbative)



Factorization

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- At a superficial level, factorization means that :

$$\mathcal{O}_{\text{hadrons}} = F \otimes \mathcal{O}_{\text{partons}}$$

- ◆ F = parton distribution
- ◆ $\mathcal{O}_{\text{partons}}$ = observable at the partonic level
(calculable in perturbation theory)

- For this to be useful, F must be **universal** (i.e. independent of the observable \mathcal{O})
- In order to test QCD experimentally, measure as many observables as possible, and try to find common F 's that fit all the data
Note : at this stage, by looking at only one observable, it is impossible to perform any meaningful test, since it is always possible to adjust F so that it works



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- Some loop corrections in $\mathcal{O}_{\text{partons}}$ are enhanced by large logarithms, e.g.

$$\alpha_s \ln \left(\frac{M^2}{m_H^2} \right), \quad \alpha_s \ln \left(\frac{s}{M^2} \right) \sim \alpha_s \ln \left(\frac{1}{x} \right)$$

Note : the log that occurs depends on the details of the kinematics

- ◆ Bjorken limit: $s, M^2 \rightarrow +\infty$ with s/M^2 fixed
 - ◆ Regge limit: $s \rightarrow +\infty, M^2$ fixed
- These logs upset a naive application of perturbation theory when $\alpha_s \ln(\cdot) \sim 1 \triangleright$ they must be resummed
 - This resummation can be performed analytically
 - ◆ the result of the resummation is universal
 - ◆ all the leading logs can be absorbed in F
 - \triangleright the factorization formula remains true
 - \triangleright this summation dictates how F evolves with M^2 or x

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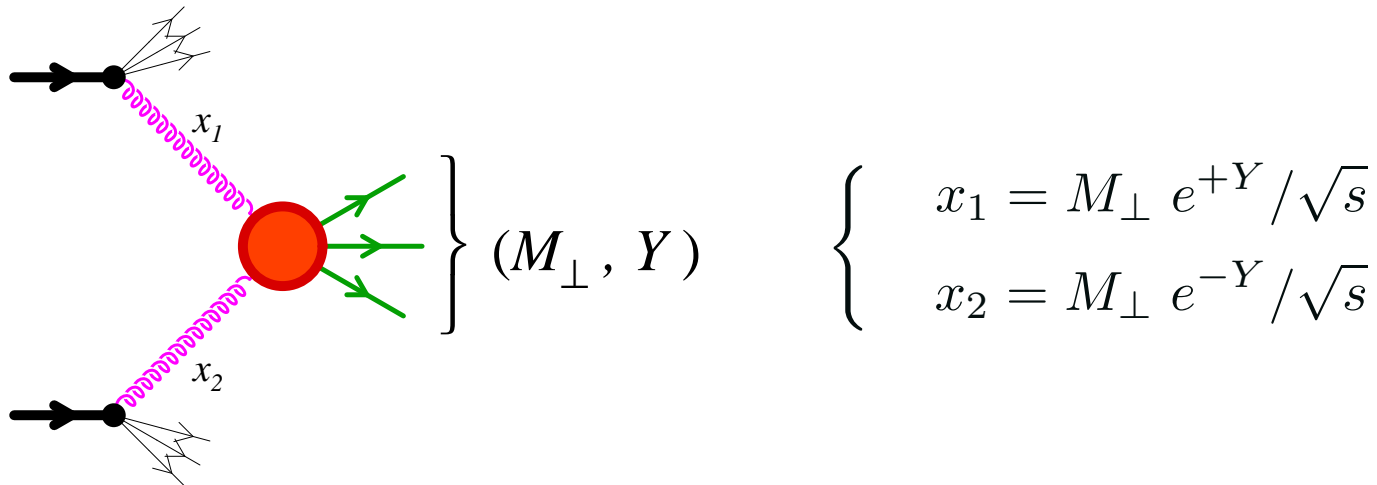
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- These logarithms tell us that the relevant parton distributions depend on the resolution scales (in time and in transverse momentum) associated to a given process

- Calculation of some process at LO :



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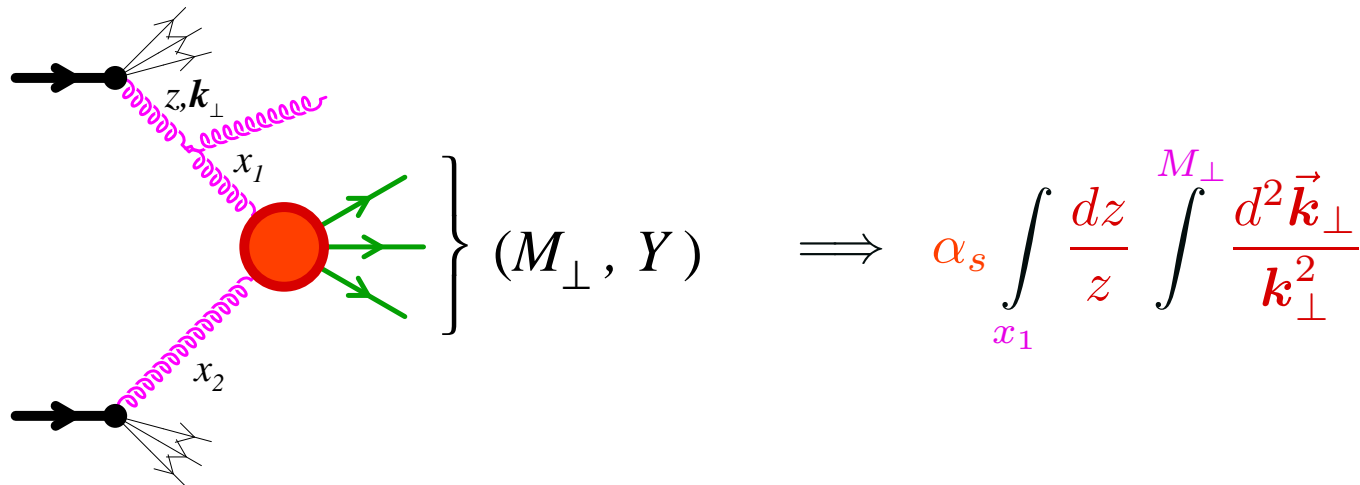
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Phenomenology of saturation

- These logarithms tell us that the relevant parton distributions depend on the resolution scales (in time and in transverse momentum) associated to a given process

- Radiation of an extra gluon :



- Practical consequence : pQCD predicts not only $\mathcal{O}_{\text{partons}}$ but also the evolution $\partial_M F$ (or $\partial_x F$)
 - ▷ the only required non-perturbative input is $F(x, M_0)$ or $F(x_0, M)$



Collinear factorization

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- Logs of $M_{\perp} \implies$ **DGLAP**. Important when :
 - ◆ $M_{\perp} \gg \Lambda_{QCD}$, while x_1, x_2 are rather large

- Cross-sections read :

$$\frac{d\sigma}{dY d^2\vec{P}_{\perp}} \propto F(x_1, M_{\perp}^2) F(x_2, M_{\perp}^2) |\mathcal{M}|^2$$

with $x_{1,2} = M_{\perp} \exp(\pm Y) / \sqrt{s}$

- Note : there are convolutions in x_1 and x_2 if some particles are integrated out in the final state
- The factorization of logarithms has been proven to all orders for sufficiently inclusive quantities (see **Collins, Soper, Sterman, 1984–1985**)



Kt-factorization

Collins, Ellis (1991), Catani, Ciafaloni, Hautmann (1991)

- Logs of $1/x \implies$ **BFKL**. Important when :
 - ◆ M_{\perp} remains moderate, while x_1 or x_2 (or both) are small
- The BFKL equation is non-local in transverse momentum
 - ▷ it applies to **non-integrated gluon distributions** $\varphi(x, \vec{k}_{\perp})$

$$xG(x, Q^2) = \int \frac{d^2 \vec{k}_{\perp}}{(2\pi)^2} \varphi(x, \vec{k}_{\perp})$$

▷ the matrix element is calculated for (off-shell) gluons with $\vec{k}_{\perp} \neq \vec{0}$

- In this framework, cross-sections read :

$$\frac{d\sigma}{dY d^2 \vec{P}_{\perp}} \propto \int_{\vec{k}_{1\perp}, \vec{k}_{2\perp}} \delta(\vec{k}_{1\perp} + \vec{k}_{2\perp} - \vec{P}_{\perp}) \varphi_1(x_1, k_{1\perp}) \varphi_2(x_2, k_{2\perp}) \frac{|\mathcal{M}|^2}{k_{1\perp}^2 k_{2\perp}^2}$$

$$(x_{1,2} = M_{\perp} e^{\pm Y} / \sqrt{s})$$

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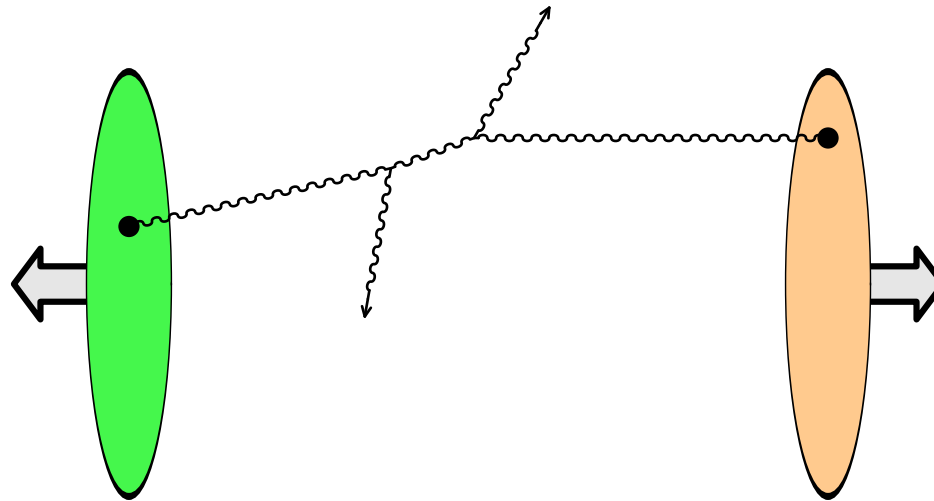
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- **Collinear or kt -factorization** : only one parton in each projectile take part in the process of interest

Multi-parton interactions?

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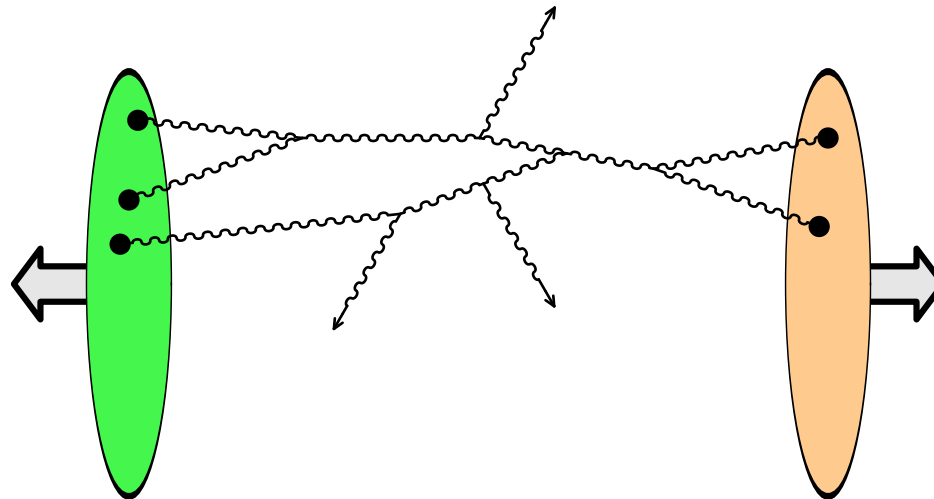
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- **Collinear or kt -factorization** : only one parton in each projectile take part in the process of interest
- **If multiparton interactions are important** : the above forms of factorization cannot work anymore, because the only information they retain about the distribution of partons is their 2-point correlations (i.e. the number of partons)



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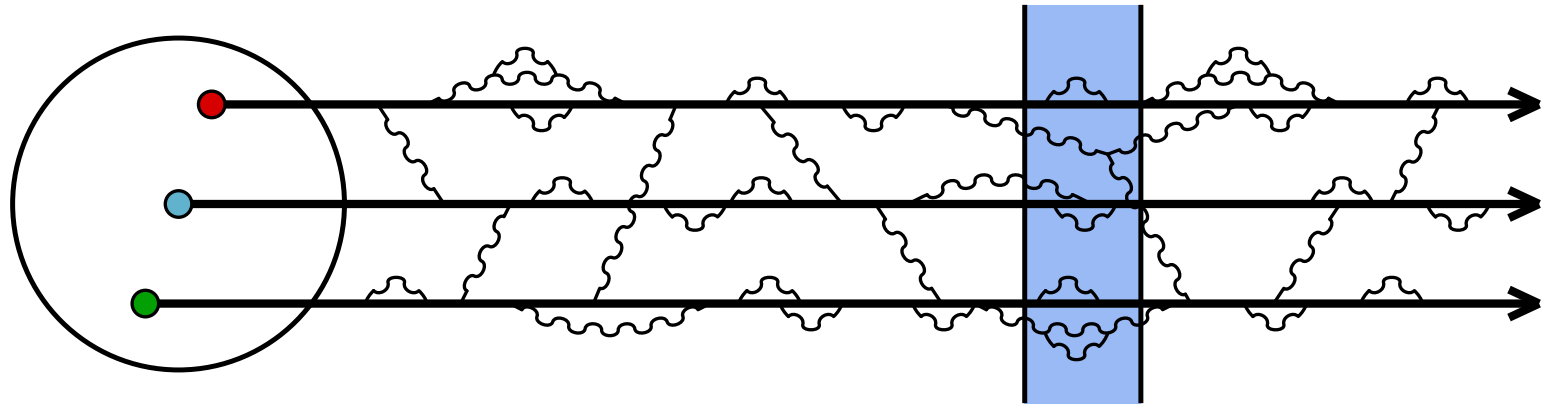
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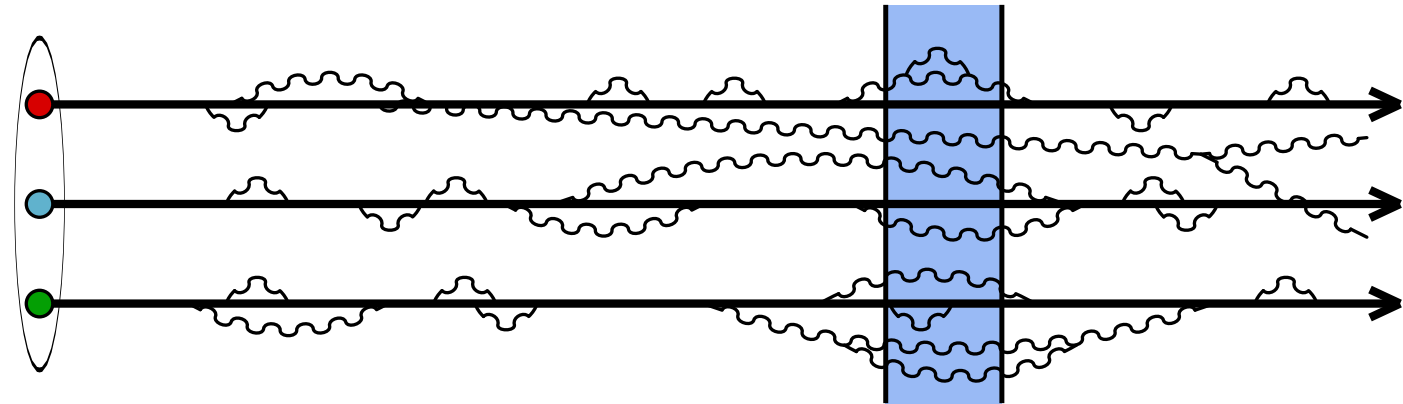
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Nucleon at low energy



- A **nucleon at rest** is a very complicated object...
- Contains **fluctuations at all space-time scales** smaller than its own size
- Only the fluctuations that are longer lived than the external probe participate in the interaction process
- The only role of short lived fluctuations is to renormalize the masses and couplings
- Interactions are very complicated if the constituents of the nucleon have a non trivial dynamics over time-scales comparable to those of the probe

Nucleon at high energy



- Dilation of all internal time-scales for a **high energy nucleon**
- Interactions among constituents now take place over time-scales that are longer than the characteristic time-scale of the probe
 - ▷ **the constituents behave as if they were free**
- Many fluctuations live long enough to be seen by the probe. The nucleon appears **denser at high energy** (it contains more gluons)



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- At the time of the interaction, the nucleon can be seen as a collection of **free constituents**, called **partons**
- It can be described by **non-perturbative parton distributions** that depend on the momentum fraction x of the partons
- One can separate the perturbative hard scattering from the non-perturbative distribution functions, because the strong interactions that are responsible for these non-perturbative aspects occur on much larger timescales (**factorization**)
- All these properties are based only on kinematics and causality, and should remain true in the saturation regime
 - ◆ what we use as the “parton distribution” must contain information about multiparton configurations
 - ◆ the calculation of the “hard” process is more involved



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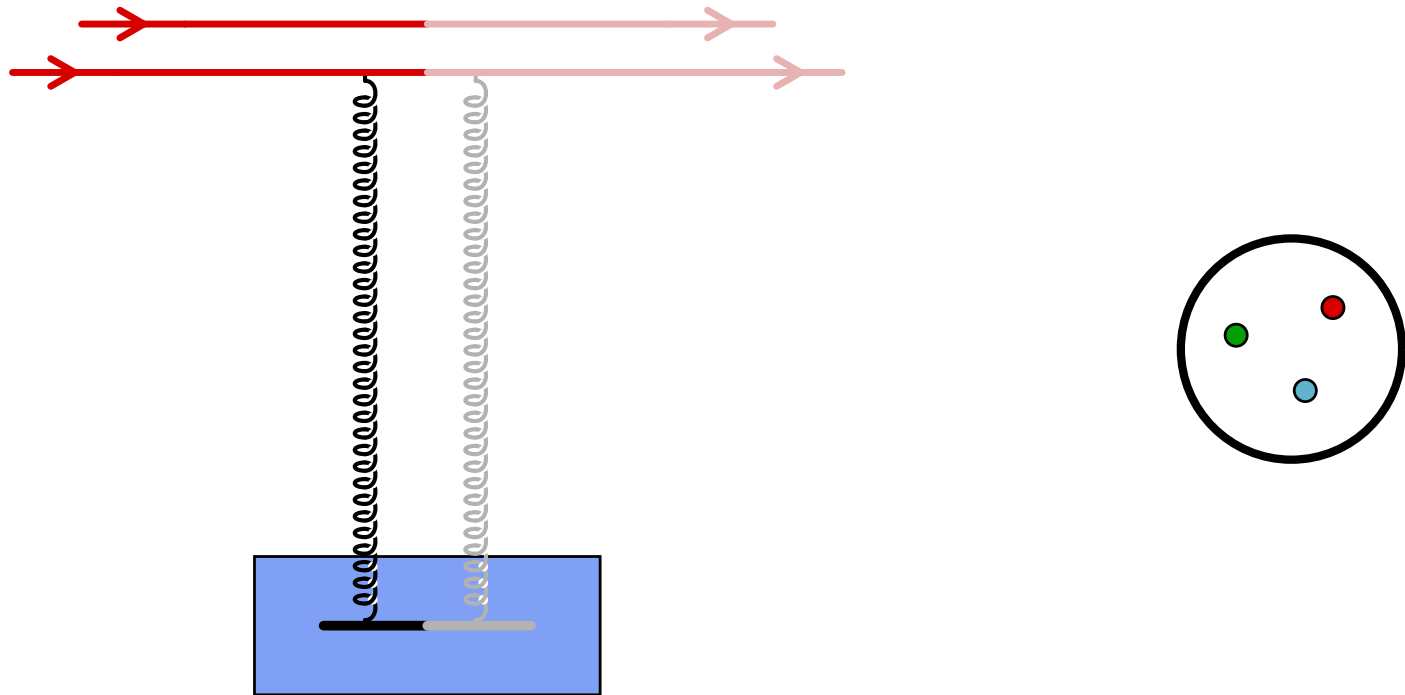
● Saturation criterion

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Phenomenology of saturation



- ▷ assume that the projectile is big, e.g. a nucleus, and has many valence quarks (only two are represented)
- ▷ on the contrary, consider a small probe, with few partons
- ▷ at low energy, only valence quarks are present in the hadron wave function

Parton evolution

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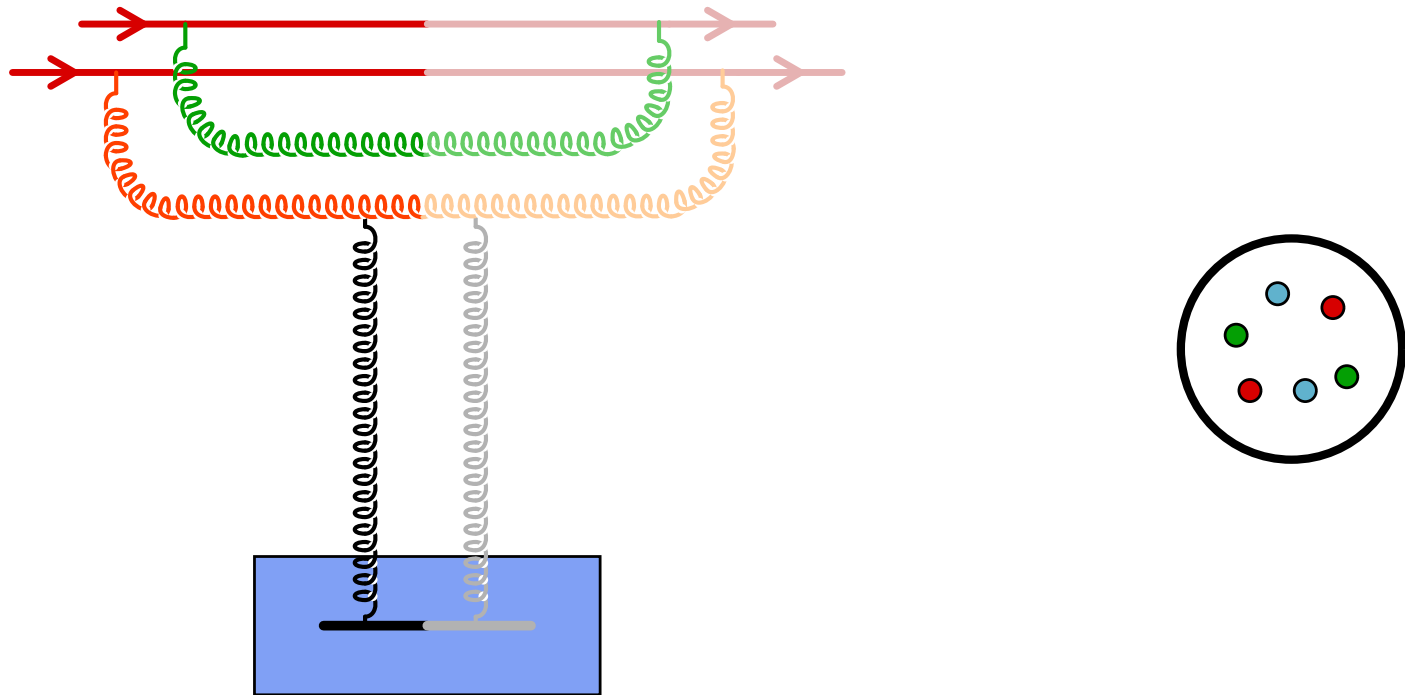
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- ▷ when energy increases, new partons are emitted
- ▷ the emission probability is $\alpha_s \int \frac{dx}{x} \sim \alpha_s \ln\left(\frac{1}{x}\right)$, with x the longitudinal momentum fraction of the gluon
- ▷ at small- x (i.e. high energy), these logs need to be resummed

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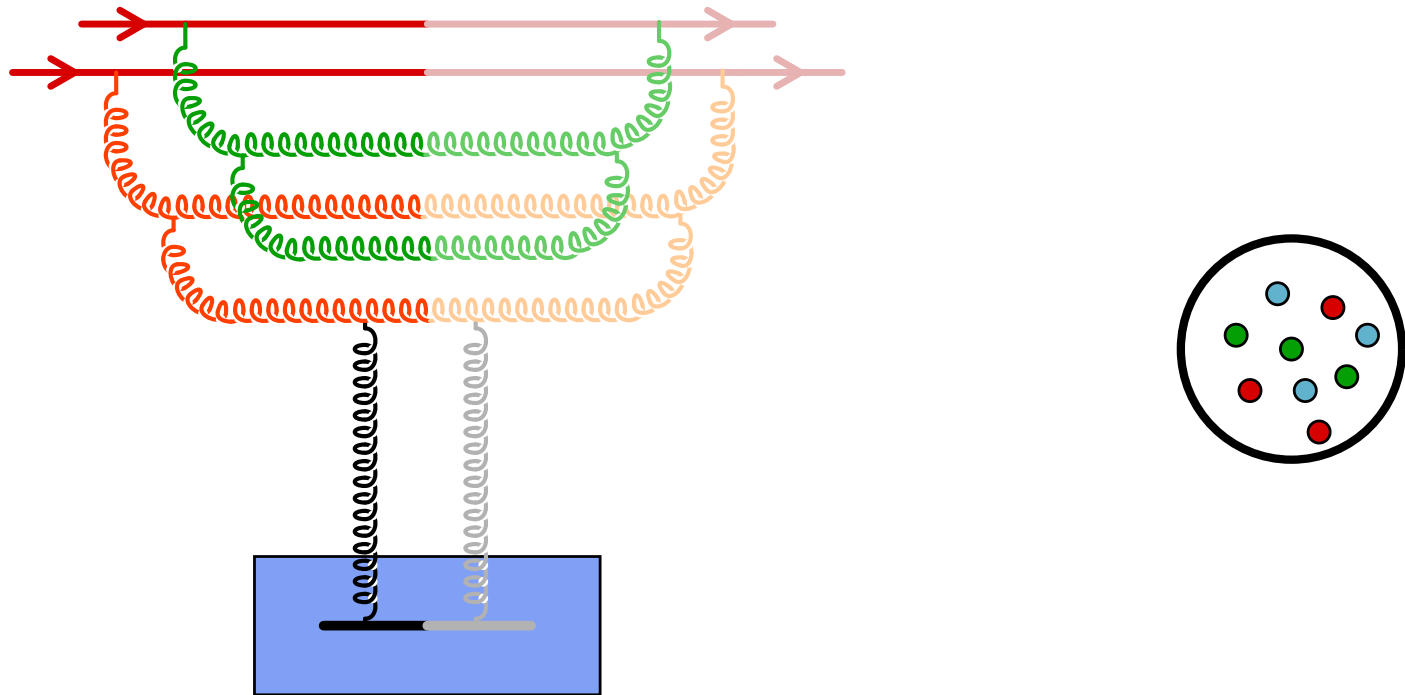
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▷ as long as the density of constituents remains small, the evolution is **linear**: the number of partons produced at a given step is proportional to the number of partons at the previous step (BFKL)

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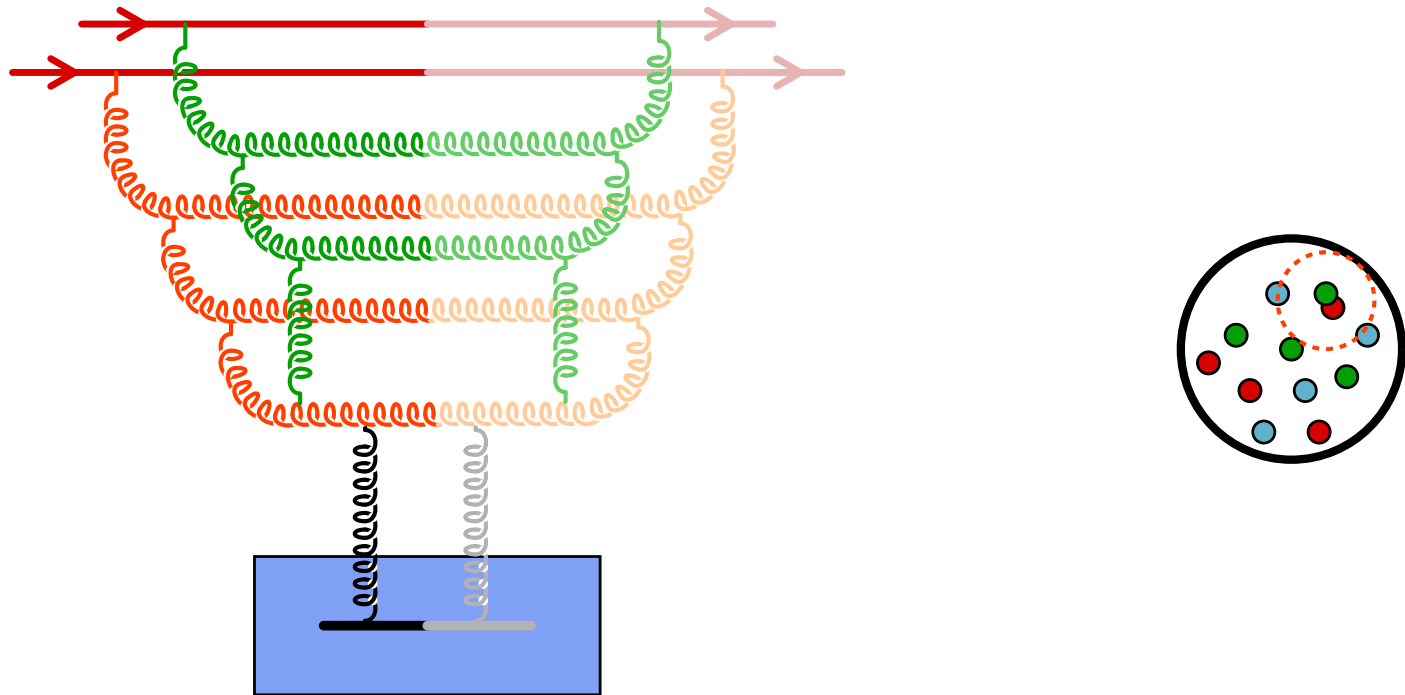
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Phenomenology of saturation



- ▷ eventually, the partons start overlapping in phase-space
- ▷ **parton recombination** becomes favorable
- ▷ after this point, the evolution is **non-linear**:
the number of partons created at a given step depends non-linearly on the number of partons present previously



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Phenomenology of saturation

Gribov, Levin, Ryskin (1983)

- Number of gluons per unit area:

$$\rho \sim \frac{xG_A(x, Q^2)}{\pi R_A^2}$$

- Recombination cross-section:

$$\sigma_{gg \rightarrow g} \sim \frac{\alpha_s}{Q^2}$$

- Recombination happens if $\rho \sigma_{gg \rightarrow g} \gtrsim 1$, i.e. $Q^2 \lesssim Q_s^2$, with:

$$Q_s^2 \sim \frac{\alpha_s xG(x, Q_s^2)}{\pi R_A^2} \sim A^{1/3} \frac{1}{x^{0.3}}$$

- At saturation, the phase-space density is:

$$\frac{dN_g}{d^2 \vec{x}_\perp d^2 \vec{p}_\perp} \sim \frac{\rho}{Q^2} \sim \frac{1}{\alpha_s}$$

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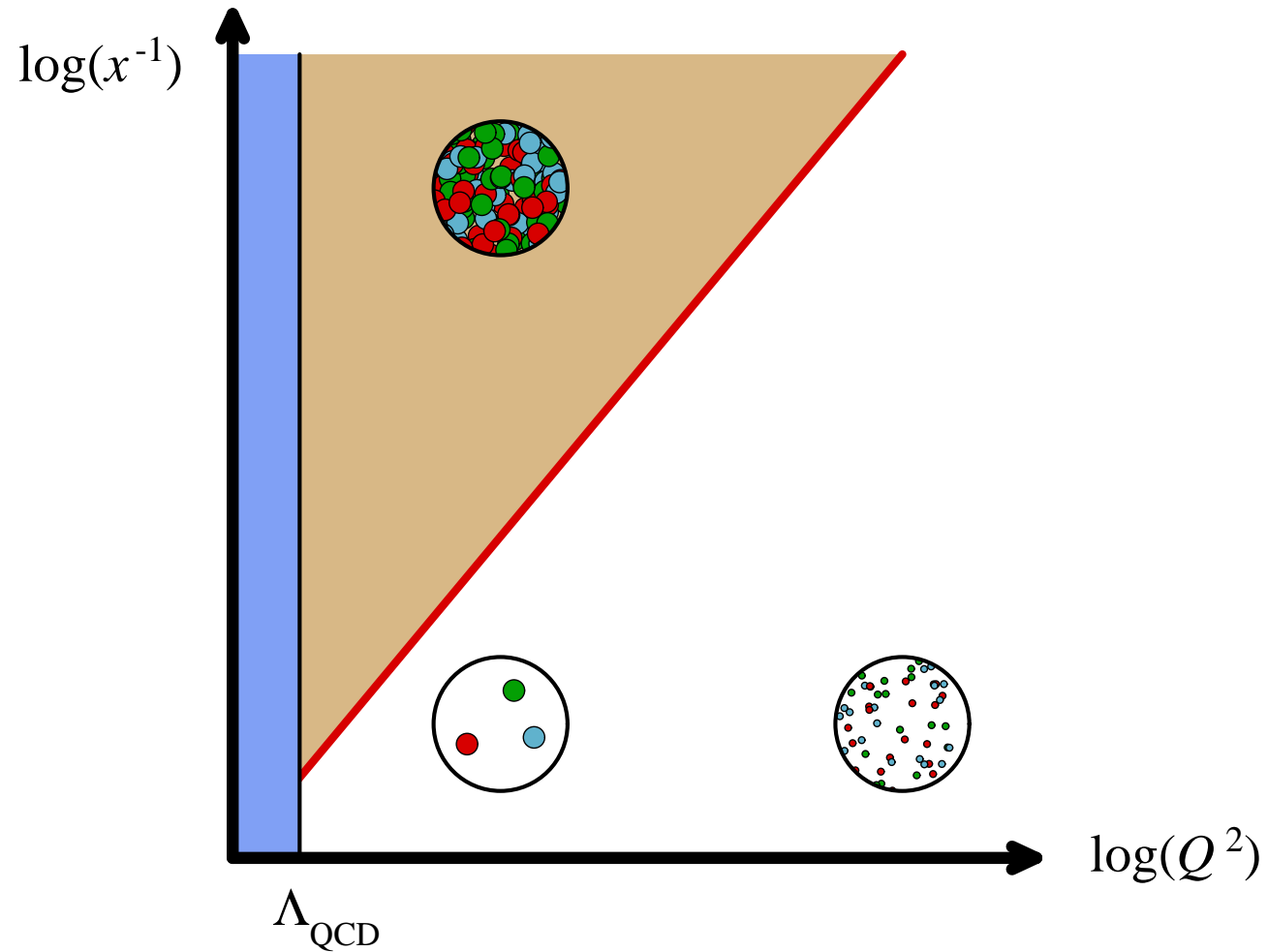
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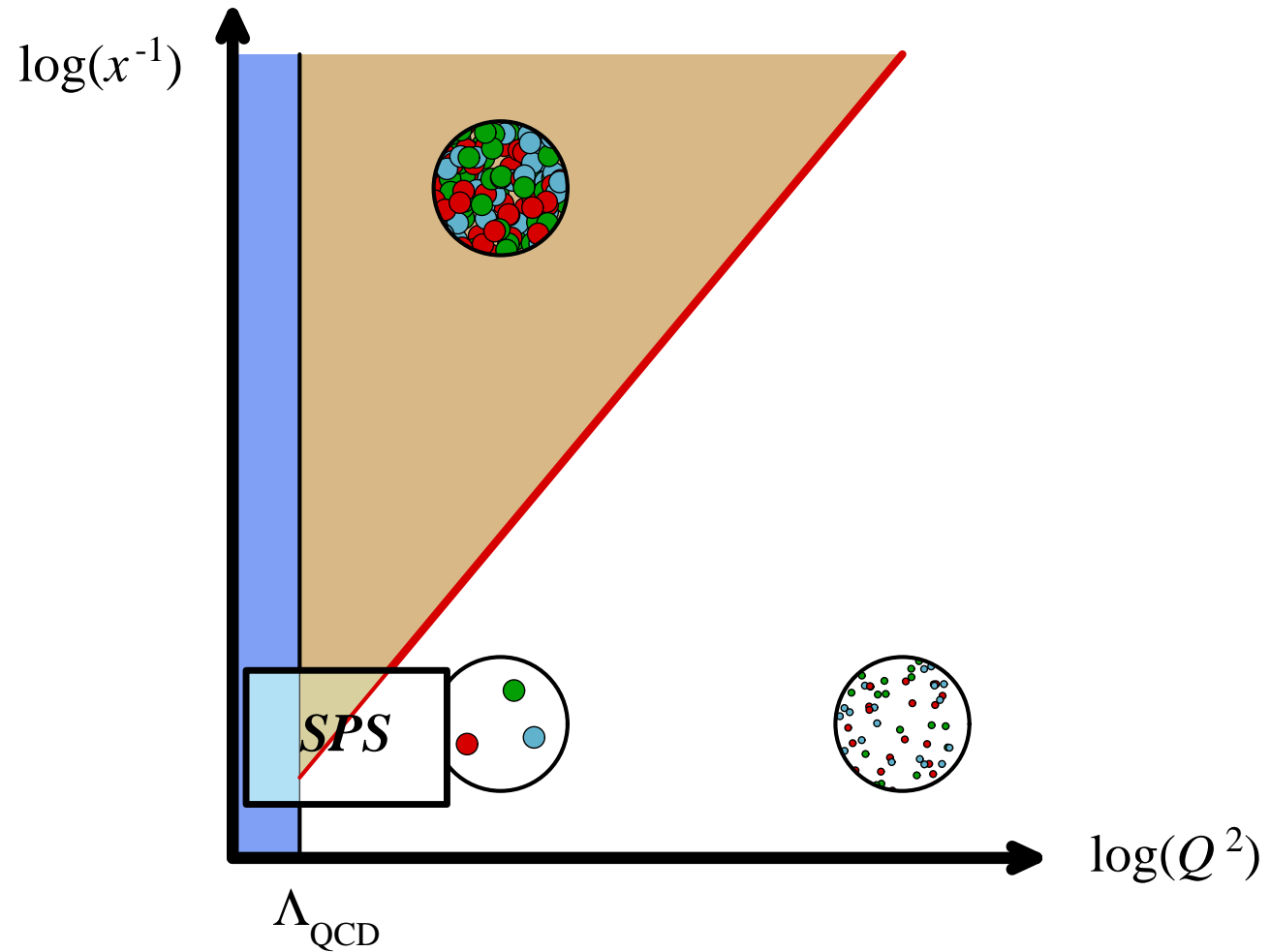
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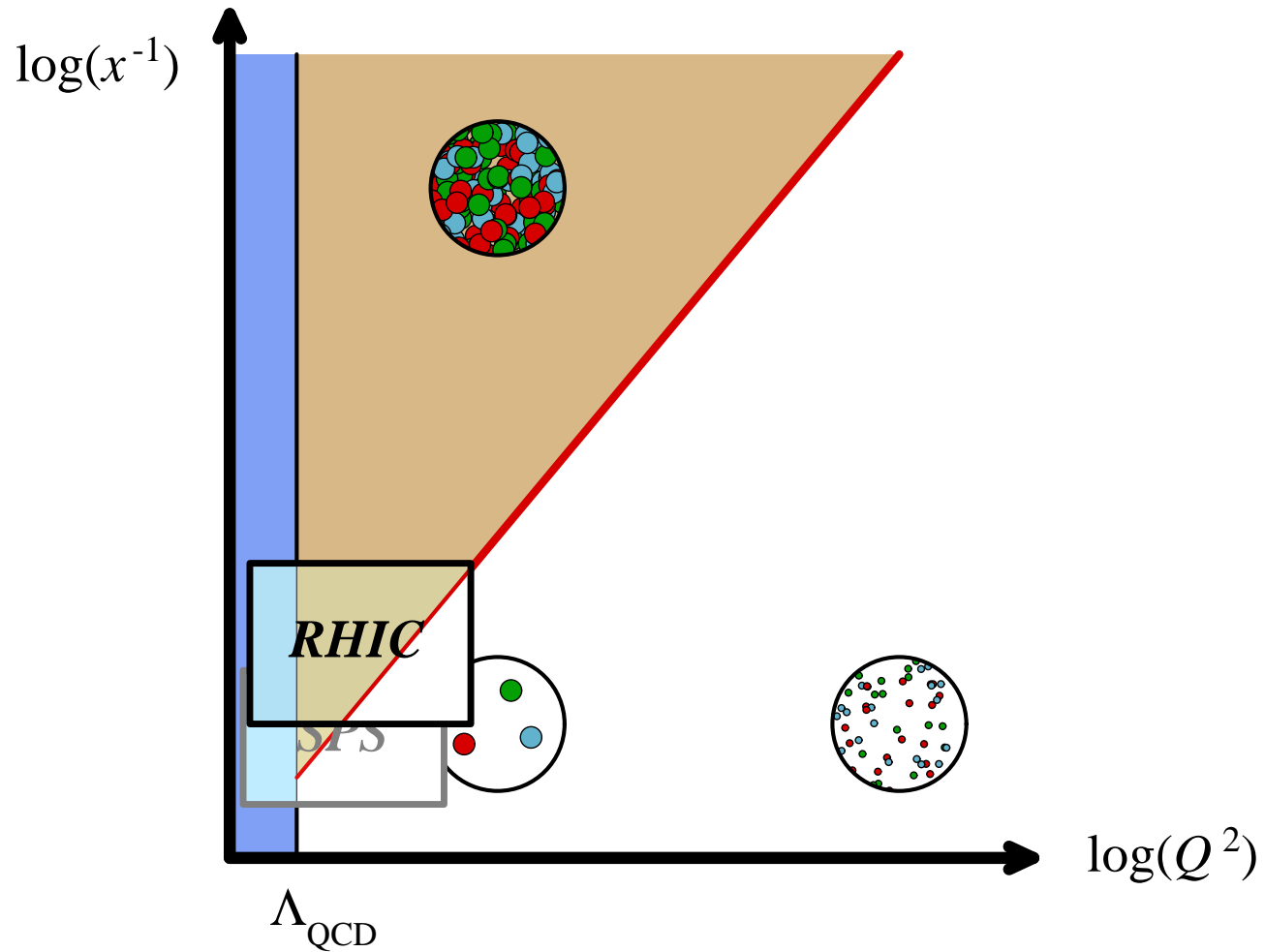
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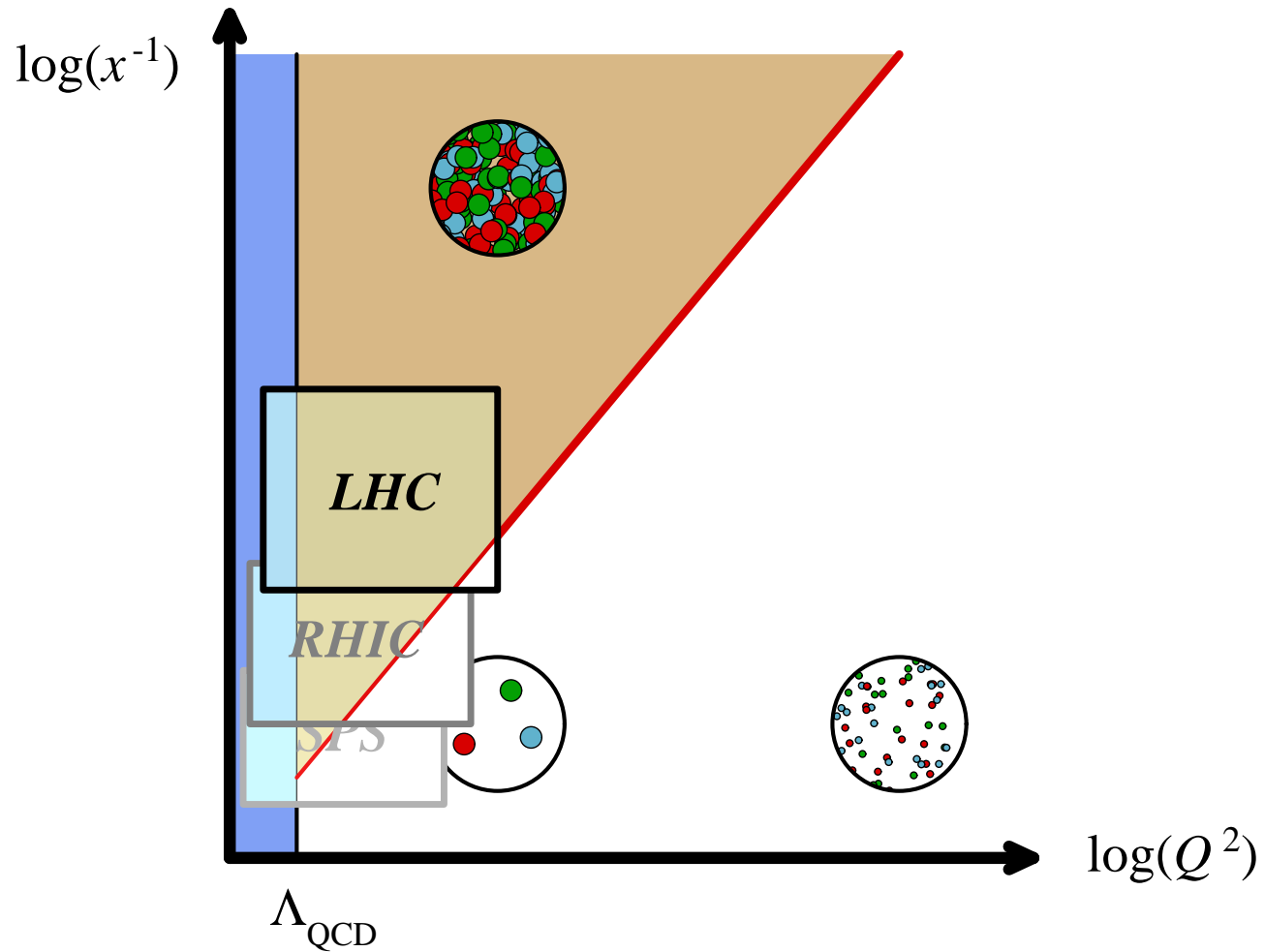
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- Saturation criterion
- **Saturation domain**
- Multiple scatterings

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Phenomenology of saturation



Saturation domain

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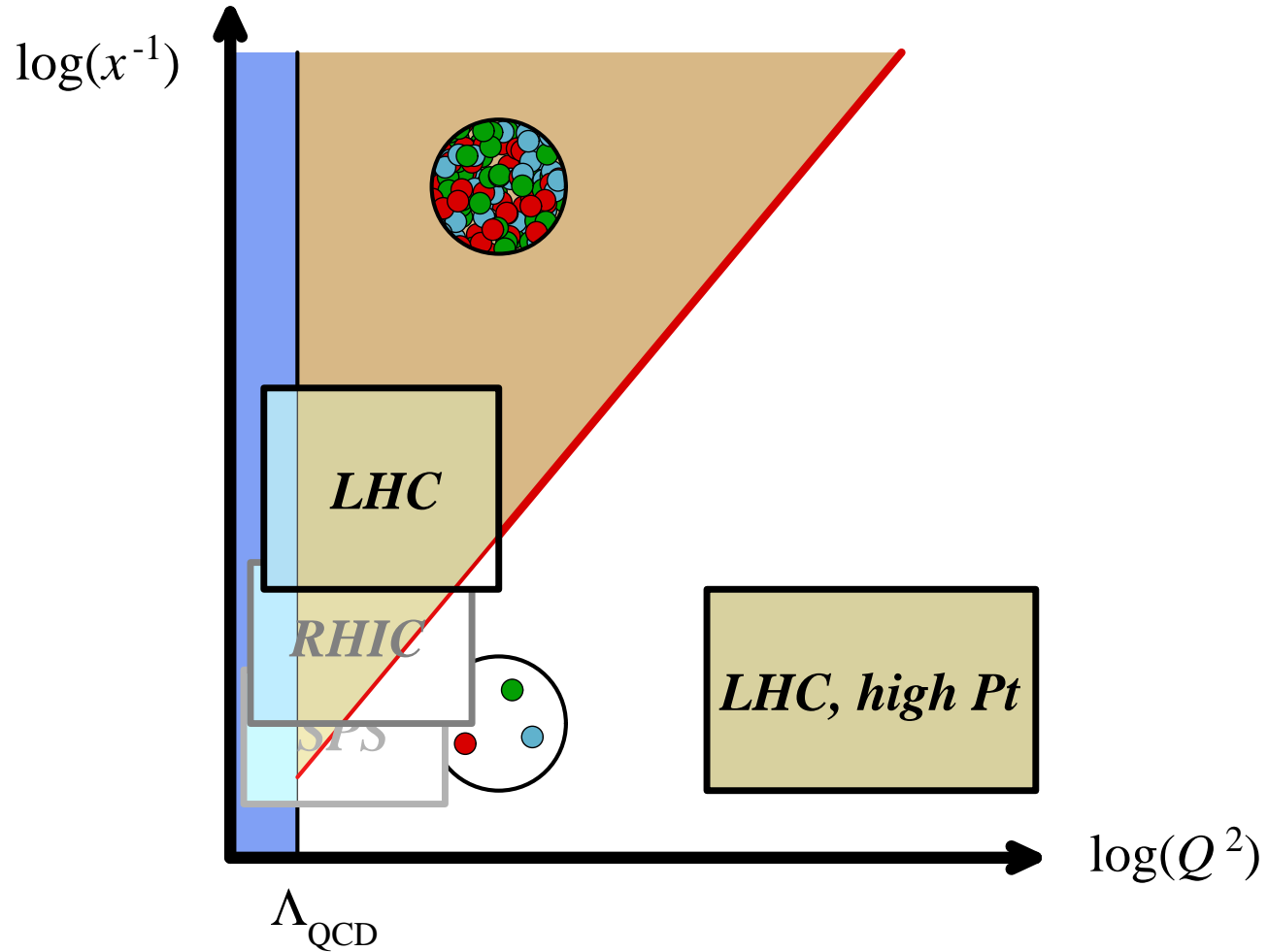
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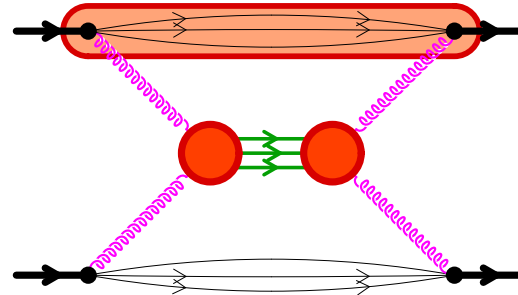
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Multiple scatterings

■ Single scattering :



▷ 2-point function in the projectile ▷ gluon number

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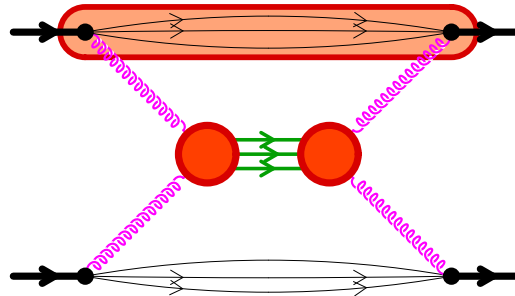
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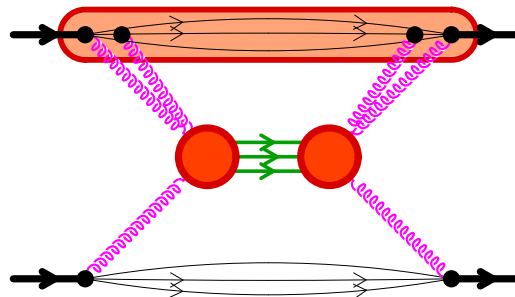
Phenomenology of saturation

■ Single scattering :



▷ 2-point function in the projectile ▷ gluon number

■ Multiple scatterings :



▷ 4-point function in the projectile ▷ higher correlation
▷ multiple scatterings in the projectile



Multiple scatterings

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- **Power counting** : rescattering corrections are suppressed by inverse powers of the typical mass scale in the process :

$$\left[\frac{\mu^2}{M_\perp^2} \right]^n$$

- The parameter μ^2 has a factor of α_s , and a factor proportional to the gluon density \triangleright **rescatterings are important at high density**
- Relative order of magnitude :

$$\frac{2 \text{ scatterings}}{1 \text{ scattering}} \sim \frac{Q_s^2}{M_\perp^2} \quad \text{with} \quad Q_s^2 \sim \alpha_s \frac{xG(x, Q_s^2)}{\pi R^2}$$

- When this ratio becomes ~ 1 , all the rescattering corrections become important \triangleright **one must resum all $[Q_s/M_\perp]^n$**
- These effects are not accounted for in DGLAP or BFKL



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- The fast partons (large x) are frozen by time dilation
▷ described as **static color sources** on the light-cone :

$$J_a^\mu = \delta^{\mu+} \delta(x^-) \rho_a(\vec{x}_\perp) \quad (x^- \equiv (t - z)/\sqrt{2})$$

- Slow partons (small x) are radiated by the fast ones. They have a large occupation number ▷ described by a **classical color field** A^μ that obeys Yang-Mills's equation:

$$[D_\nu, F^{\nu\mu}]_a = J_a^\mu$$

- The color sources ρ_a are **random**, and described by a **distribution functional** $W_Y[\rho]$, with Y the rapidity that separates “soft” and “hard”

Deep Inelastic Scattering

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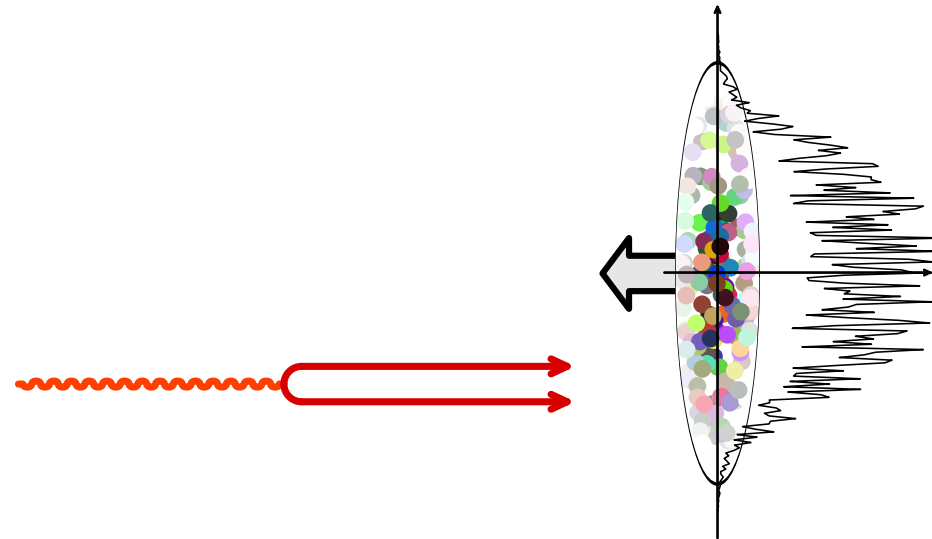
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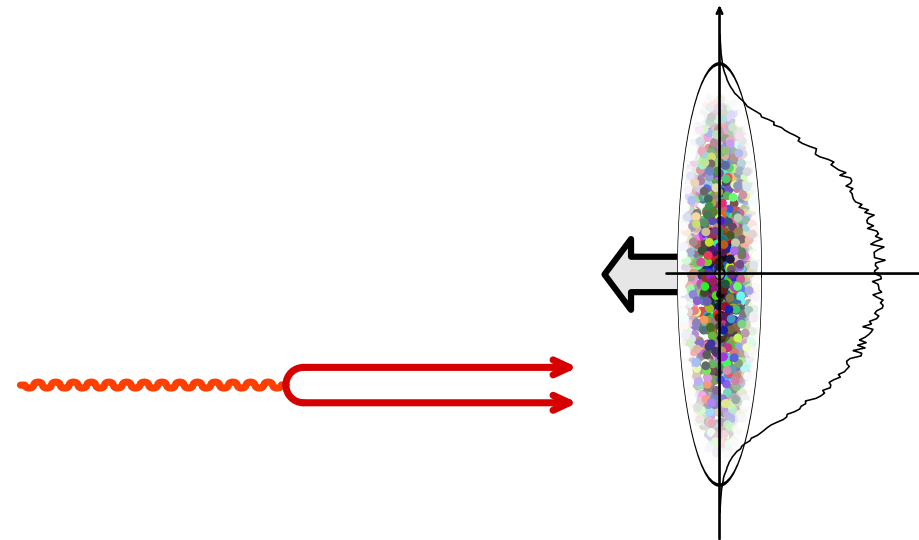
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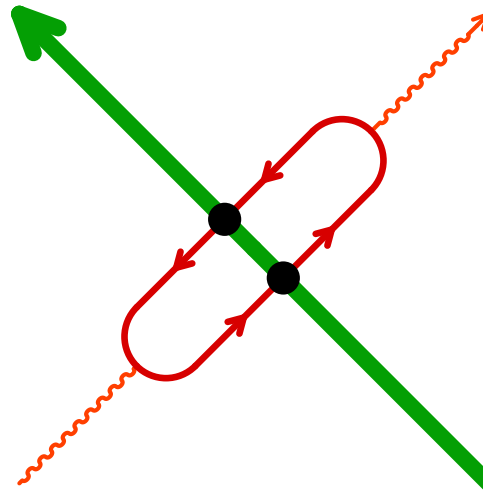
Phenomenology of saturation



100 configurations

Deep Inelastic Scattering

- Reactions involving a hadron or nucleus and an “elementary” projectile are fairly straightforward to study
- Example : **forward DIS amplitude** :



$$\langle \mathbf{T}(\vec{x}_\perp, \vec{y}_\perp) \rangle = \int [D\rho] W_Y[\rho] \left[1 - \frac{1}{N_c} \text{tr}(U(\vec{x}_\perp)U^\dagger(\vec{y}_\perp)) \right]$$

▷ this formula resums all the $[\alpha_s \ln(1/x)]^m [Q_s/p_\perp]^n$ for the inclusive DIS cross-section



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- $U(\vec{x}_\perp)$ is a **Wilson line** that represents the scattering at high energy between a quark and the color field of the nucleus (moving in the $-z$ direction) :

$$U(\vec{x}_\perp) \equiv P_+ \exp ig \int_{-\infty}^{+\infty} dz^+ A^-(z^+, \vec{x}_\perp)$$

with

$$-\vec{\nabla}_\perp^2 A^-(x^+, \vec{x}_\perp) = \delta(x^+) \rho(\vec{x}_\perp)$$

- The scattering of the antiquark is represented by $U^\dagger(\vec{y}_\perp)$

Balitsky-Kovchegov equation

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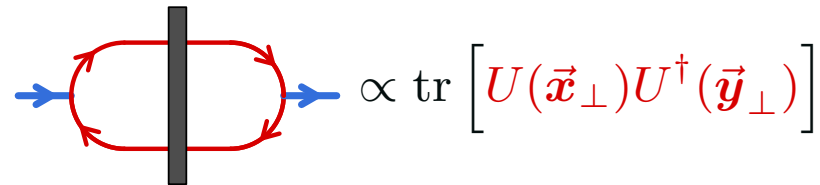
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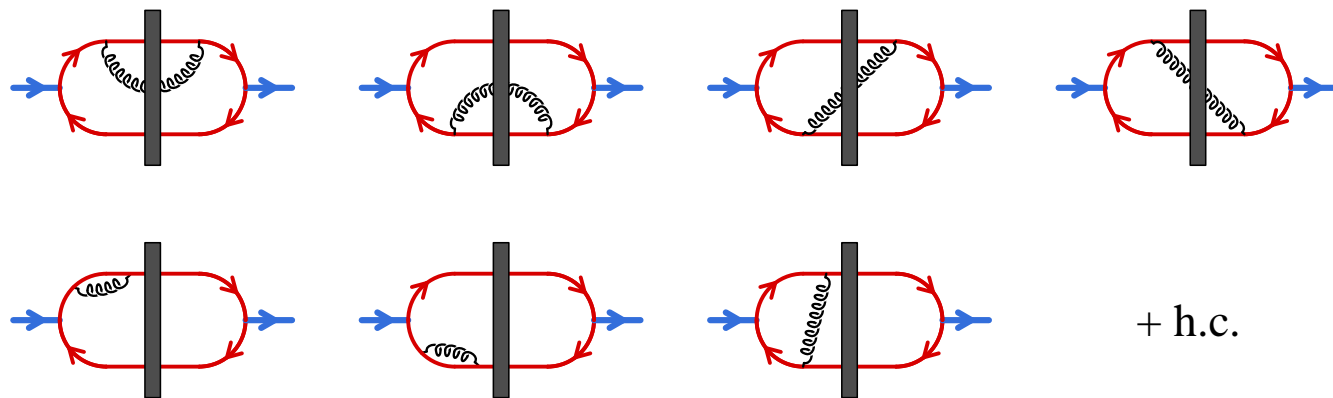
- The bare scattering amplitude of a **color singlet** $Q\bar{Q}$ dipole can be written as :



$$\propto \text{tr} [U(\vec{x}_\perp)U^\dagger(\vec{y}_\perp)]$$

Note : this bare dipole amplitude is independent of energy. The energy dependence comes from higher-order corrections

- At one loop, the following diagrams must be evaluated :



Balitsky-Kovchegov equation

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- The NLO corrections lead to an equation that drives the dependence of $T(\vec{x}_\perp, \vec{y}_\perp) \equiv 1 - \frac{1}{N_c} \text{tr} (U(\vec{x}_\perp)U^\dagger(\vec{y}_\perp))$ with respect to $Y \sim \ln(\sqrt{s})$:

$$\frac{\partial T(\vec{x}_\perp, \vec{y}_\perp)}{\partial Y} = \frac{\alpha_s N_c}{2\pi^2} \int d^2 \vec{z}_\perp \frac{(\vec{x}_\perp - \vec{y}_\perp)^2}{(\vec{x}_\perp - \vec{z}_\perp)^2 (\vec{y}_\perp - \vec{z}_\perp)^2} \times \left\{ T(\vec{x}_\perp, \vec{z}_\perp) + T(\vec{z}_\perp, \vec{y}_\perp) - \underline{T(\vec{x}_\perp, \vec{y}_\perp)} - \underline{T(\vec{x}_\perp, \vec{z}_\perp)T(\vec{z}_\perp, \vec{y}_\perp)} \right\}$$

(Balitsky-Kovchegov equation)

- Both $T = 0$ and $T = 1$ are fixed points of this equation

$$T = \epsilon : \quad \text{r.h.s.} > 0 \quad \Rightarrow \quad T = 0 \text{ is unstable}$$

$$T = 1 - \epsilon : \quad \text{r.h.s.} > 0 \quad \Rightarrow \quad T = 1 \text{ is stable}$$

- Without the non-linear term (underlined), one would have the BFKL equation, that has only an unstable fixed point at $T = 0$



JIMWLK equation

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- An alternate – equivalent – point of view is to keep the bare Wilson lines, and to say that one boosts the target so that the distribution of its color sources changes

- Evolution equation for $W_Y[\rho]$ (JIMWLK) :

$$\frac{\partial W_Y[\rho]}{\partial Y} = \mathcal{H}[\rho] W_Y[\rho]$$

$$\mathcal{H}[\rho] = \int_{\vec{x}_\perp} \sigma(\vec{x}_\perp) \frac{\delta}{\delta \rho(\vec{x}_\perp)} + \frac{1}{2} \int_{\vec{x}_\perp, \vec{y}_\perp} \chi(\vec{x}_\perp, \vec{y}_\perp) \frac{\delta^2}{\delta \rho(\vec{x}_\perp) \delta \rho(\vec{y}_\perp)}$$

- σ and χ are non-linear functionals of ρ
- Note : this point of view is more general because it also applies to situations where the observable cannot be written as a certain combination of Wilson lines – e.g. the gluon inclusive spectrum in AA collisions

Initial condition - MV model

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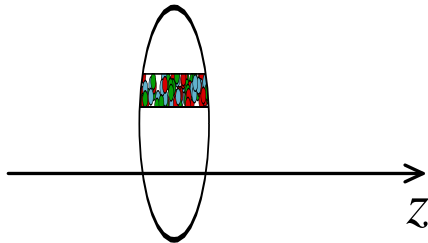
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● MV model

Phenomenology of saturation

- The JIMWLK equation must be completed by an initial condition, given at some moderate x_0
- As with DGLAP, the problem of finding the initial condition is in general non-perturbative
- The **McLerran-Venugopalan** model is often used as an initial condition at moderate x_0 for a **large nucleus** :



- ◆ partons distributed randomly
- ◆ many partons in a small tube
- ◆ no correlations at different \vec{x}_\perp

- The MV model assumes that the density of color charges $\rho(\vec{x}_\perp)$ has a **Gaussian** distribution :

$$W_{x_0}[\rho] = \exp \left[- \int d^2 \vec{x}_\perp \frac{\rho_a(\vec{x}_\perp) \rho_a(\vec{x}_\perp)}{2\mu^2(\vec{x}_\perp)} \right]$$



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- Shadowing

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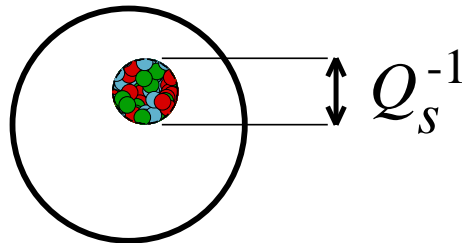
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- In a nucleon at low energy, the typical correlation length among color charges is of the order of the nucleon size, i.e. $\Lambda_{QCD}^{-1} \sim 1 \text{ fm}$. This is because the typical color screening distance is Λ_{QCD}^{-1} . At low energy, color screening is due to confinement, and thus non-perturbative
- At high energy (small x), partons are much more densely packed, and it can be shown that color neutralization occurs in fact over distances of the order of $Q_s^{-1} \ll \Lambda_{QCD}^{-1}$



- This implies that all hadrons, and nuclei, behave in the same way at high energy. In this sense, the small x regime described by the CGC is universal

Multiple scatterings

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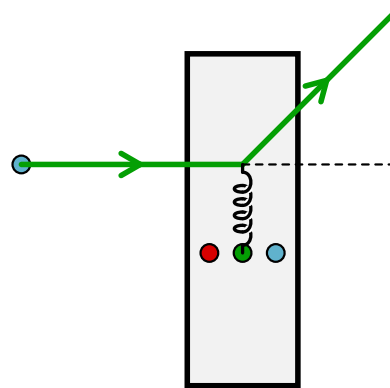
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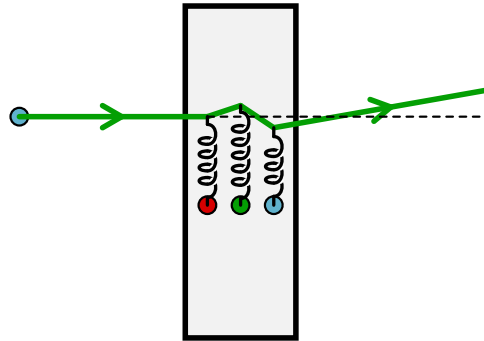
- Single scattering dominates at high p_{\perp} :



- ◆ Differential cross-sections between a parton and a nucleus at high p_{\perp} should scale like the atomic number A (**volume scaling**)

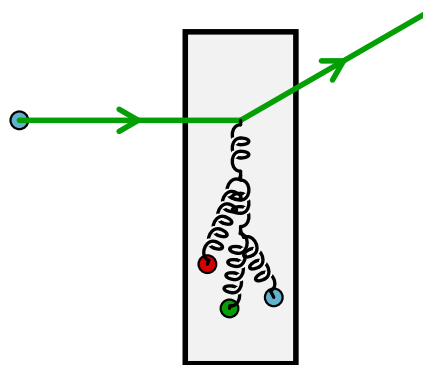
Multiple scatterings

■ Multiple scatterings at low p_{\perp} :



- ◆ One of the scatterings “produces” the final state, while the others merely change its momentum
- ◆ Each extra scattering corresponds to a correction $\alpha_s A^{1/3} \Lambda^2 / p_{\perp}^2$
 - ▷ important correction at low p_{\perp} , despite the α_s suppression
- ◆ When this effect is extremal, differential cross-sections at low p_{\perp} scale like $A^{2/3}$ (**area scaling**)
- ◆ Multiple scatterings only affect the momentum distribution of the final states, not the yield ▷ the suppression at low p_{\perp} is compensated by an increase at higher p_{\perp} (**Cronin effect**)

- Interactions among the partons in the nuclear target (**shadowing**) :



- ◆ Modification of the single scattering contribution due to the non-linear interactions of partons inside the target
- ◆ At low x , this effect induces a suppression of the differential cross-section : $d\sigma_{pA}/d^2\vec{p}_\perp \sim A^\alpha$ with $\alpha < 1$



Lecture II : Initial particle production

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Outline of lecture II

- Introduction to nucleus-nucleus collisions
- Power counting and bookkeeping
- Inclusive gluon spectrum
- Loop corrections and factorization



Lecture III : Instabilities, thermalization

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Outline of lecture III

- Reminder on initial gluon production
- Glasma instabilities
- Instabilities in anisotropic plasmas
- Thermalization ?



Lecture IV : Kinetic theory

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Outline of lecture IV

- Collisionless kinetic equations
- Boltzmann equation
- Transport coefficients
- From kinetic theory to hydrodynamics



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Complements



Light-cone coordinates

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- **Light-cone coordinates** are defined by choosing a privileged axis (generally the z axis) along which particles have a large momentum. Then, for any 4-vector a^μ , one defines :

$$a^+ \equiv \frac{a^0 + a^3}{\sqrt{2}} \quad , \quad a^- \equiv \frac{a^0 - a^3}{\sqrt{2}}$$

$$a^{1,2} \text{ unchanged.} \quad \text{Notation : } \vec{a}_\perp \equiv (a^1, a^2)$$

- Under a Lorentz boost in the z direction :

$$a^+ \rightarrow \Lambda a^+ \quad , \quad a^- \rightarrow \Lambda^{-1} a^- \quad , \quad a^{1,2} \rightarrow a^{1,2}$$

- Some useful formulas :

$$x \cdot y = x^+ y^- + x^- y^+ - \vec{x}_\perp \cdot \vec{y}_\perp$$

$$d^4x = dx^+ dx^- d^2\vec{x}_\perp$$

$$\square = 2\partial^+ \partial^- - \vec{\nabla}_\perp^2 \quad \text{Notation : } \partial^+ \equiv \frac{\partial}{\partial x^-} \quad , \quad \partial^- \equiv \frac{\partial}{\partial x^+}$$



Solution of Yang Mills eq. for 1 nucleus

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- The Yang-Mills equations are the classical equations that give the color field induced by a given current. They are the analogue in QCD of the Maxwell equations in electrodynamics :

$$[D_\mu, F^{\mu\nu}] = J^\nu$$

where

- ◆ $D^\mu \equiv \partial^\mu - igA^\mu$ is the covariant derivative
- ◆ $F^{\mu\nu} \equiv i[D^\mu, D^\nu]/g = \partial^\mu A^\nu - \partial^\nu A^\mu - ig[A^\mu, A^\nu]$

- For a single nucleus moving in the $+z$ direction

$$J^\nu = \delta^{\nu+} \delta(x^-) \rho(\vec{x}_\perp)$$

- This current must be covariantly conserved :

$$[D_\nu, J^\nu] = 0$$

- ▷ the formula for J^ν may have higher order corrections in ρ



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- Reminder: in QED, we would simply have :

$$\square A^\nu - \partial^\nu (\partial_\mu A^\mu) = J^\nu$$

which in the gauge $\partial_\mu A^\mu = 0$ (Lorenz) simplifies into

$$\square A^\nu = J^\nu$$

- For QCD in the Lorenz gauge, one can first rewrite the Yang-Mills equations as :

$$\square A^\nu - ig [A_\mu, F^{\mu\nu} + \partial^\mu A^\nu] = J^\nu$$

- It is useful to expand all the quantities in powers of the color density ρ :

$$A^\mu \equiv \sum_{n=0}^{\infty} A_{(n)}^\mu, \quad J^\mu \equiv \sum_{n=0}^{\infty} J_{(n)}^\mu$$

where $A_{(n)}^\mu \sim J_{(n)}^\mu \sim \rho^n$



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- At the order $n = 1$, the equations are simply :

$$\partial_\nu J_{(1)}^\nu = 0 \quad , \quad \square A_{(1)}^\nu = J_{(1)}^\nu$$

and their solution reads

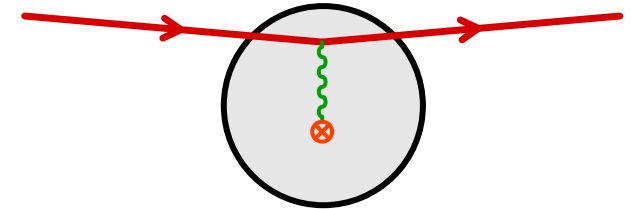
$$J_{(1)}^\nu = \delta^{\nu+} \delta(x^-) \rho(\vec{x}_\perp)$$

$$A_{(1)}^\nu = \delta^{\nu+} \delta(x^-) \alpha(\vec{x}_\perp) \quad , \quad -\vec{\nabla}_\perp^2 \alpha(\vec{x}_\perp) = \rho(\vec{x}_\perp)$$

- At this order, the only non-zero component of $F^{\mu\nu}$ is
 $F^{i+} = -F^{+i} = \partial^i A_{(1)}^+$
- By writing the equations for the corrections of order ρ^2 , we find that all the non-linear terms cancel and that all these corrections are zero
- This feat can be repeated to all orders in ρ \triangleright the complete solution of the non-linear Yang-Mills equations is linear in ρ !
Note : this result is only valid for Lorenz gauge and for the kind of current we have considered

Eikonal scattering

- Consider the scattering amplitude off an external potential :



$$S_{\beta\alpha} \equiv \langle \beta_{\text{out}} | \alpha_{\text{in}} \rangle = \langle \beta_{\text{in}} | U(+\infty, -\infty) | \alpha_{\text{in}} \rangle$$

where $U(+\infty, -\infty)$ is the evolution operator from $t = -\infty$ to $t = +\infty$

$$U(+\infty, -\infty) = T \exp \left[i \int d^4x \mathcal{L}_{\text{int}}(\phi_{\text{in}}(x)) \right]$$

Note : \mathcal{L}_{int} contains the self-interactions of the fields and their interactions with the external potential

- We want to calculate its high energy limit :

$$S_{\beta\alpha}^{(\infty)} \equiv \lim_{\omega \rightarrow +\infty} \langle \beta_{\text{in}} | e^{i\omega K^3} U(+\infty, -\infty) e^{-i\omega K^3} | \alpha_{\text{in}} \rangle$$

where K^3 is the generator of boosts in the $+z$ direction



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- In a scattering at high energy, the collision time goes to zero as $s^{-1/2}$
 - With **scalar interactions**, this implies a decrease of the scattering amplitude as $s^{-1/2}$
 - With **vectorial interactions**, this decrease is compensated by the growth of the component J^+ of the vector current
- ▷ the **eikonal approximation** gives the finite limit of the scattering amplitude in the case of vectorial interactions when $s \rightarrow +\infty$



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- Consider an external vector potential, that couples via $e \mathcal{A}_\mu(x) J^\mu(x)$ (J^μ is the Noether current associated to some conserved charge)
- We will assume that the external potential is non-zero only in a finite range in x^+ , $x^+ \in [-L, +L]$
- The action of K^3 on states and (scalar) fields is :

$$e^{-i\omega K^3} |\vec{p} \cdots \text{in}\rangle = |(e^\omega p^+, \vec{p}_\perp) \cdots \text{in}\rangle$$

$$e^{i\omega K^3} \phi_{\text{in}}(x) e^{-i\omega K^3} = \phi_{\text{in}}(e^{-\omega} x^+, e^\omega x^-, \vec{x}_\perp)$$

- K^3 does not change the ordering in x^+ . Hence,

$$e^{i\omega K^3} U(+\infty, -\infty) e^{-i\omega K^3} = T_+ \exp i \int d^4x \mathcal{L}_{\text{int}}(e^{i\omega K^3} \phi_{\text{in}}(x) e^{-i\omega K^3})$$

where $\mathcal{L}_{\text{int}} = \mathcal{L}_{\text{self}}(\phi) - e \mathcal{A}_\mu J^\mu$



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- Split the evolution operator $U(+\infty, -\infty)$ into three factors :

$$U(+\infty, -\infty) = U(+\infty, +L)U(+L, -L)U(-L, -\infty)$$

Upon application of K^3 , this becomes :

$$\begin{aligned} e^{i\omega K^3} U(+\infty, -\infty) e^{-i\omega K^3} &= e^{i\omega K^3} U(+\infty, +L) e^{-i\omega K^3} \\ &\times e^{i\omega K^3} U(+L, -L) e^{-i\omega K^3} e^{i\omega K^3} U(-L, -\infty) e^{-i\omega K^3} \end{aligned}$$

- The external potential $\mathcal{A}_\mu(x)$ is unaffected by K^3
- The components of $J^\mu(x)$ are changed as follows :

$$e^{i\omega K^3} J^i(x) e^{-i\omega K^3} = J^i(e^{-\omega} x^+, e^\omega x^-, \vec{x}_\perp)$$

$$e^{i\omega K^3} J^-(x) e^{-i\omega K^3} = e^{-\omega} J^-(e^{-\omega} x^+, e^\omega x^-, \vec{x}_\perp)$$

$$e^{i\omega K^3} J^+(x) e^{-i\omega K^3} = e^\omega J^+(e^{-\omega} x^+, e^\omega x^-, \vec{x}_\perp)$$



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- The factors $U(+\infty, +L)$ and $U(-L, -\infty)$ do not contain the external potential. In order to deal with these factors, it is sufficient to change variables : $e^{-\omega}x^+ \rightarrow x^+$, $e^{\omega}x^- \rightarrow x^-$. This leads to :

$$\lim_{\omega \rightarrow +\infty} e^{i\omega K^3} U(+\infty, +L) e^{-i\omega K^3} = U_{\text{self}}(+\infty, 0)$$

$$\lim_{\omega \rightarrow +\infty} e^{i\omega K^3} U(-L, -\infty) e^{-i\omega K^3} = U_{\text{self}}(0, -\infty)$$

where U_{self} is the same as U , but with the self-interactions only

- For the factor $U(L, -L)$, the change $e^{\omega}x^- \rightarrow x^-$ leads to :

$$\begin{aligned} e^{i\omega K^3} U(+L, -L) e^{-i\omega K^3} &= \\ &= T_+ \exp i \int_{-L}^{+L} d^4x e^{-\omega} \left[e \mathcal{A}^-(x^+, e^{-\omega}x^-, \vec{x}_\perp) \right. \\ &\quad \left. \times e^{\omega} J^+(e^{-\omega}x^+, x^-, \vec{x}_\perp) + \mathcal{O}(1) \right] \end{aligned}$$



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- Therefore, in the limit $\omega \rightarrow +\infty$, we have :

$$\lim_{\omega \rightarrow +\infty} e^{i\omega K^3} U(+L, -L) e^{-i\omega K^3} = \exp \left[i e \int d^2 \vec{x}_\perp \chi(\vec{x}_\perp) \rho(\vec{x}_\perp) \right]$$

$$\text{with} \quad \begin{cases} \chi(\vec{x}_\perp) \equiv \int dx^+ \mathcal{A}^-(x^+, 0, \vec{x}_\perp) \\ \rho(\vec{x}_\perp) \equiv \int dx^- J^+(0, x^-, \vec{x}_\perp) \end{cases}$$

- The high-energy limit of the scattering amplitude is :

$$S_{\beta\alpha}^{(\infty)} = \langle \beta_{\text{in}} | U_{\text{self}}(+\infty, 0) \exp \left[i e \int_{\vec{x}_\perp} \chi(\vec{x}_\perp) \rho(\vec{x}_\perp) \right] U_{\text{self}}(0, -\infty) | \alpha_{\text{in}} \rangle$$

- ◆ Only the – component of the **vector potential** matters
- ◆ The self-interactions and the interactions with the external potential are factorized \triangleright **parton model**
- ◆ This is an exact result when $s \rightarrow +\infty$



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- The previous formula still contains all the self-interactions of the fields. In order to perform the perturbative expansion, it is convenient to write first :

$$S_{\beta\alpha}^{(\infty)} = \sum_{\gamma,\delta} \langle \beta_{\text{in}} | U_{\text{self}}(+\infty, 0) | \gamma_{\text{in}} \rangle \\ \times \langle \gamma_{\text{in}} | \exp \left[i e \int_{\vec{x}_{\perp}} \chi(\vec{x}_{\perp}) \rho(\vec{x}_{\perp}) \right] | \delta_{\text{in}} \rangle \langle \delta_{\text{in}} | U_{\text{self}}(0, -\infty) | \alpha_{\text{in}} \rangle$$

- The factor

$$\sum_{\delta} | \delta_{\text{in}} \rangle \langle \delta_{\text{in}} | U_{\text{self}}(0, -\infty) | \alpha_{\text{in}} \rangle$$

is the **Fock expansion** of the initial state: the state prepared at $x^+ = -\infty$ may have fluctuated into another state before it interacts with the external potential



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- We need to calculate matrix elements such as $\langle \gamma_{\text{in}} | \mathbf{F} | \delta_{\text{in}} \rangle$, with :

$$\mathbf{F} \equiv \exp i e \int \chi_a(\vec{x}_\perp) \rho^a(\vec{x}_\perp)$$

- ◆ having QCD in mind, we have reinstated the color indices
- ◆ the contribution of quarks and antiquarks to $\rho^a(\vec{x}_\perp)$ is :

$$\rho^a(\vec{x}_\perp) = t_{ij}^a \int \frac{dp^+}{4\pi p^+} \frac{d^2 \vec{p}_\perp}{(2\pi)^2} \frac{d^2 \vec{q}_\perp}{(2\pi)^2} \left\{ b_{\text{in}}^\dagger(p^+, \vec{p}_\perp; i) b_{\text{in}}(p^+, \vec{q}_\perp; j) e^{i(\vec{p}_\perp - \vec{q}_\perp) \cdot \vec{x}_\perp} \right. \\ \left. - d_{\text{in}}^\dagger(p^+, \vec{p}_\perp; i) d_{\text{in}}(p^+, \vec{q}_\perp; j) e^{-i(\vec{p}_\perp - \vec{q}_\perp) \cdot \vec{x}_\perp} \right\}$$

- ◆ Note : one should keep the ordering of the exponential in x^+
 - ◆ the contribution of gluons is similar, with a color matrix in the adjoint representation
- The action of \mathbf{F} on a state $|\delta_{\text{in}}\rangle$ gives a state with the same particle content, the same $+$ components for the momenta, but modified transverse momenta and colors

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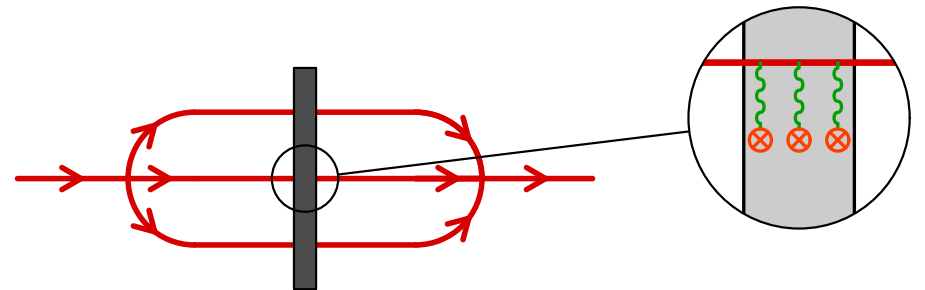
- For each intermediate state $\langle \delta_{\text{in}} | \equiv \langle \{k_i^+, \vec{k}_{i\perp}\} |$, define the corresponding **light-cone wave function** by :

$$\Psi_{\delta\alpha}(\{k_i^+, \vec{x}_{i\perp}\}) \equiv \prod_i \int \frac{d^2 \vec{k}_{i\perp}}{(2\pi)^2} e^{-i\vec{k}_{i\perp} \cdot \vec{x}_{i\perp}} \langle \delta_{\text{in}} | U_{\text{self}}(0, -\infty) | \alpha_{\text{in}} \rangle$$

- Each charged particle going through the external field acquires a **phase proportional to its charge** (antiparticles get an opposite phase) :

$$\Psi_{\delta\alpha}(\{k_i^+, \vec{x}_{i\perp}\}) \longrightarrow \Psi_{\delta\alpha}(\{k_i^+, \vec{x}_{i\perp}\}) \prod_i U_i(\vec{x}_{i\perp})$$

$$U_i(\vec{x}_{i\perp}) \equiv T_+ \exp \left[ig_i \int dx^+ \mathcal{A}_a^-(x^+, 0, \vec{x}_{i\perp}) t^a \right]$$



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- We have seen that the number and the nature of the particles is unchanged under the action of the operator F . Moreover, in terms of the transverse coordinates, we simply have

$$\langle \gamma_{\text{in}} | F | \delta_{\text{in}} \rangle = \delta_{NN'} \prod_i \left[4\pi k_i^+ \delta(k_i^+ - k_i^{+'}) \delta(\vec{x}_{i\perp} - \vec{x}'_{i\perp}) U_{R_i}(\vec{x}_{i\perp}) \right]$$

where $U_R(\vec{x}_\perp)$ is a Wilson line operator, in the representation R appropriate for the particle going through the target

- Therefore, the high energy scattering amplitude can be written as :

$$S_{\beta\alpha}^{(\infty)} = \sum_\delta \int \left[\prod_{i \in \delta} d\Phi_i \right] \Psi_{\delta\beta}^\dagger(\{k_i^+, \vec{x}_{i\perp}\}) \left[\prod_{i \in \delta} U_{R_i}(\vec{x}_{i\perp}) \right] \Psi_{\delta\alpha}(\{k_i^+, \vec{x}_{i\perp}\})$$

- As we shall see shortly, some loop corrections are enhanced by logs of the energy. They must be resummed and drive the energy evolution of the amplitude

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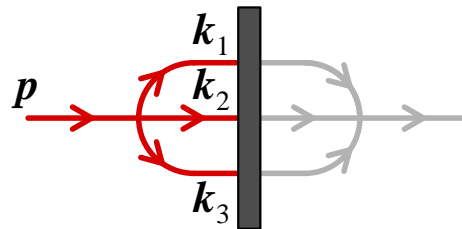
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- The calculation of $\langle \delta_{\text{in}} | U_{\text{self}}(0, -\infty) | \alpha_{\text{in}} \rangle$ is similar to that of scattering amplitudes in the vacuum. The only difference is that the integration over x^+ at each vertex runs only over half of the real axis $[-\infty, 0]$
 - ◆ In Fourier space, this means that the $-$ component of the momentum is not conserved at the vertices
 - ◆ Instead of a δ function, one gets an energy denominator
- Example with a single interaction :



$$\begin{aligned}
 \langle \vec{k}_1 \vec{k}_2 \vec{k}_3 | U(0, -\infty) | \vec{p}_{\text{in}} \rangle &= -ig \int_{-\infty}^0 d^4x e^{i(k_1 + k_2 + k_3 - p) \cdot x} \\
 &= -g \frac{(2\pi)^3 \delta(\vec{k}_{1\perp} + \vec{k}_{2\perp} + \vec{k}_{3\perp} - \vec{p}_{\perp}) \delta(k_1^+ + k_2^+ + k_3^+ - p^+)}{k_1^- + k_2^- + k_3^- - p^- - i\epsilon}
 \end{aligned}$$

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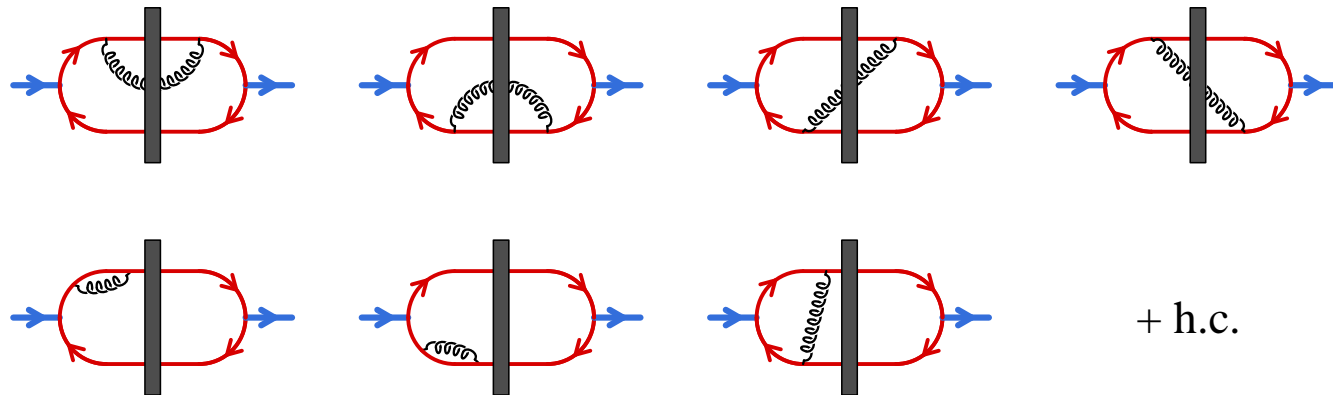
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- Assume that the initial and final states are a **color singlet** $Q\bar{Q}$ dipole. The bare scattering amplitude can be written as :

$$\propto \left| \Psi^{(0)}(\vec{x}_\perp, \vec{y}_\perp) \right|^2 \text{tr} \left[U(\vec{x}_\perp) U^\dagger(\vec{y}_\perp) \right]$$

- At one loop, the following diagrams must be evaluated :



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- In the gauge $A^+ = 0$, the emission of a gluon of momentum k by a quark can be written as :



$$= 2gt^a \frac{\vec{\epsilon}_\lambda \cdot \vec{k}_\perp}{k_\perp^2}$$

- In coordinate space, this reads :

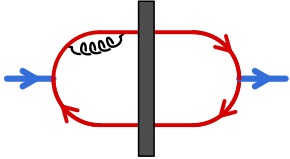
$$\int \frac{d^2 \vec{k}_\perp}{(2\pi)^2} e^{i\vec{k}_\perp \cdot (\vec{x}_\perp - \vec{z}_\perp)} 2gt^a \frac{\vec{\epsilon}_\lambda \cdot \vec{k}_\perp}{k_\perp^2} = \frac{2ig}{2\pi} t^a \frac{\vec{\epsilon}_\lambda \cdot (\vec{x}_\perp - \vec{z}_\perp)}{(\vec{x}_\perp - \vec{z}_\perp)^2}$$

- When connecting two gluons, one must use :

$$\sum_\lambda \vec{\epsilon}_\lambda^i \vec{\epsilon}_\lambda^j = -g^{ij}$$

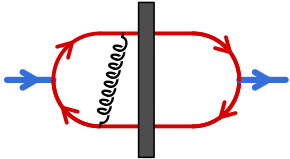
Virtual corrections

- Consider first the loop corrections inside the wavefunction of the incoming or outgoing dipole
- Examples :



$$= \left| \Psi^{(0)}(\vec{x}_\perp, \vec{y}_\perp) \right|^2 \text{tr} \left[t^a t^a U(\vec{x}_\perp) U^\dagger(\vec{y}_\perp) \right]$$

$$\times -2\alpha_s \int \frac{dk^+}{k^+} \int \frac{d^2 \vec{z}_\perp}{(2\pi)^2} \frac{(\vec{x}_\perp - \vec{z}_\perp) \cdot (\vec{x}_\perp - \vec{z}_\perp)}{(\vec{x}_\perp - \vec{z}_\perp)^2 (\vec{x}_\perp - \vec{z}_\perp)^2}$$



$$= \left| \Psi^{(0)}(\vec{x}_\perp, \vec{y}_\perp) \right|^2 \text{tr} \left[t^a U(\vec{x}_\perp) U^\dagger(\vec{y}_\perp) t^a \right]$$

$$\times 4\alpha_s \int \frac{dk^+}{k^+} \int \frac{d^2 \vec{z}_\perp}{(2\pi)^2} \frac{(\vec{x}_\perp - \vec{z}_\perp) \cdot (\vec{y}_\perp - \vec{z}_\perp)}{(\vec{x}_\perp - \vec{z}_\perp)^2 (\vec{y}_\perp - \vec{z}_\perp)^2}$$

- Reminder : $t^a t^a = (N_c^2 - 1)/2N_c \equiv C_F$

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- The sum of all virtual corrections is :

$$-\frac{C_F \alpha_s}{\pi^2} \int \frac{dk^+}{k^+} \int d^2 \vec{z}_\perp \frac{(\vec{x}_\perp - \vec{y}_\perp)^2}{(\vec{x}_\perp - \vec{z}_\perp)^2 (\vec{y}_\perp - \vec{z}_\perp)^2} \times \left| \Psi^{(0)}(\vec{x}_\perp, \vec{y}_\perp) \right|^2 \text{tr} \left[U(\vec{x}_\perp) U^\dagger(\vec{y}_\perp) \right]$$

- The integral over k^+ is divergent. It should have an upper bound at p^+ :

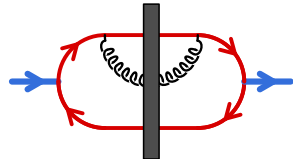
$$\int^{p^+} \frac{dk^+}{k^+} = \ln(p^+) = Y$$

▷ When Y is large, $\alpha_s Y$ may not be small. By differentiating with respect to Y , we will get an evolution equation in Y whose solution resums all the powers $(\alpha_s Y)^n$

- The integral over \vec{z}_\perp is divergent when $\vec{z}_\perp = \vec{x}_\perp$ or \vec{y}_\perp

Real corrections

- There are also real corrections, for which the state that interacts with the target has an extra gluon
- Example :



$$= \left| \Psi^{(0)}(\vec{x}_\perp, \vec{y}_\perp) \right|^2 \text{tr} \left[t^a U(\vec{x}_\perp) t^b U^\dagger(\vec{y}_\perp) \right]$$

$$\times 4\alpha_s \int \frac{dk^+}{k^+} \int \frac{d^2 \vec{z}_\perp}{(2\pi)^2} \tilde{U}_{ab}(\vec{z}_\perp) \frac{(\vec{x}_\perp - \vec{z}_\perp) \cdot (\vec{x}_\perp - \vec{z}_\perp)}{(\vec{x}_\perp - \vec{z}_\perp)^2 (\vec{x}_\perp - \vec{z}_\perp)^2}$$

◆ $\tilde{U}_{ab}(\vec{z}_\perp)$ is a Wilson line in the **adjoint representation**

- In order to simplify the color structure, first recall that :

$$t^a \tilde{U}_{ab}(\vec{z}_\perp) = U(\vec{z}_\perp) t^b U^\dagger(\vec{z}_\perp)$$

- Then use the $SU(N_c)$ **Fierz identity** :

$$t_{ij}^b t_{kl}^b = \frac{1}{2} \delta_{il} \delta_{jk} - \frac{1}{2N_c} \delta_{ij} \delta_{kl}$$

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- The Wilson lines can be rearranged into :

$$\text{tr} \left[t^a U(\vec{x}_\perp) t^b U^\dagger(\vec{y}_\perp) \right] \tilde{U}_{ab}(\vec{z}_\perp) = \frac{1}{2} \text{tr} \left[U^\dagger(\vec{z}_\perp) U(\vec{x}_\perp) \right] \text{tr} \left[U(\vec{z}_\perp) U^\dagger(\vec{y}_\perp) \right] - \frac{1}{2N_c} \text{tr} \left[U(\vec{x}_\perp) U^\dagger(\vec{y}_\perp) \right]$$

- ◆ The term in $1/2N_c$ cancels against a similar term in the virtual contribution
- ◆ All the real terms have the same color structure
- When we sum all the real terms, we generate the same kernel as in the virtual terms :

$$\frac{(\vec{x}_\perp - \vec{y}_\perp)^2}{(\vec{x}_\perp - \vec{z}_\perp)^2 (\vec{y}_\perp - \vec{z}_\perp)^2}$$

- In order to simplify the notations, let us denote :

$$S(\vec{x}_\perp, \vec{y}_\perp) \equiv \frac{1}{N_c} \text{tr} \left[U(\vec{x}_\perp) U^\dagger(\vec{y}_\perp) \right]$$



Evolution equation

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- The 1-loop scattering amplitude reads :

$$-\frac{\alpha_s N_c^2 Y}{2\pi^2} \left| \Psi^{(0)}(\vec{x}_\perp, \vec{y}_\perp) \right|^2 \int d^2 \vec{z}_\perp \frac{(\vec{x}_\perp - \vec{y}_\perp)^2}{(\vec{x}_\perp - \vec{z}_\perp)^2 (\vec{y}_\perp - \vec{z}_\perp)^2} \times \left\{ \mathbf{S}(\vec{x}_\perp, \vec{y}_\perp) - \mathbf{S}(\vec{x}_\perp, \vec{z}_\perp) \mathbf{S}(\vec{z}_\perp, \vec{y}_\perp) \right\}$$

- Reminder: the bare scattering amplitude was :

$$\left| \Psi^{(0)}(\vec{x}_\perp, \vec{y}_\perp) \right|^2 N_c \mathbf{S}(\vec{x}_\perp, \vec{y}_\perp)$$

- Hence, we have :

$$\frac{\partial \mathbf{S}(\vec{x}_\perp, \vec{y}_\perp)}{\partial Y} = -\frac{\alpha_s N_c}{2\pi^2} \int d^2 \vec{z}_\perp \frac{(\vec{x}_\perp - \vec{y}_\perp)^2}{(\vec{x}_\perp - \vec{z}_\perp)^2 (\vec{y}_\perp - \vec{z}_\perp)^2} \times \left\{ \mathbf{S}(\vec{x}_\perp, \vec{y}_\perp) - \mathbf{S}(\vec{x}_\perp, \vec{z}_\perp) \mathbf{S}(\vec{z}_\perp, \vec{y}_\perp) \right\}$$

- ◆ since $\mathbf{S}(\vec{x}_\perp, \vec{x}_\perp) = 1$, the integral over \vec{z}_\perp is now regular



BFKL equation

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Kuraev, Lipatov, Fadin (1977), Balitsky, Lipatov (1978)

- The BFKL equation can be obtained by linearizing the previous equation
- Write $S(\vec{x}_\perp, \vec{y}_\perp) \equiv 1 - T(\vec{x}_\perp, \vec{y}_\perp)$ and assume that we are in the **dilute regime**, so that the scattering amplitude T is small. Drop the terms that are non-linear in T :

$$\frac{\partial T(\vec{x}_\perp, \vec{y}_\perp)}{\partial Y} = \frac{\alpha_s N_c}{2\pi^2} \int d^2 \vec{z}_\perp \frac{(\vec{x}_\perp - \vec{y}_\perp)^2}{(\vec{x}_\perp - \vec{z}_\perp)^2 (\vec{y}_\perp - \vec{z}_\perp)^2} \times \left\{ T(\vec{x}_\perp, \vec{z}_\perp) + T(\vec{z}_\perp, \vec{y}_\perp) - T(\vec{x}_\perp, \vec{y}_\perp) \right\}$$

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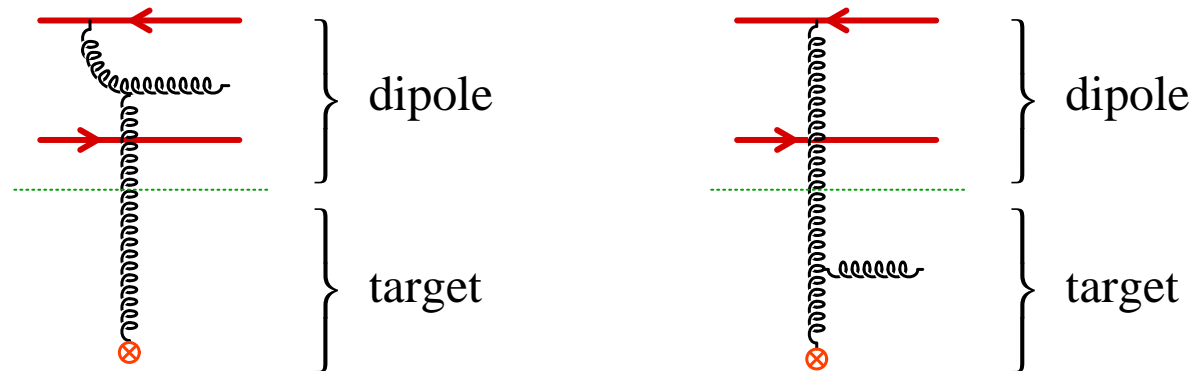
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- Note : $T(\vec{x}_\perp, \vec{y}_\perp)$ is independent on the frame. In particular, it depends only on the rapidity difference between the dipole and the target
 - ▷ in a frame where the dipole is held fixed, the target has to evolve in such a way as to reproduce the Y dependence of T



- The corresponding evolution in the target is the radiation of a gluon



Unitarity problem

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- The solution of this equation grows exponentially when $Y \rightarrow +\infty$ \triangleright serious unitarity problem...

- In perturbation theory, the forward scattering amplitude between a small dipole and a target made of gluons reads :

$$T(\vec{x}_\perp, \vec{y}_\perp) \propto |\vec{x}_\perp - \vec{y}_\perp|^2 xG(x, |\vec{x}_\perp - \vec{y}_\perp|^{-2})$$

where $Y \equiv \ln(1/x)$

- Therefore, the exponential behavior of T implies an increase of the gluon distribution at small x

$$T \sim e^{\lambda Y} \quad \longleftrightarrow \quad xG(x, Q^2) \sim \frac{1}{x^\lambda}$$



Non-linear evolution equation

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- In fact, the first evolution equation we derived has a bounded solution. The unbounded solutions of BFKL are due to dropping the non-linear term. The full equation reads :

$$\frac{\partial \mathbf{T}(\vec{x}_\perp, \vec{y}_\perp)}{\partial Y} = \frac{\alpha_s N_c}{2\pi^2} \int d^2 \vec{z}_\perp \frac{(\vec{x}_\perp - \vec{y}_\perp)^2}{(\vec{x}_\perp - \vec{z}_\perp)^2 (\vec{y}_\perp - \vec{z}_\perp)^2} \times \left\{ \mathbf{T}(\vec{x}_\perp, \vec{z}_\perp) + \mathbf{T}(\vec{z}_\perp, \vec{y}_\perp) - \mathbf{T}(\vec{x}_\perp, \vec{y}_\perp) - \mathbf{T}(\vec{x}_\perp, \vec{z}_\perp) \mathbf{T}(\vec{z}_\perp, \vec{y}_\perp) \right\}$$

(Balitsky-Kovchegov equation)

- The r.h.s. vanishes when \mathbf{T} reaches 1, and the growth stops. The non-linear term lets both dipoles interact after the splitting of the original dipole
- Both $\mathbf{T} = 0$ and $\mathbf{T} = 1$ are fixed points of this equation

$$\mathbf{T} = \epsilon : \quad \text{r.h.s.} > 0 \quad \Rightarrow \quad \mathbf{T} = 0 \text{ is unstable}$$

$$\mathbf{T} = 1 - \epsilon : \quad \text{r.h.s.} > 0 \quad \Rightarrow \quad \mathbf{T} = 1 \text{ is stable}$$



Caveats and improvements

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- So far, we have studied the scattering amplitude between a color dipole and a “god given” patch of color field. This is too naive to describe any realistic situation
- We need to improve the treatment of the target
- An experimentally measured cross-section is an **average over many collisions**, and there is no reason why these fields should be the same in different collisions :

$$\mathbf{T} \rightarrow \langle \mathbf{T} \rangle$$

$\langle \dots \rangle$ denotes the average over the target configurations, i.e.

$$\langle \dots \rangle = \int [D\rho] W_Y[\rho] \dots$$



Balitsky hierarchy

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- Because of this average over the target configurations, the evolution equation we have derived should be written as :

$$\frac{\partial \langle \mathbf{T}(\vec{x}_\perp, \vec{y}_\perp) \rangle}{\partial Y} = \frac{\alpha_s N_c}{2\pi^2} \int d^2 \vec{z}_\perp \frac{(\vec{x}_\perp - \vec{y}_\perp)^2}{(\vec{x}_\perp - \vec{z}_\perp)^2 (\vec{y}_\perp - \vec{z}_\perp)^2} \times \left\{ \langle \mathbf{T}(\vec{x}_\perp, \vec{z}_\perp) \rangle + \langle \mathbf{T}(\vec{z}_\perp, \vec{y}_\perp) \rangle - \langle \mathbf{T}(\vec{x}_\perp, \vec{y}_\perp) \rangle - \langle \mathbf{T}(\vec{x}_\perp, \vec{z}_\perp) \mathbf{T}(\vec{z}_\perp, \vec{y}_\perp) \rangle \right\}$$

- As one can see, the equation is no longer a closed equation, since the equation for $\langle \mathbf{T} \rangle$ depends on a new object, $\langle \mathbf{T} \mathbf{T} \rangle$
- One can derive an evolution equation for $\langle \mathbf{T} \mathbf{T} \rangle$. Its right hand side contains $\langle \mathbf{T} \mathbf{T} \mathbf{T} \rangle$. And so on...
- There is in fact an infinite hierarchy of nested evolution equations : **Balitsky hierarchy**

Note : this hierarchy is equivalent to the JIMWLK equation for $W_Y[\rho]$



Balitsky-Kovchegov equation

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- In order to truncate the hierarchy of equations, one may assume the following simplification

$$\langle \mathbf{T T} \rangle \approx \langle \mathbf{T} \rangle \langle \mathbf{T} \rangle$$

- This approximation gives for $\langle \mathbf{T} \rangle$ the same evolution equation as the one we had for a fixed configuration of the target



Analogy with reaction-diffusion

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Munier, Peschanski (2003,2004)

- Assume rotation invariance, and define :

$$N(Y, k_{\perp}) \equiv 2\pi \int d^2 \vec{x}_{\perp} e^{i\vec{k}_{\perp} \cdot \vec{x}_{\perp}} \frac{\langle \mathbf{T}(0, \vec{x}_{\perp}) \rangle_Y}{x_{\perp}^2}$$

- From the Balitsky-Kovchegov equation for $\langle \mathbf{T} \rangle_Y$, we obtain the following equation for N :

$$\frac{\partial N(Y, k_{\perp})}{\partial Y} = \frac{\alpha_s N_c}{\pi} \left[\chi(-\partial_L) N(Y, k_{\perp}) - N^2(Y, k_{\perp}) \right]$$

with

$$L \equiv \ln(k_{\perp}^2 / k_0^2)$$

$$\chi(\gamma) \equiv 2\psi(1) - \psi(\gamma) - \psi(1 - \gamma)$$



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- Expand the function $\chi(\gamma)$ to second order around its minimum $\gamma = 1/2$

- Introduce new variables :

$$t \sim Y$$

$$z \sim L + \frac{\alpha_s N_c}{2\pi} \chi''(1/2) Y$$

- The equation for N becomes :

$$\partial_t N = \partial_z^2 N + N - N^2$$

(known as the Fisher-Kolmogorov-Petrov-Piscounov equation)



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- **Interpretation** : this equation is typical for all the **diffusive systems** in which a **reaction** $A \longleftrightarrow A + A$ takes place
 - ◆ $\partial_z^2 N$: diffusion term (the quantity under consideration can hop from a site to the neighboring sites)
 - ◆ $+N$: gain term corresponding to $A \rightarrow A + A$
 - ◆ $-N^2$: loss term corresponding to $A + A \rightarrow A$
- **Note** : this equation has two fixed points :
 - ◆ $N = 0$: unstable
 - ◆ $N = 1$: stable
- The stable fixed point at $N = 1$ exists only if one keeps the loss term. In other words, one would not have it from the BFKL equation

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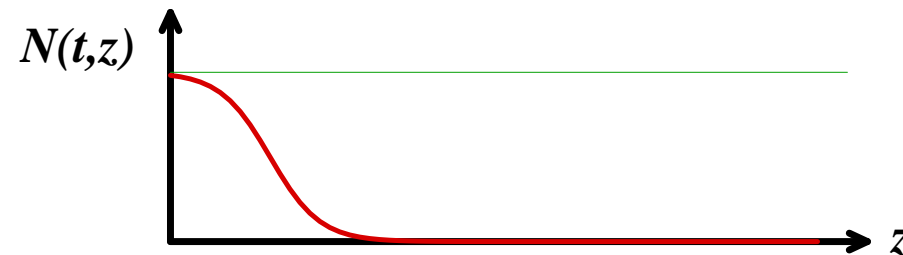
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- Assume an initial condition $N(t_0, z)$ that goes smoothly from 1 at $z = -\infty$ to 0 at $z = +\infty$, and behaves like $\exp(-\beta z)$ when $z \gg 1$



- The solution of the **F-KPP equation** is known to behave like a **traveling wave** at asymptotic times (**Bramson, 1983**) :

$$N(t, z) \underset{t \rightarrow +\infty}{\sim} N(z - m_\beta(t))$$

with $m_\beta(t) = 2t - 3 \ln(t)/2 + \mathcal{O}(1)$ if $\beta > 1$

▷ universal front velocity

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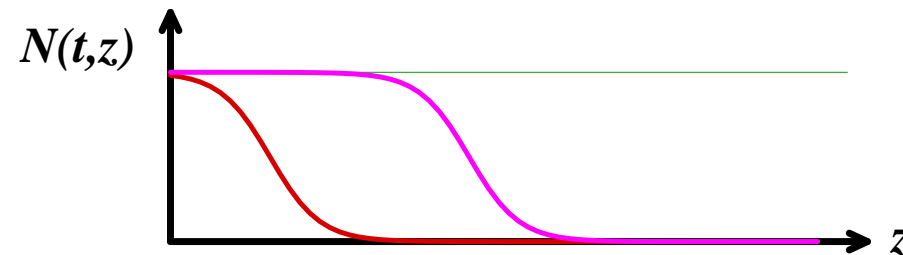
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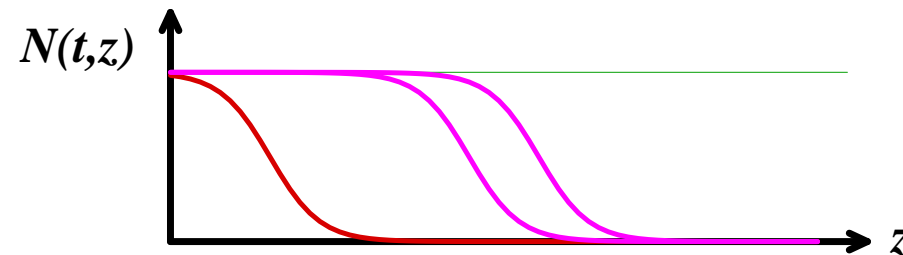
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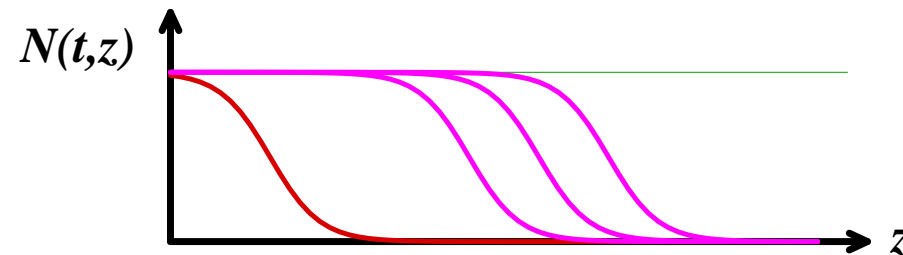
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Iancu, Itakura, McLerran (2002)

Mueller, Triantafyllopoulos (2002)

Munier, Peschanski (2003)

- In QCD, the initial condition is of the required form, with $\beta > 1$
 - ▷ front velocity independent of the initial condition
- Going back to the original variables, one gets :

$$N(Y, k_{\perp}) = N(k_{\perp}/Q_s(Y))$$

with

$$Q_s^2(Y) = k_0^2 Y^{-\delta} e^{\lambda Y}$$

- Going from $N(Y, k_{\perp})$ to $\langle T(0, \vec{x}_{\perp}) \rangle_Y$, we obtain :

$$\langle T(0, \vec{x}_{\perp}) \rangle_Y = T(Q_s(Y) x_{\perp})$$



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- The **total γ^*p cross-section**, measured in **Deep Inelastic Scattering**, can be written in terms of N :

$$\sigma_{\gamma^*p}^{\text{tot}}(Y, Q^2) = 2\pi R^2 \int d^2\vec{x}_\perp \int_0^1 dz |\psi(z, \vec{x}_\perp, Q^2)|^2 \langle \mathbf{T}(0, \vec{x}_\perp) \rangle_Y$$

- ◆ The photon wavefunction ψ is calculable in QED. It depends on the dipole size \vec{x}_\perp only via

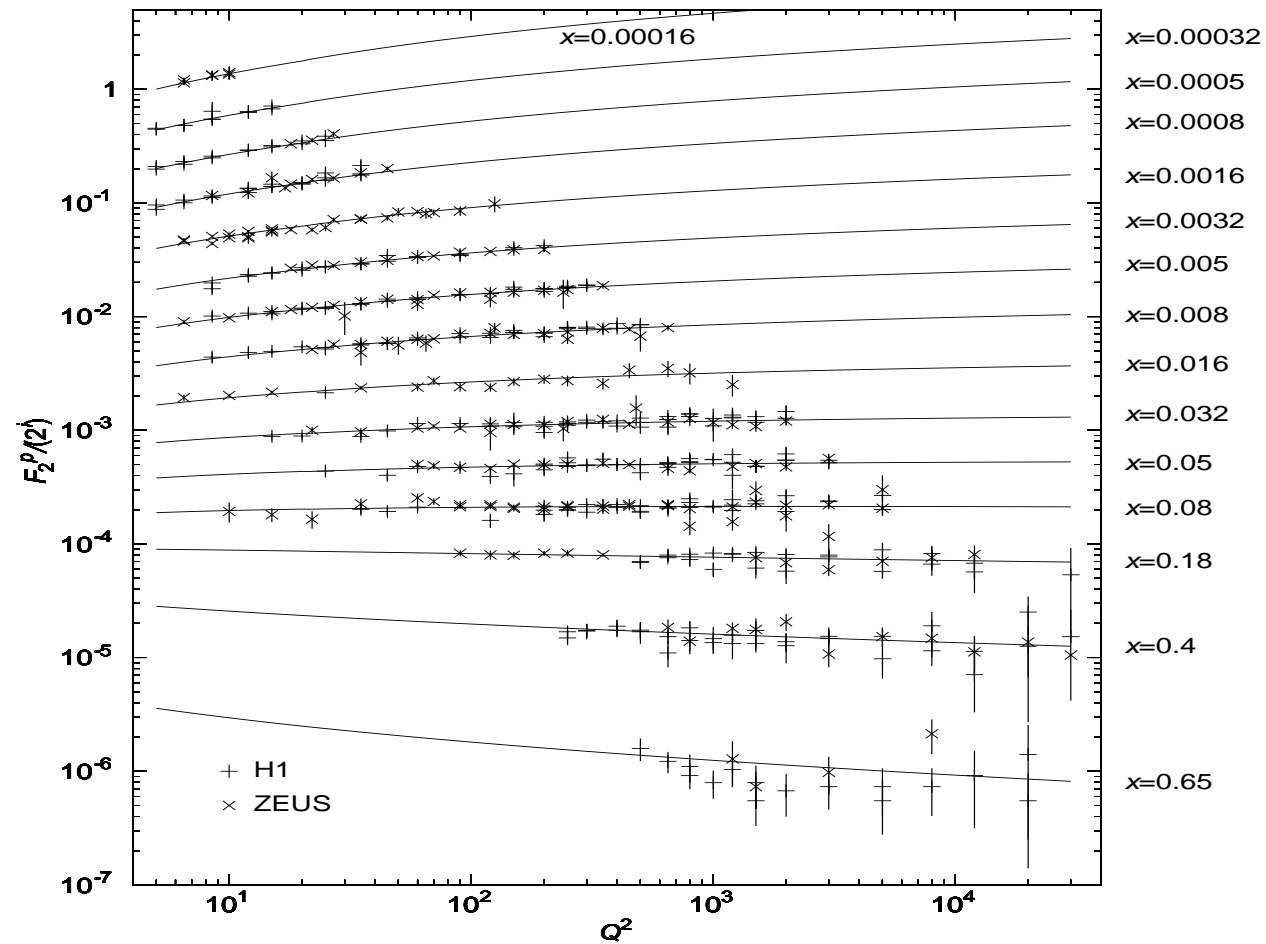
$$|\psi(z, \vec{x}_\perp, Q^2)|^2 = f(\bar{Q}_f \vec{x}_\perp)$$

$$\text{with } \bar{Q}_f^2 \equiv m_f^2 + Q^2 z^2(1 - z^2)$$

- If one neglects the quark masses, the scaling properties of $\langle \mathbf{T} \rangle_Y$ imply that σ_{γ^*p} depends only on the ratio $Q^2/Q_s^2(Y)$, rather than on Q^2 and Y separately

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■ HERA data as a function of Q^2 and x :



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Stasto, Golec-Biernat, Kwiecinski (2000)

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