

From QCD kinetics to hydrodynamics

Derek Teaney
Stony Brook University



Stony Brook University

- ▶ G. Basar, DT, Phys. Rev. C (2013)
- ▶ L. Keegan, A. Kurkela, A. Mazeliauskas, DT, JHEP (2016)
- ▶ A. Kurkela, A. Mazeliauskas, J.F. Paquet, S. Schlichting, DT: 1805.01604
- ▶ A. Kurkela, A. Mazeliauskas, J.F. Paquet, S. Schlichting, DT: 1805.00961

Two Questions:

1. How does v_2 depends on multiplicity?
2. At what multiplicity does hydro set in?

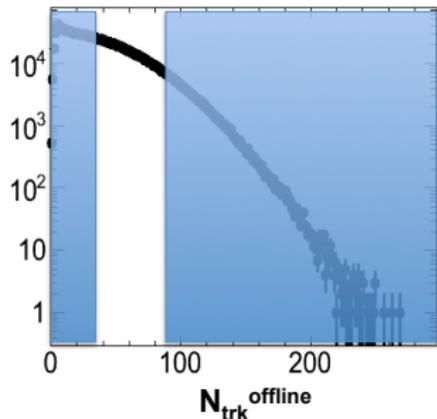
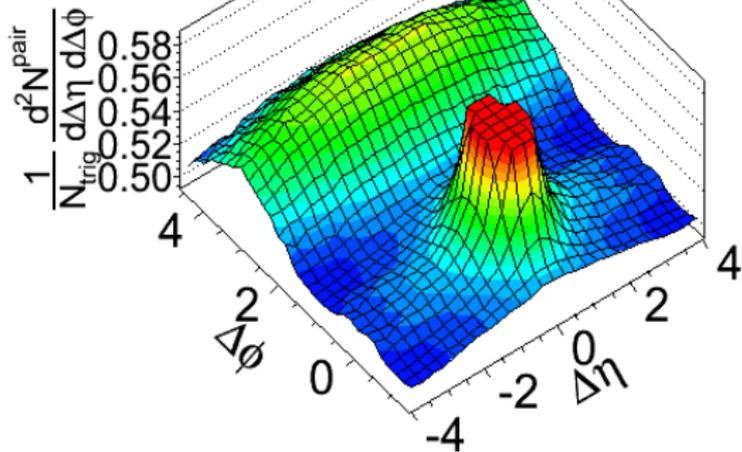
A typical pPb event

(figure G. Roland)

CMS pPb $\sqrt{s} = 5.02$ TeV $35 \leq N < 90$

$1 < p_{\text{T}}^{\text{trig}} < 2$ GeV/c

$1 < p_{\text{T}}^{\text{assoc}} < 2$ GeV/c



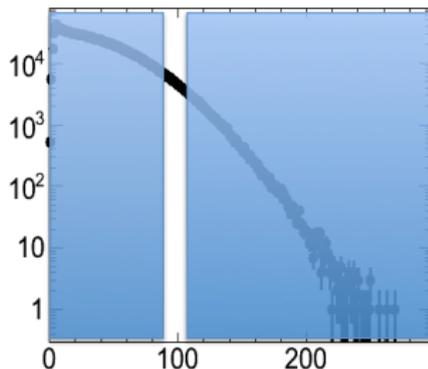
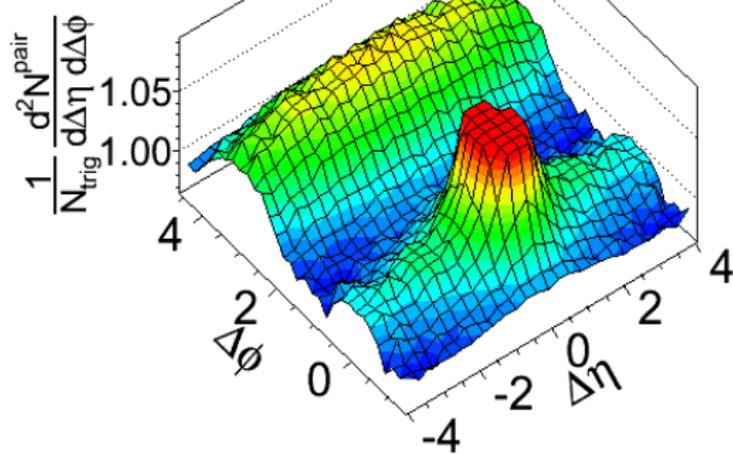
Somewhat rare event 3%

(figure G. Roland)

CMS pPb $\sqrt{s} = 5.02$ TeV $90 \leq N < 110$

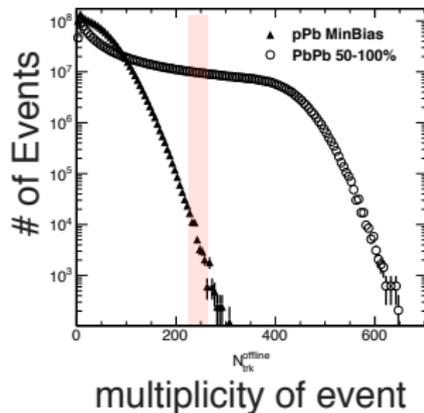
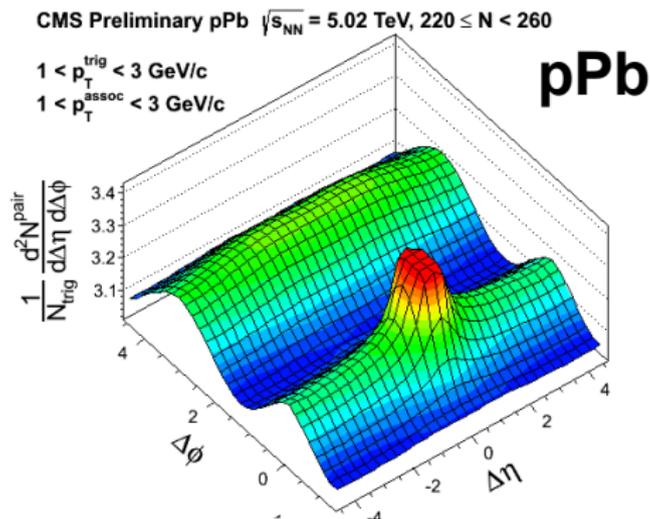
$1 < p_T^{\text{trig}} < 2$ GeV/c

$1 < p_T^{\text{assoc}} < 2$ GeV/c



Rare event – one in 20 thousand

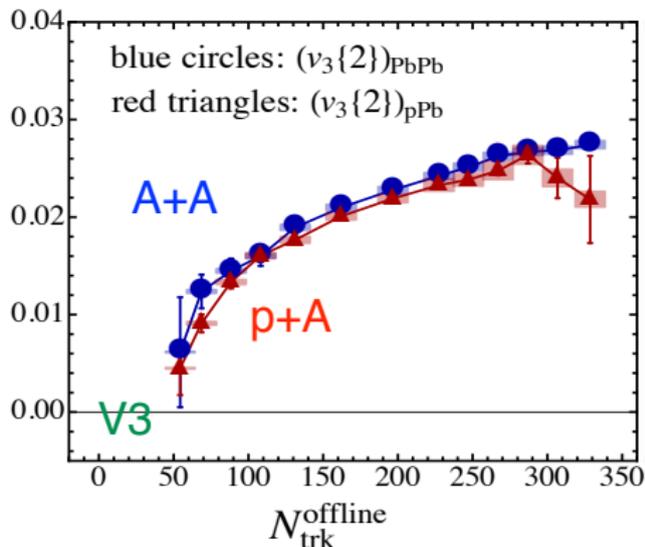
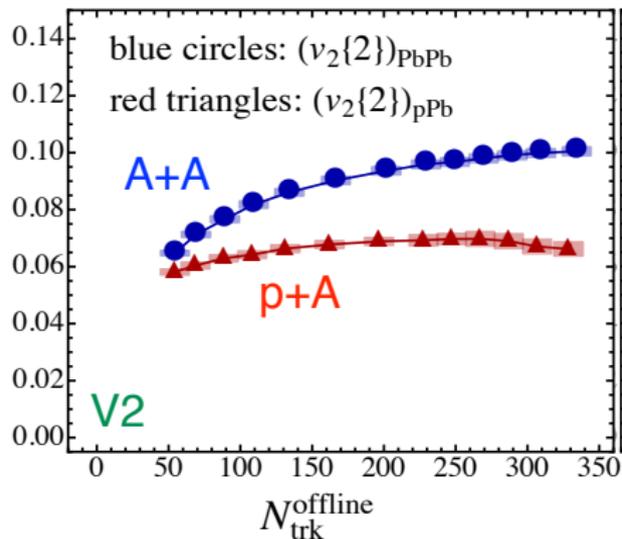
(figure G. Roland)



Want to understand this *transition*!

Can extract a v_2 and a v_3 from these data ...

(CMS Data)



The v_3 are *really the same!*

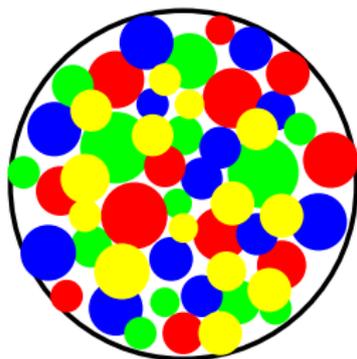
Is the mechanism for the correlation the same in pPb and PbPb?
Is hydrodynamics/kinetics the origin of pPb correlations?

Flow in pA and AA from hydro and kinetics:

$$\underbrace{v_2}_{\text{Measured E by E flow}} = \underbrace{k_2}_{\text{Response coefficient}} \times \underbrace{\epsilon_2}_{\text{geometry}}$$

1. Are the response coefficients the same?
2. Is the geometry or ϵ_2 the same? ϵ_3 ?

What is the ℓ_{mfp}/R in high multiplicity p+A and A+A?



1. Throw N_{clust} clusters in in the transverse plane:

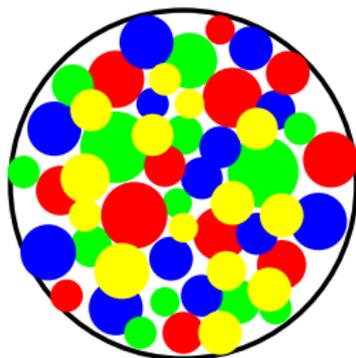
- This defines a momentum scale (the saturation momentum)

$$Q_s^2 \propto \frac{N_{\text{clust}}}{\pi R^2} \quad (R = \text{transverse size})$$

2. Assume Q_s is the only scale for the 1D expansion, $1/Q_s \ll \tau \ll R$

$$\ell_{\text{mfp}} \propto \frac{1}{Q_s} \propto \frac{1}{T_0}$$

What is the ℓ_{mfp}/R in high multiplicity p+A and A+A?



3. Assume the multiplicity is proportional to the number of clusters

$$\frac{dN}{dy} \propto N_{\text{clust}}$$

4. Find that ℓ_{mfp}/R (at early time) is fixed by dN/dy

$$\frac{\ell_{\text{mfp}}}{R} \propto \frac{1}{T_o R} \propto \frac{1}{Q_s R} \propto \frac{1}{\sqrt{dN/dy}}$$

$\frac{\ell_{\text{mfp}}}{R}$ is the same in pA and AA at fixed multiplicity!

What is the ℓ_{mfp}/R when the flow develops?

1. Need to estimate ℓ_{mfp}/R when the flow develops, $1/Q_s \ll \tau \sim R$

$$\tau \sim \frac{R}{c_s} \quad \Leftarrow \quad \text{elliptic flow develops at this time}$$

2. Using the Bjorken estimate:

$$\ell_{\text{mfp}} \propto \frac{1}{T(\tau)} \propto \frac{1}{T_o} \left(\frac{\tau}{\tau_o} \right)^{1/3}$$

Find:

$$\frac{1}{w(R)} \equiv \frac{\ell_{\text{mfp}}}{R} \Big|_{\tau \sim R} \propto \frac{1}{\left(\frac{dN}{dy} \right)^{1/3}}$$

Again $\frac{\ell_{\text{mfp}}}{R}$ is the same in pA and AA at *fixed* multiplicity

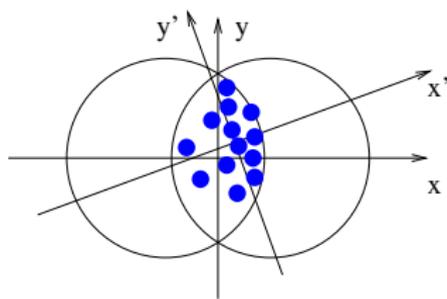
Flow in pA and AA from hydro and kinetics:

$$\underbrace{v_2}_{\text{Measured E by E flow}} = \underbrace{k_2}_{\text{Response coefficient}} \times \underbrace{\epsilon_2}_{\text{geometry}}$$

1. Are the response coefficients the same?

Yes. The ℓ_{mfp}/R is the same at fixed multiplicity

2. Is the geometry or ϵ_2 the same? ϵ_3 ?



► Throw N_{clust} sampled from the *average/smooth* geometry

► Calculate $\epsilon_2\{2\} = \sqrt{\langle \epsilon_2^2 \rangle}$

$$\langle \epsilon_2^2 \rangle_{AA} = \underbrace{\epsilon_s^2}_{\text{ave. geometry}} + \underbrace{\langle \delta \epsilon_2^2 \rangle}_{\text{fluctuations}} \quad \langle \delta \epsilon_2^2 \rangle_{AA} \equiv \frac{\langle r^4 \rangle_{AA}}{N_{\text{clust}} \langle r^2 \rangle_{AA}^2}$$

Reproduces the results of more complex Glauber models

Understanding the fluctuating geometry in p+A:

- ▶ In pA the average geometric eccentricity is zero, $\epsilon_s = 0$

$$\underbrace{\langle \epsilon_2^2 \rangle_{pA}^2}_{\text{only fluctuates}} = \underbrace{\langle \delta \epsilon_2^2 \rangle_{pA}}_{\text{only fluctuates}} \equiv \frac{\langle r^4 \rangle_{pA}}{N_{\text{clust}} \langle r^2 \rangle_{pA}^2}$$

- ▶ So to compare to pA, we scale out the average geometry out of the AA system

$$\underbrace{(v_2\{2\})_{\text{PbPb,rscl}}}_{\text{only fluctuates}} \equiv \sqrt{1 - \frac{\epsilon_s^2}{\langle \epsilon_2^2 \rangle_{AA}}} (v_2\{2\})_{\text{PbPb}}$$

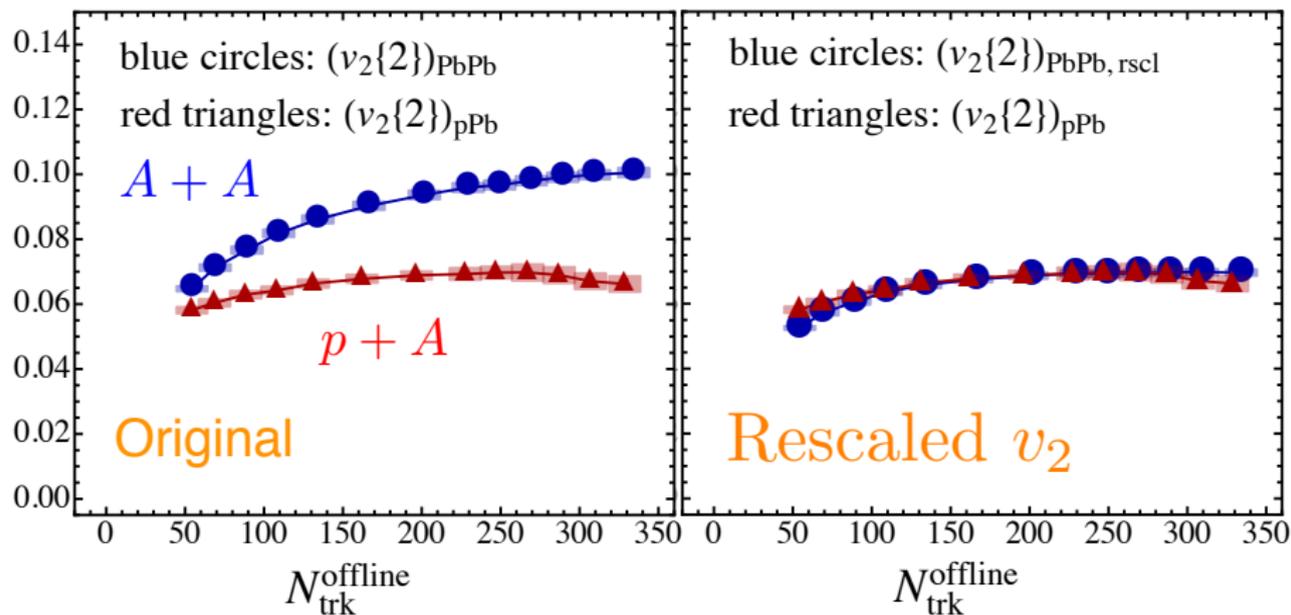
- ▶ With this rescaling we expect

$$(v_2\{2\})_{\text{PbPb,rscl}} \equiv k_2 \sqrt{\langle \delta \epsilon_2^2 \rangle_{AA}}$$
$$(v_2\{2\})_{pA} \equiv k_2 \sqrt{\langle \delta \epsilon_2^2 \rangle_{pA}}$$

Lets compare $(v_2\{2\})_{\text{PbPb,rscl}}$ to $(v_2\{2\})_{pPb} \dots$

Comparing fluctuation driven $v_2 \dots$

(CMS Data)



Fluctuation driven v_2 's are equal to 5% accuracy!
 Strong indication that v_2 is a response to geometry in $p + A$

Ratio of fluctuation v_2 's and eccentricities

- ▶ Fluctuation driven v_2

$$\frac{(v_2\{2\})_{pPb}}{(v_2\{2\})_{PbPb}} = \frac{k_2}{k_2} \sqrt{\frac{\langle \delta\epsilon_2^2 \rangle_{pA}}{\langle \delta\epsilon_2^2 \rangle_{AA}}} = \sqrt{\frac{(\langle r^4 \rangle / \langle r^2 \rangle^2)_{pA}}{(\langle r^4 \rangle / \langle r^2 \rangle^2)_{AA}}}$$

- ▶ Don't know the radial profile in pA :

But, it doesn't matter:

- ▶ Ratio of fluct-driven v_2 is determined by a root of a double ratio!
 - ▶ Very different profiles give the almost same answer

$$\sqrt{\frac{\langle \delta\epsilon_2^2 \rangle_{\text{hard-sphere}}}{\langle \delta\epsilon_2^2 \rangle_{\text{Gaussian}}}} \approx 0.85$$

Flow in pA and AA from hydro and kinetics:

$$\underbrace{v_2}_{\text{Measured E by E flow}} = \underbrace{k_2}_{\text{Response coefficient}} \times \underbrace{\epsilon_2}_{\text{geometry}}$$

1. Are the response coefficients the same? Yes.
2. Is the fluctuating geometry is the same? Yes.

v_2 scales with multiplicity as:

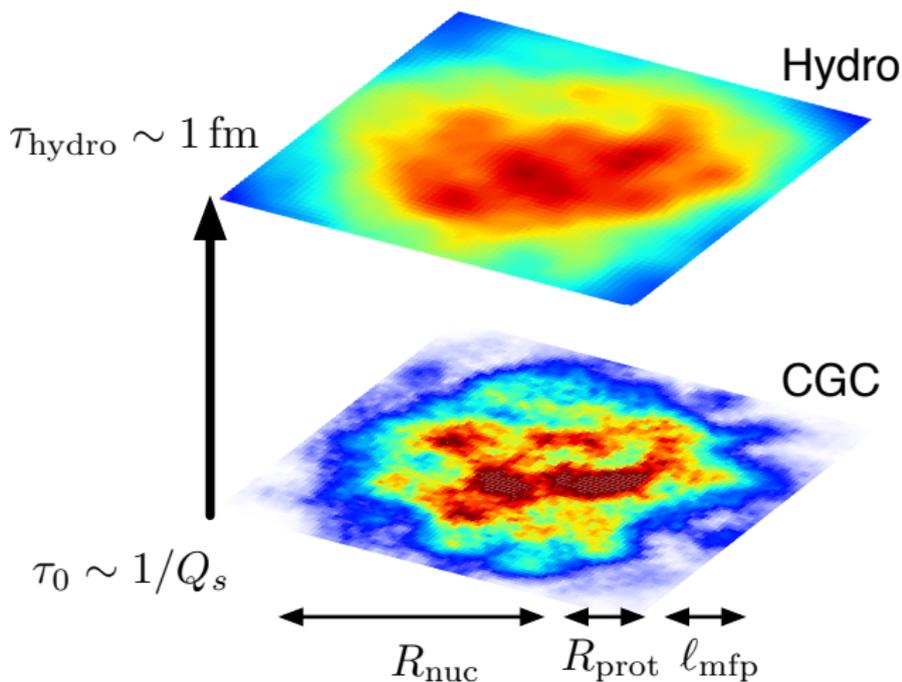
$$v_2 \propto \underbrace{\left(\frac{dN}{dy}\right)^{1/3}}_{\text{response}} \times \underbrace{\left(\frac{dN}{dy}\right)^{-1/2}}_{\text{geometry}} \propto \left(\frac{dN}{dy}\right)^{-1/6}$$

v_2 from fluctuations slowly *increases* with *decreasing* multiplicity

Two Questions:

1. How does v_2 depends on multiplicity ?
2. At what multiplicity does hydro set in?
 - ▶ Studied the approach to hydro in AA

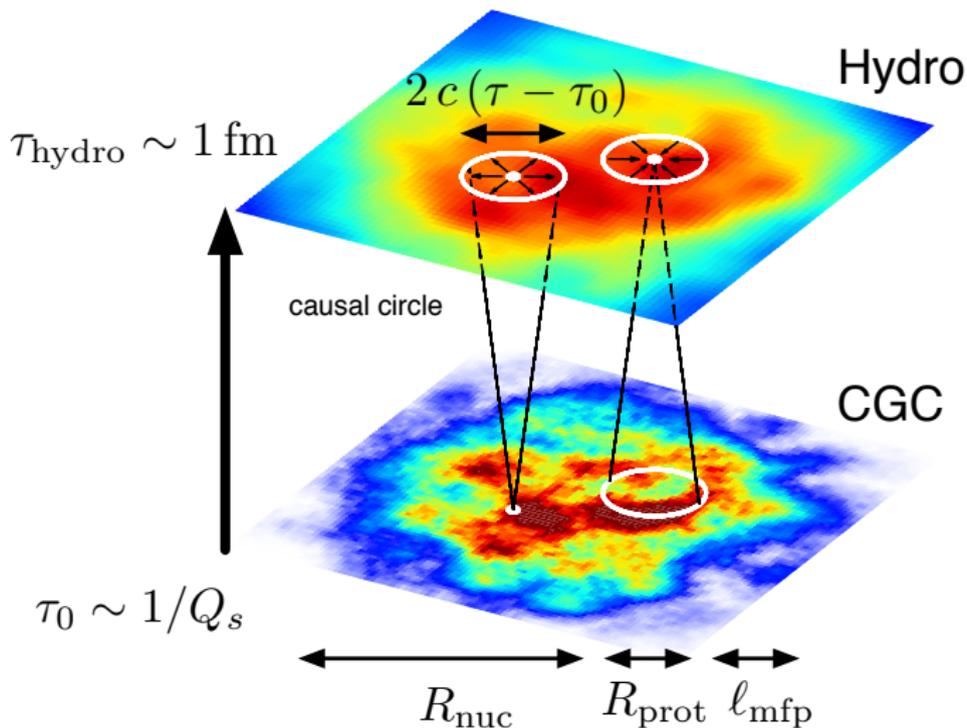
Mapping the CGC fluctuating initial conditions to hydro



Use QCD kinetic theory to map the CGC initial state to hydrodynamics

$$R_{\text{nuc}} \gg R_{\text{prot}} \sim l_{\text{mfp}} \gg 1/Q_s$$

Mapping the CGC fluctuating initial conditions to hydro

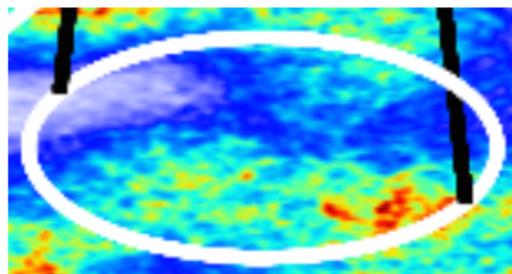


Causality limits the equilibration dynamics within a causal circle

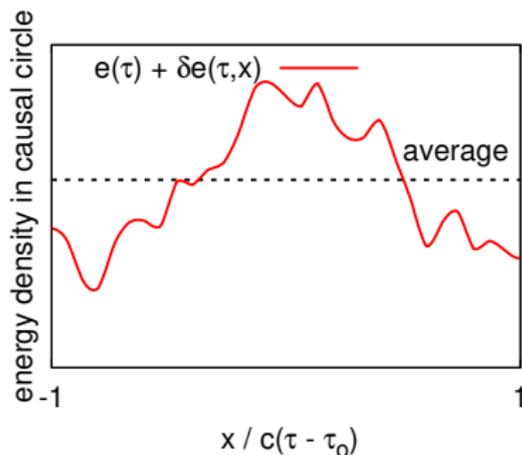
$$R_{\text{nuc}} \gg R_{\text{prot}} \sim \ell_{\text{mfp}} \sim c\tau_{\text{hydro}} \gg 1/Q_s$$

An approximation scheme for the equilibration dynamics:

look in causal circle



$$2c(\tau - \tau_0)$$



1. Determine the evolution of the average (homogeneous) background
Bottom-Up Thermalization!
2. Construct a Green function to propagate the linearized fluctuations.

$$\underbrace{\frac{\delta e(\tau, \mathbf{x})}{e(\tau)}}_{\text{final energy perturb}} = \int d^2 \mathbf{x}' G(\mathbf{x} - \mathbf{x}'; \tau, \tau_0) \underbrace{\frac{\delta e(\tau_0, \mathbf{x}')}{e(\tau_0)}}_{\text{initial energy perturb}}$$

final energy perturb

initial energy perturb

How to compute the background and perturbations:

$$\partial_\tau f + \frac{\mathbf{p}}{|\mathbf{p}|} \cdot \nabla f - \underbrace{\frac{p_z}{\tau} \partial_{p_z} f}_{\text{Bjorken expansion}} = -\underbrace{\mathcal{C}_{2 \leftrightarrow 2}[f]}_{\text{diagram}} - \underbrace{\mathcal{C}_{1 \leftrightarrow 2}[f]}_{\text{diagram}},$$

Gluon distribution function for background and perturbations

$$f = \underbrace{\bar{f}_{\mathbf{p}}}_{\text{uniform background}} + \underbrace{\delta f_{\mathbf{k}_\perp, \mathbf{p}} e^{i\mathbf{k}_\perp \cdot \mathbf{x}_\perp}}_{\text{transverse perturbations}}.$$

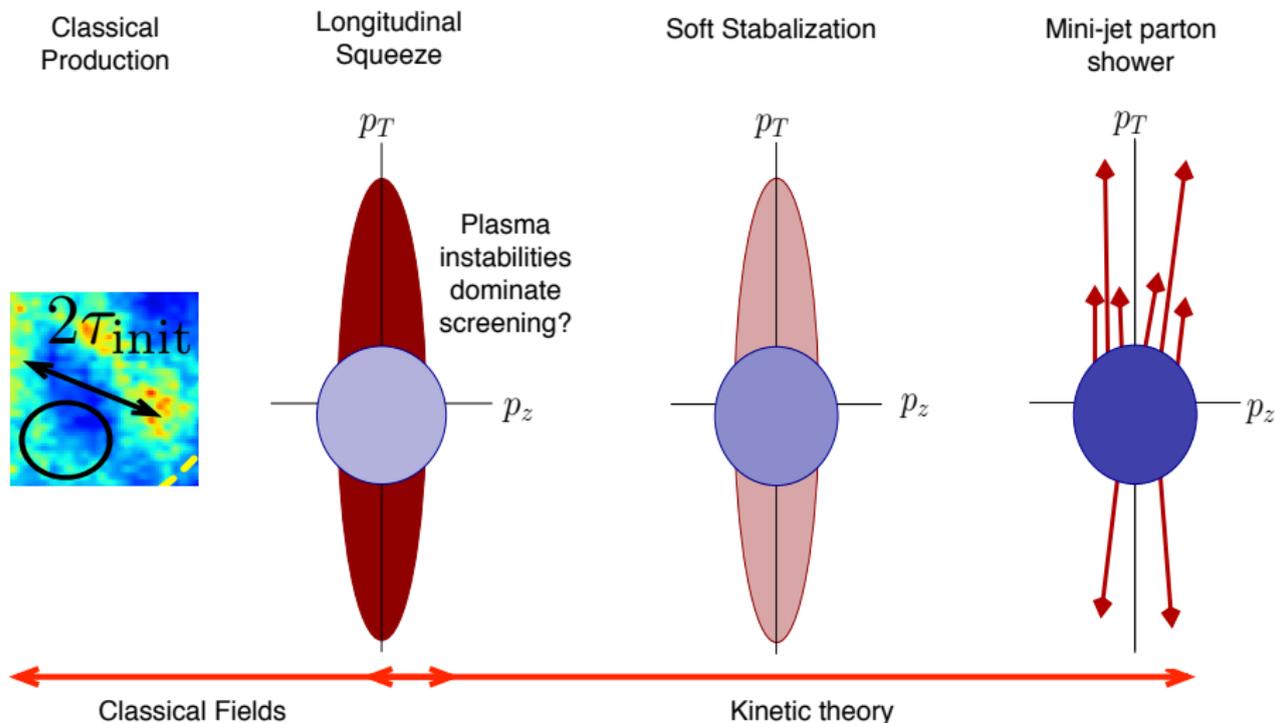
$$(\partial_\tau - \frac{p_z}{\tau} \partial_{p_z}) \bar{f}_{\mathbf{p}} = -\mathcal{C}[\bar{f}] \quad \text{background}$$

$$(\partial_\tau - \frac{p_z}{\tau} \partial_{p_z} + \frac{i\mathbf{p}_\perp \cdot \mathbf{k}_\perp}{p}) \delta f_{\mathbf{k}_\perp, \mathbf{p}} = -\delta\mathcal{C}[\bar{f}, \delta f] \quad \text{perturbation}$$

We will discuss the background and perturbations separately

The background and “bottom-up” thermalization

Baier, Mueller, Schiff, Son

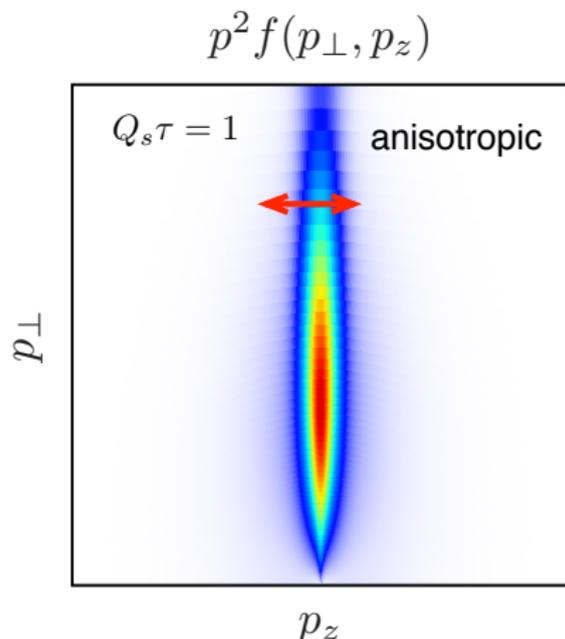


Reach a thermal state in $\tau_{\text{hydro}} \sim 1/(\alpha_s^{13/5} Q_s)$

A numerical realization of bottom-up

- Builds upon the first numerical realization

Kurkela, Zhu PRL (2015)



Initialization:

- Partons are initialized with:

$$\langle p_\perp^2 \rangle \sim Q_s^2 \quad \langle p_z^2 \rangle \simeq 0$$

- Take a coupling of $\alpha_s = 0.3$

$$\lambda \equiv \underbrace{4\pi\alpha_s N_c}_{\text{theorists version of } \alpha_s = 0.3} = 10$$

theorists version of $\alpha_s = 0.3$

corresponding to

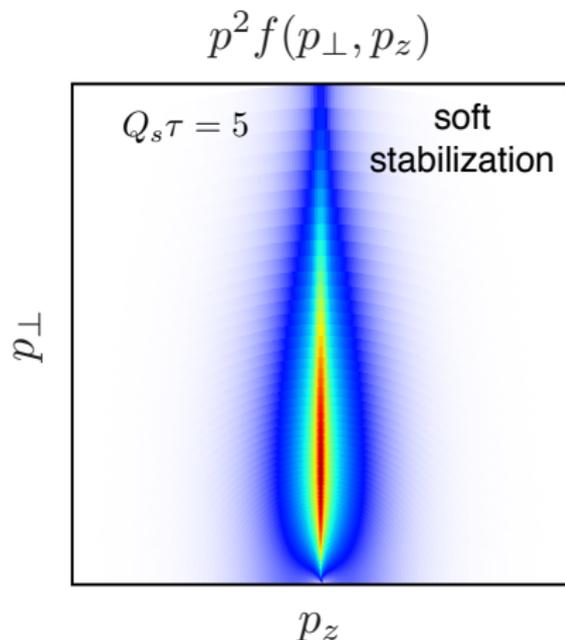
$$\frac{\eta}{s} = 0.6 = \frac{7.5}{4\pi}$$

We see “Bottom-Up” in the computer code.

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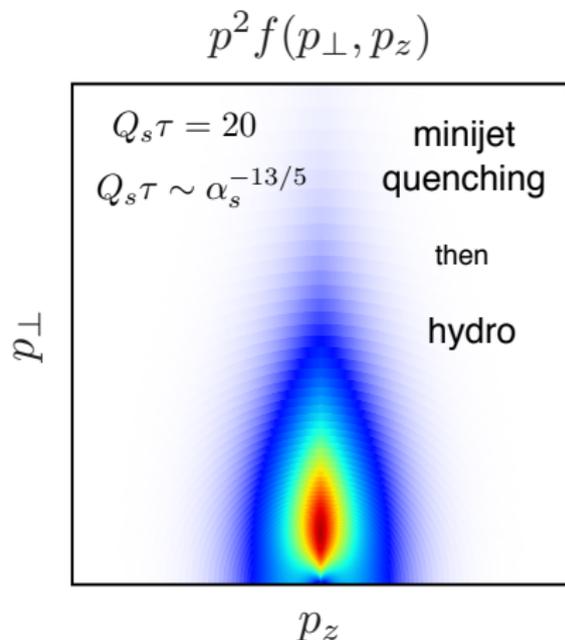
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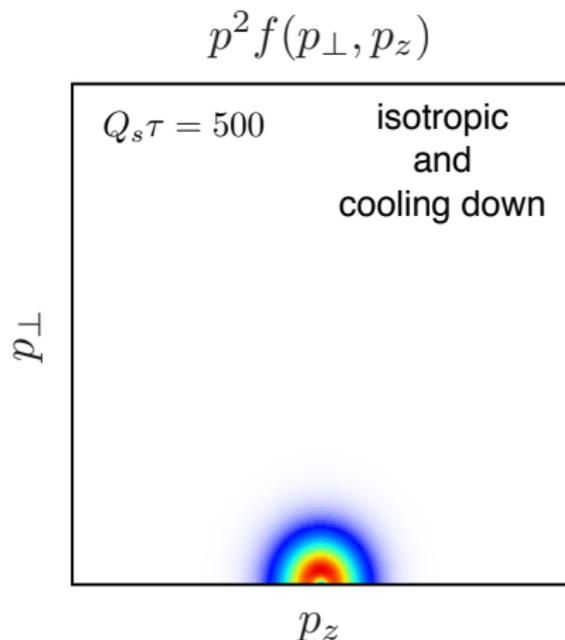
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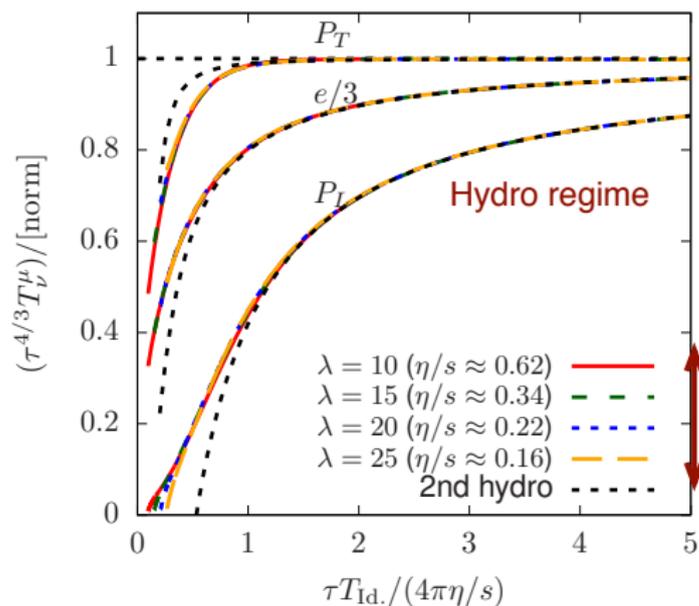
theorists version of $\alpha_s = 0.3$

corresponding to

$$\frac{\eta}{s} = 0.6 = \frac{7.5}{4\pi}$$

We see “Bottom-Up” in the computer code.

When does the stress tensor approach 2nd order hydrodynamics?



Different values of coupling
give different η/s

In terms of η/s , all couplings
thermalize at same scaled time
Keegan, Kurkela, Romatschke, Schee, Zhu

Gives a basis for interpolating
from weak coupling results to
stronger coupling

Measure time in a physical relaxation time given by $\tau_R \equiv 4\pi\eta/s T_{\text{eff}}(\tau)$:

$$\frac{\tau}{\tau_R} \equiv \frac{\tau T_{\text{eff}}(\tau)}{4\pi\eta/s} \quad \text{with} \quad T_{\text{Id}} \equiv T_{\text{eff}}(\tau) \equiv \Lambda_T / (\Lambda_T \tau)^{1/3}$$

Can start hydro when $\tau T_{\text{eff}}(\tau) / 4\pi\eta/s \sim 1$

Translating earliest hydro starting time into physical units:

- ▶ At late times the dynamics is ideal hydro: $T_{\text{eff}}(\tau) = \Lambda_T / (\Lambda_T \tau)^{1/3}$
- ▶ Hydro fits to multiplicity give:

$$\langle \tau s \rangle |_{\tau=1.2 \text{ fm}} = \underbrace{4.1 \text{ GeV}^2}_{\text{highly constrained by } \frac{dN}{dy}!} \propto \Lambda_T^2$$

- ▶ The estimate for τ_{hydro} :

$$\frac{\tau_{\text{hydro}} T_{\text{eff}}(\tau_{\text{hydro}})}{4\pi(\eta/s)} = 1$$

Find that hydrodynamics is applicable for times later than:

$$\tau_{\text{hydro}} \approx 1.1 \text{ fm} \left(\frac{4\pi(\eta/s)}{2} \right)^{\frac{3}{2}} \left(\frac{4.1 \text{ GeV}}{\langle \tau s \rangle} \right)^{-1/2} \left(\frac{\nu_{\text{eff}}}{40} \right)^{1/2}$$

Estimate for the applicability of hydro for system of radius R :

- ▶ For a hydrodynamic regime the system should equilibrate by $\tau \sim R$:

$$R/\ell_{\text{mfp}} \text{ at time } R: \quad w(R) \equiv \left. \frac{\tau T_{\text{id}}(\tau)}{4\pi(\eta/s)} \right|_{\tau=R} > 1$$

- ▶ Since $T \sim (dN/dy)/R^{1/3}$, then $w(R) \propto RT$ is independent of R .

$$w(R) = \left(\frac{4\pi(\eta/s)}{2} \right)^{-1} \left(\frac{dN_{\text{ch}}/d\eta}{63} \right)^{1/3} > 1$$

- We used the EOS to relate dS/dy to measured dN_{ch}/dy

- ▶ Corrections to ideal hydro response coefficients come in powers of $w(R)^{-1}$.

$$k_2(w) = k_2^{\text{ideal}} \left(1 - \underbrace{\frac{C_1}{w(R)}}_{\text{first order hydro}} - \underbrace{\frac{C_2}{w(R)^2}}_{\text{second order}} + \dots \right)$$

Two Questions with Answers:

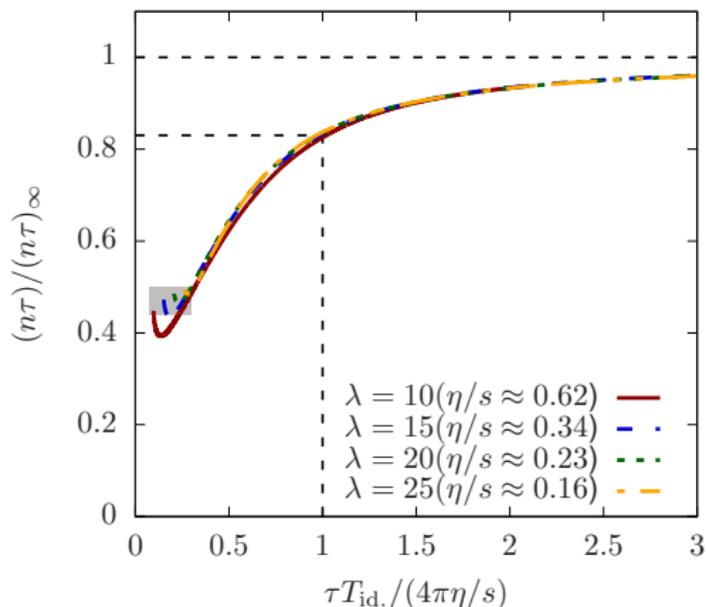
1. How does v_2 in small systems depends on multiplicity ?

$$v_2 \propto \left(\frac{dN}{dy} \right)^{-1/6}$$

2. At what multiplicity does hydro set in?

$$\left(\frac{dN}{dy} \right)_{\text{hydro}} = 63 \left(\frac{4\pi(\eta/s)}{2} \right)^3$$

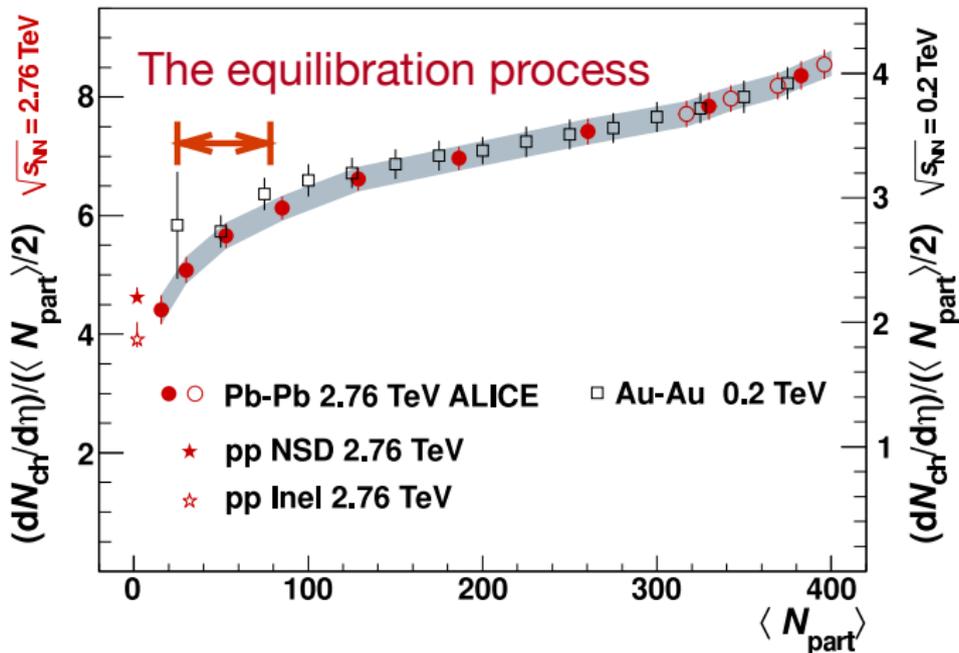
Changes during the equilibration process: increase in multiplicity



$$n\tau = \frac{1}{A} \frac{dN}{dy}$$

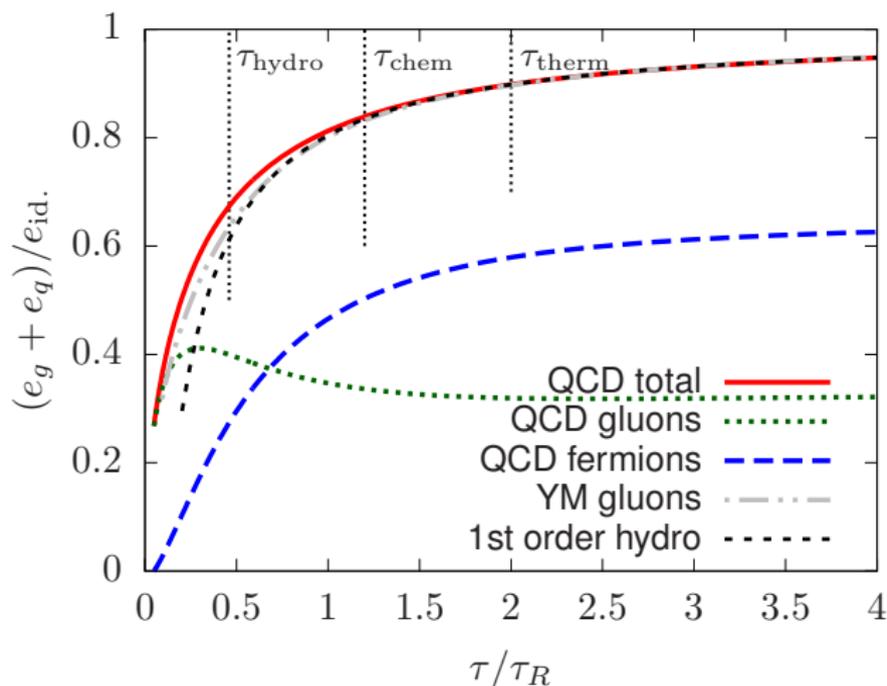
The final gluon multiplicity is 2.5 times the initial gluon multiplicity independent of the coupling or η/s !

Multiplicity increase in peripheral collisions:



Changes during the equilibration process: change in chemistry

► Kurkela and Mazeliauskas:1811.03040

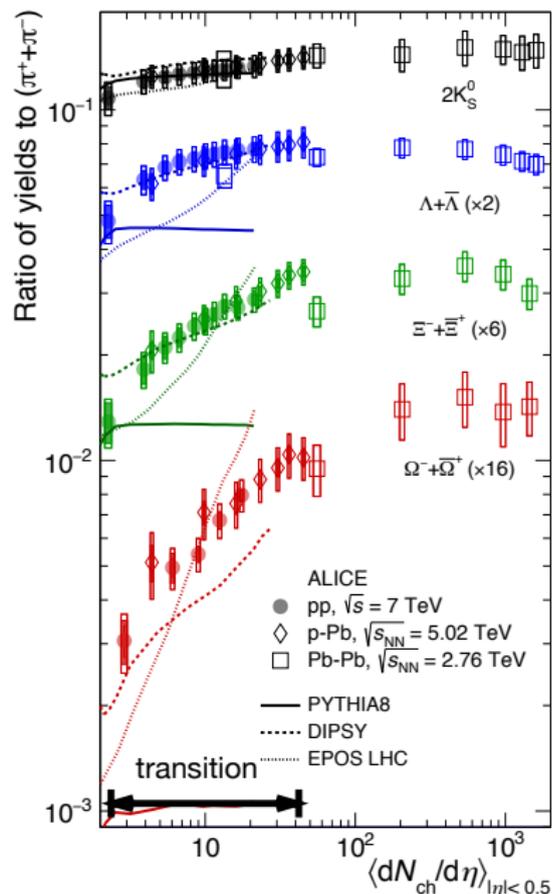


During the first one or two relaxation times the energy density becomes more than half quarks

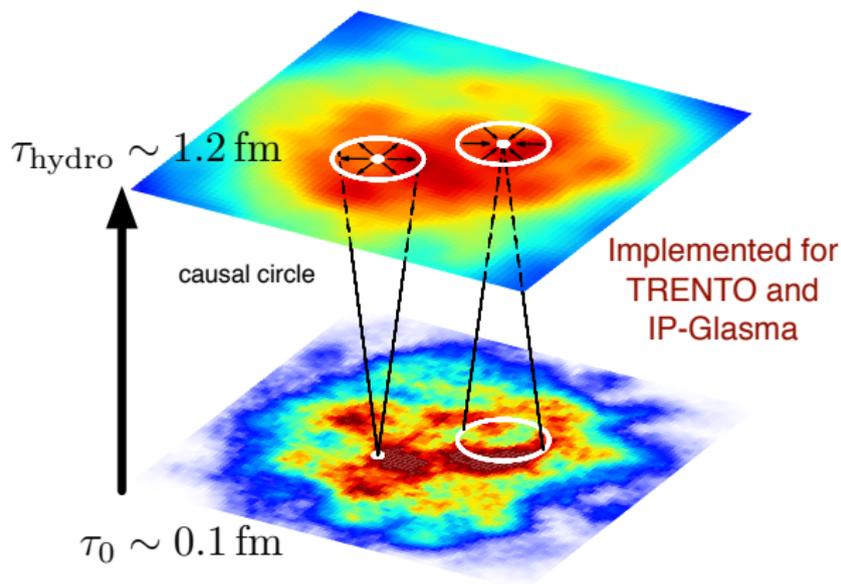
Chemical changes:

Strangeness enhancement

Alice: 1606.07424



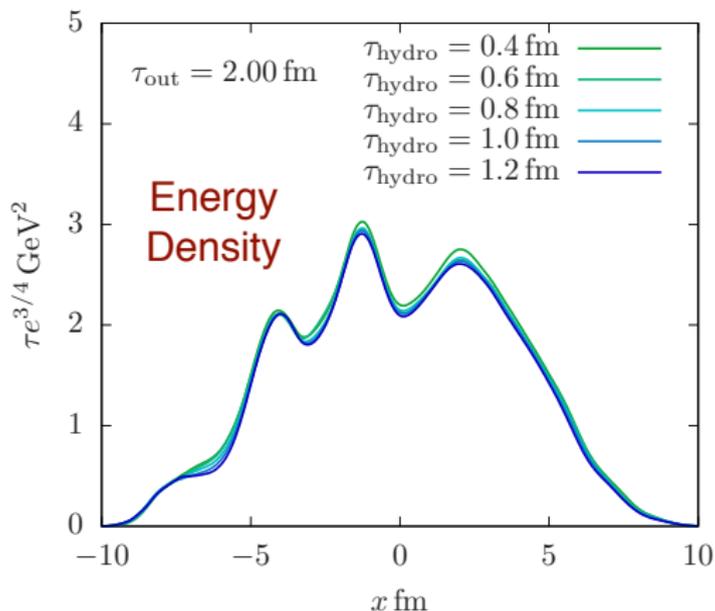
A practical algorithm for pre-hydro evolution in A+A:



- (i) For each point, average the energy in causal circle:
 - ★ Determine τ_{hydro} in units of the scaled time, $\tau_{\text{hydro}}/\tau_R$
 - ★ Evolve the background in scaled time.
- (ii) Propagate background and perturbations in scaled time
 - ★ Sometimes need to regulate the response

Do hydro results depend on τ_{hydro} ?

1. Implementation in TRENTO. $\eta/s = 2/4\pi$. Central LHC.
2. Kinetics runs from $\tau_0 = 0.1$ up to τ_{hydro} , then hydro up to τ_{out} .

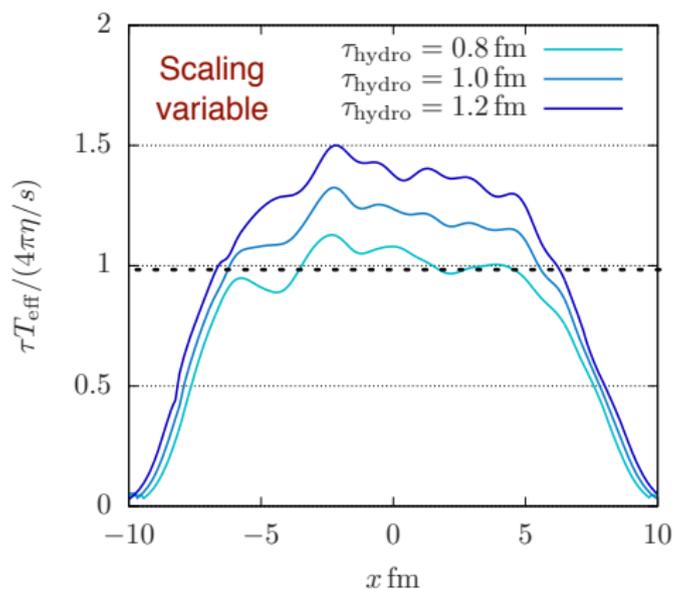


Remarkably insensitive to τ_{hydro} as we want !

Are the constitutive relations are satisfied at late times?

- For times sufficiently late times Navier-Stokes should be valid:

$$\pi^{\mu\nu} = \underbrace{-\eta\sigma^{\mu\nu}}_{\text{navier stokes}} \quad \text{for} \quad \frac{\tau T_{\text{eff}}}{4\pi\eta/s} > 1$$

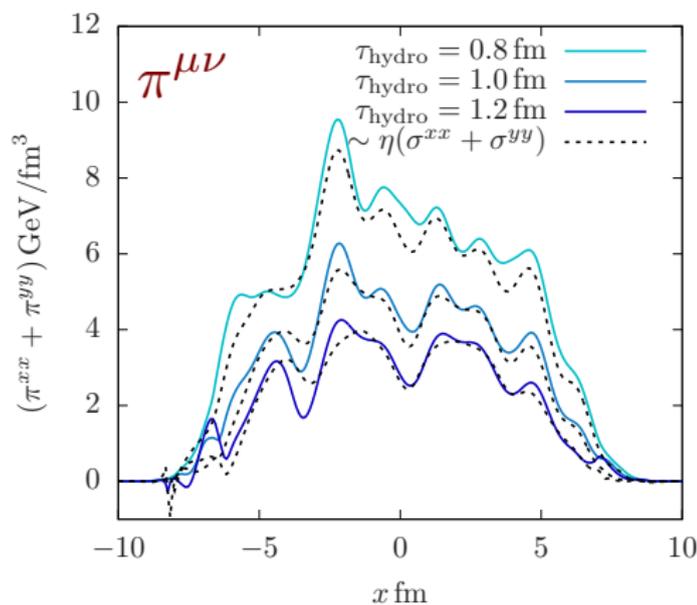


Expect the cells near the line to be equilibrated and obey constitutive equations

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$$\pi^{\mu\nu} = \underbrace{-\eta\sigma^{\mu\nu}}_{\text{navier stokes}} \quad \text{for} \quad \frac{\tau T_{\text{eff}}}{4\pi\eta/s} > 1$$



Black lines
are navier
stokes.

Color lines
are kinetics

Constitutive
relations are satisfied!

Summary:

- ▶ Want to see a transition to flow:
 - ▶ Counter productive to only study flow, and to ignore non-flow
 - ▶ Can become impossible to define them separately
- ▶ At what multiplicity does hydro set in?

$$\left(\frac{dN}{dy}\right)_{\text{hydro}} = 63 \left(\frac{4\pi(\eta/s)}{2}\right)^3$$

- ▶ Equilibration leaves fingerprints on multiplicities and strangeness
- ▶ For the AA system we have constructed a useful tool for pre-flow.