

The shear viscosity of parton matter under two-body scatterings

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Outline

- 1) Motivation
- 2) Comparison of η and η/s from different methods
- 3) Application to parton matter in the AMPT model
- 4) Conclusions

Mainly based on
Noah MacKay & ZWL,
Eur Phys J C 82, 918 (2022)



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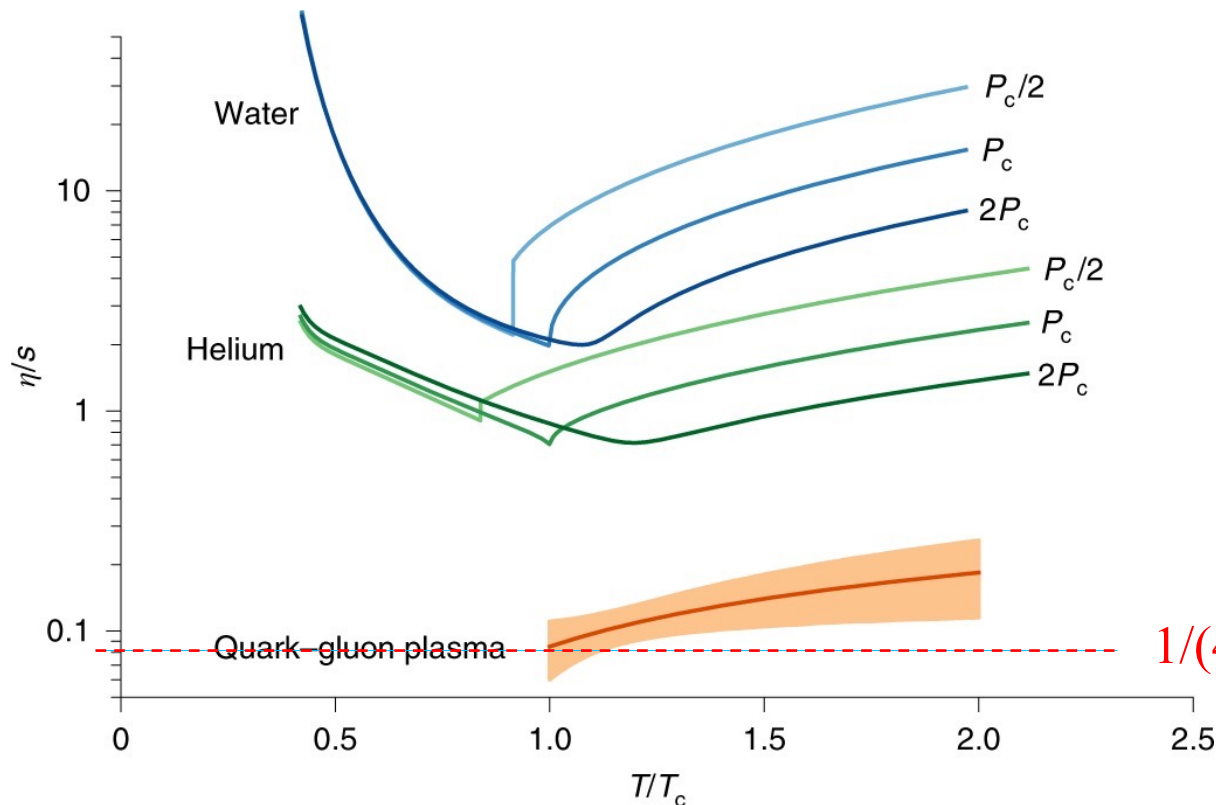
1) Motivation

Shear viscosity η is an important property of the quark–gluon plasma:

$$T^{\mu\nu} = eu^\mu u^\nu - (p + \Pi)\Delta^{\mu\nu} + \pi^{\mu\nu}$$

η or η/s : is an input function to viscous hydrodynamics;
is generated by interactions in kinetic theory: relation?

η/s can be extracted from data/model comparisons:



Bernhard, Moreland & Bass,
Nature Phys (2019)

1/(4π) bound from AdS/CFT:

Kovtun, Son & Starinets,
Phys Rev Lett (2005)

Isotropic versus forward-angle scatterings

For parton 2-to-2 elastic scatterings:

- Isotropic scattering: $\frac{d\sigma}{d\Omega} = \text{constant} = \frac{\sigma}{4\pi}$

- Forward-angle scattering:

As the example, we take the parton cross section used in AMPT/ZPC/MPC:

$$\frac{d\sigma}{d\hat{t}} = \frac{9\pi\alpha_s^2}{2} (1 + a) \frac{1}{(\hat{t} - \mu^2)^2}$$

$a \equiv \frac{\mu^2}{\hat{s}}$ is added to obtain a \hat{s} -independent cross section: $\sigma = \frac{9\pi\alpha_s^2}{2\mu^2}$

This is based on the pQCD gg-gg cross section:

$$\frac{d\sigma}{d\hat{t}} \propto 3 \left[-\frac{\hat{t}\hat{u}}{\hat{s}^2} - \frac{\hat{s}\hat{u}}{\hat{t}^2} - \frac{\hat{s}\hat{t}}{\hat{u}^2} \right] + \text{screening mass } \mu$$

Isotropic versus forward-angle scatterings

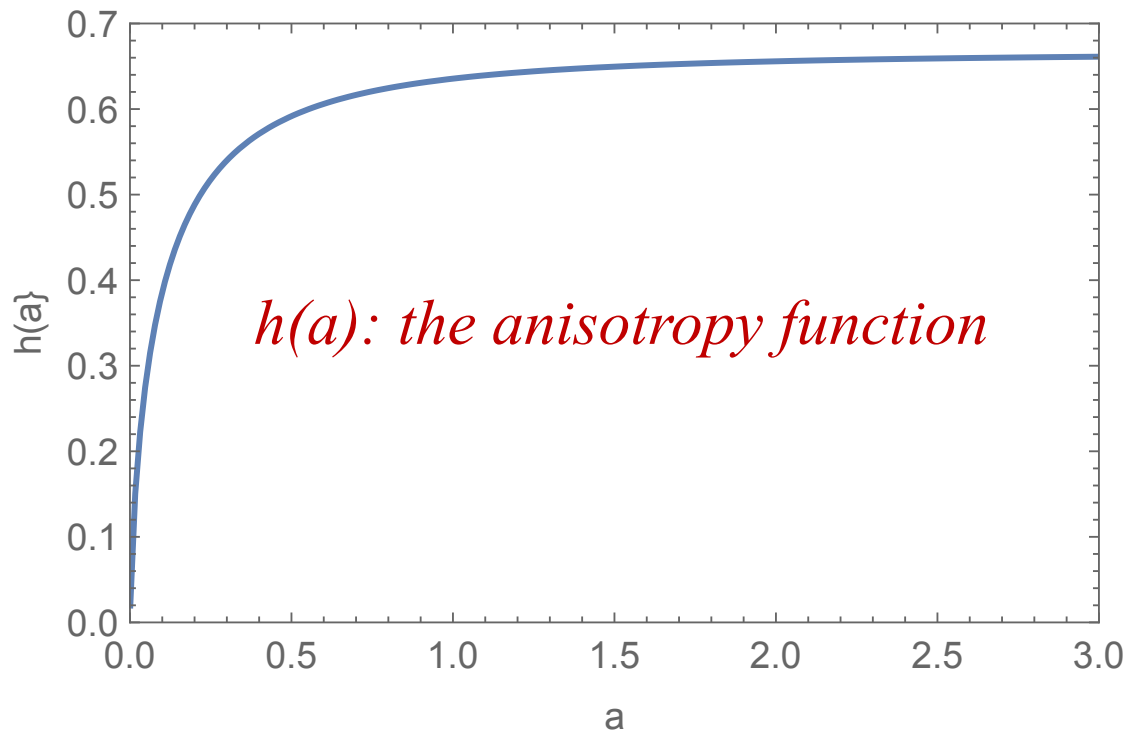
Transport cross section σ_{tr} often appears in shear viscosity expressions:

$$\sigma_{tr} \equiv \int d\sigma \sin^2 \theta_{cm}$$

θ_{cm} : scattering angle
in the 2-parton CM frame

Molnar & Gyulassy, Nucl Phys A (2002)

- Isotropic scattering: $\sigma_{tr} = \frac{2}{3} \sigma$
- Forward-angle scattering: $\sigma_{tr} = 4a(1+a) \left[(1+2a) \ln \left(1 + \frac{1}{a} \right) - 2 \right] \sigma$
 $\equiv h(a) \sigma$



$$a \equiv \frac{\mu^2}{\hat{s}}$$

$\ll 1$ very forward,
 $h(a) \rightarrow 0$: scatterings
are less effective

$\gg 1$ \sim isotropic, $h(a) \rightarrow 2/3$

Isotropic versus forward-angle scatterings

Thermal average:

even if σ is a constant, σ_{tr} is not since it depends $a \equiv \frac{\mu^2}{\hat{s}}$.

For a parton matter in thermal equilibrium at temperature T ,

the thermal average (for Boltzmann statistics) is

Kolb & Raby, Phys Rev D (1983)

$$\langle \sigma_{tr} \rangle = \frac{\sigma}{32} \int_0^\infty du [u^4 K_1(u) + 2u^3 K_2(u)] h\left(\frac{w^2}{u^2}\right)$$
$$\equiv \sigma h_0(w)$$

K_n : Bessel functions

$$w \equiv \frac{\mu}{T}, u \equiv \frac{\sqrt{\hat{s}}}{T}$$

$h_0(w)$ is just an average of the anisotropy function $h(a)$,

$$h_0(w) \rightarrow 0 \quad \text{for } w \ll 1$$

$$h_0(w) \rightarrow 2/3 \quad \text{for } w \gg 1$$

2) Comparison of η and η/s from different methods

Analytical:

- Israel–Stewart (IS) method: $\eta^{IS} = \frac{6T}{5\sigma}$ for isotropic scatt.
Huovinen & Molnar, Phys Rev C (2009)
- Navier–Stokes (NS) method: $\eta^{NS} \approx 1.2654 \frac{T}{\sigma}$ for isotropic scatt.
de Groot, van Leeuwen & Weert book (1980)
- Relaxation time approximation (RTA) & modified version (RTA*):
 $\eta^{RTA} = \frac{4T}{5\sigma}$ $\eta^{RTA*} = \frac{6T}{5\sigma}$ for isotropic scatt.
Anderson & Witting, Physica (1974) Plumari, Puglisi, Scardina & Greco, Phys Rev C (2012)
- Chapman–Enskog (CE) method: $\eta^{CE} = \frac{T \gamma_0^2}{10 c_{00}}$, ... Wiranata & Prakash, Phys Rev C (2012)

Numerical:

- Green–Kubo relation: $\eta = \frac{V}{T} \int_0^\infty dt \langle \bar{\pi}^{xy}(t+t') \bar{\pi}^{xy}(t') \rangle = \frac{4}{15} \varepsilon \tau$

τ : relaxation time extracted from correlation $\langle \dots \rangle(t)$

2) Comparison of η and η/s from different methods

Analytical:

- Israel–Stewart (IS) method: $\eta^{IS} = \frac{4T}{5\langle\sigma_{tr}\rangle} = \frac{4T}{5\sigma h_0(w)}$ ↑
Huovinen & Molnar,
Phys Rev C (2009)
- Navier–Stokes (NS) method: $\eta^{NS} \approx 0.8436 \frac{T}{\langle\sigma_{tr}\rangle}$ ↓
**generalized to
anisotropic scatt.
using σ_{tr}
instead of σ**
- Relaxation time approximation (modified version RTA*):
$$\eta^{RTA^*} = \frac{4T}{5\langle\sigma_{tr}v_{rel}\rangle}$$
 ↓
Plumari, Puglisi, Scardina & Greco, Phys Rev C (2012)
- Chapman–Enskog (CE) method: $\eta^{CE} = \frac{T \gamma_0^2}{10 c_{00}}, \dots$ Wiranata & Prakash,
Phys Rev C (2012)

Numerical:

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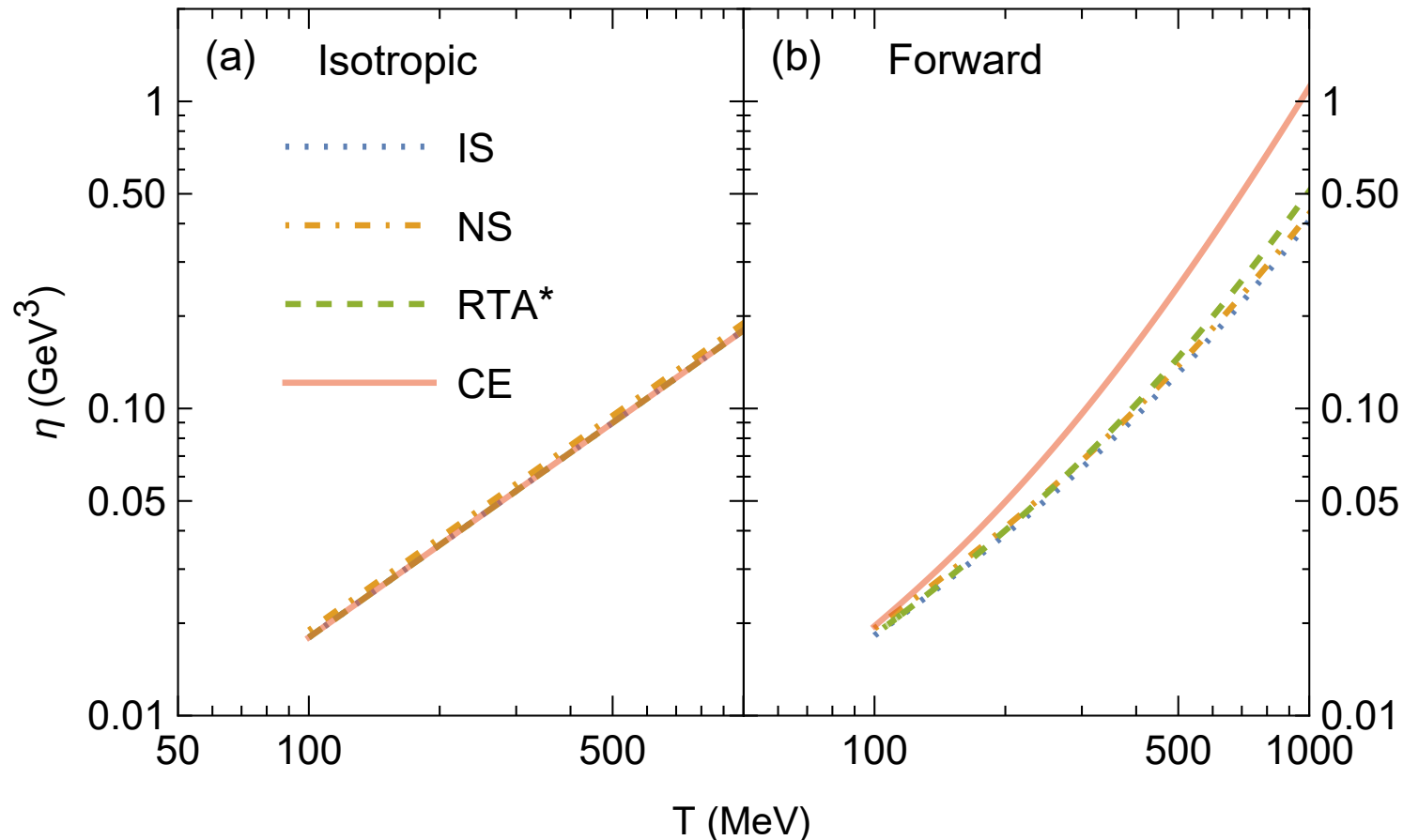
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2) Comparison of η and η/s from different methods

Analytical results of η

MacKay & ZWL, Eur Phys J C (2022)

for massless gluons & $\sigma=2.6$ mb (or $\mu\sim 0.7$ GeV at $\alpha_s \approx 0.47$):



- For isotropic scatterings:
 $IS=RTA^*=CE$
 $\approx NS$ ($\sim 5\%$ higher)

- For forward scatterings:
 $IS \approx RTA^* \approx NS < CE$ *mostly*
 $T \ll \mu \rightarrow$ almost isotropic

2) Comparison of η and η/s from different methods

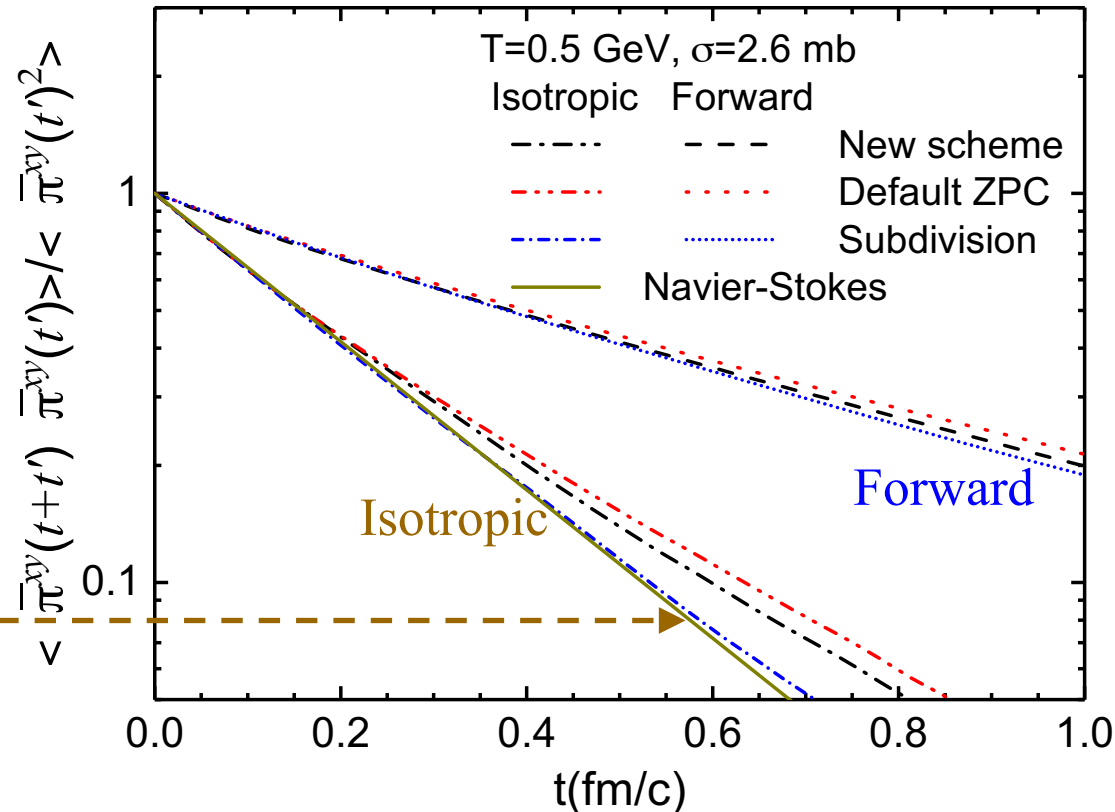
Q: which analytical result of η is accurate?

A: compare with numerical results from Green-Kubo:

$$\eta = \frac{V}{T} \int_0^\infty dt \langle \bar{\pi}^{xy}(t+t') \bar{\pi}^{xy}(t') \rangle = \frac{4}{15} \varepsilon \tau$$

With ZPC parton cascade, we have calculated η of gluons in a box with the Green-Kubo relation.

Subdivision (*with factor $l=10^6$*): results should be accurate (*no causality violation from cascade*), agree well with NS expectation for isotropic scatterings.

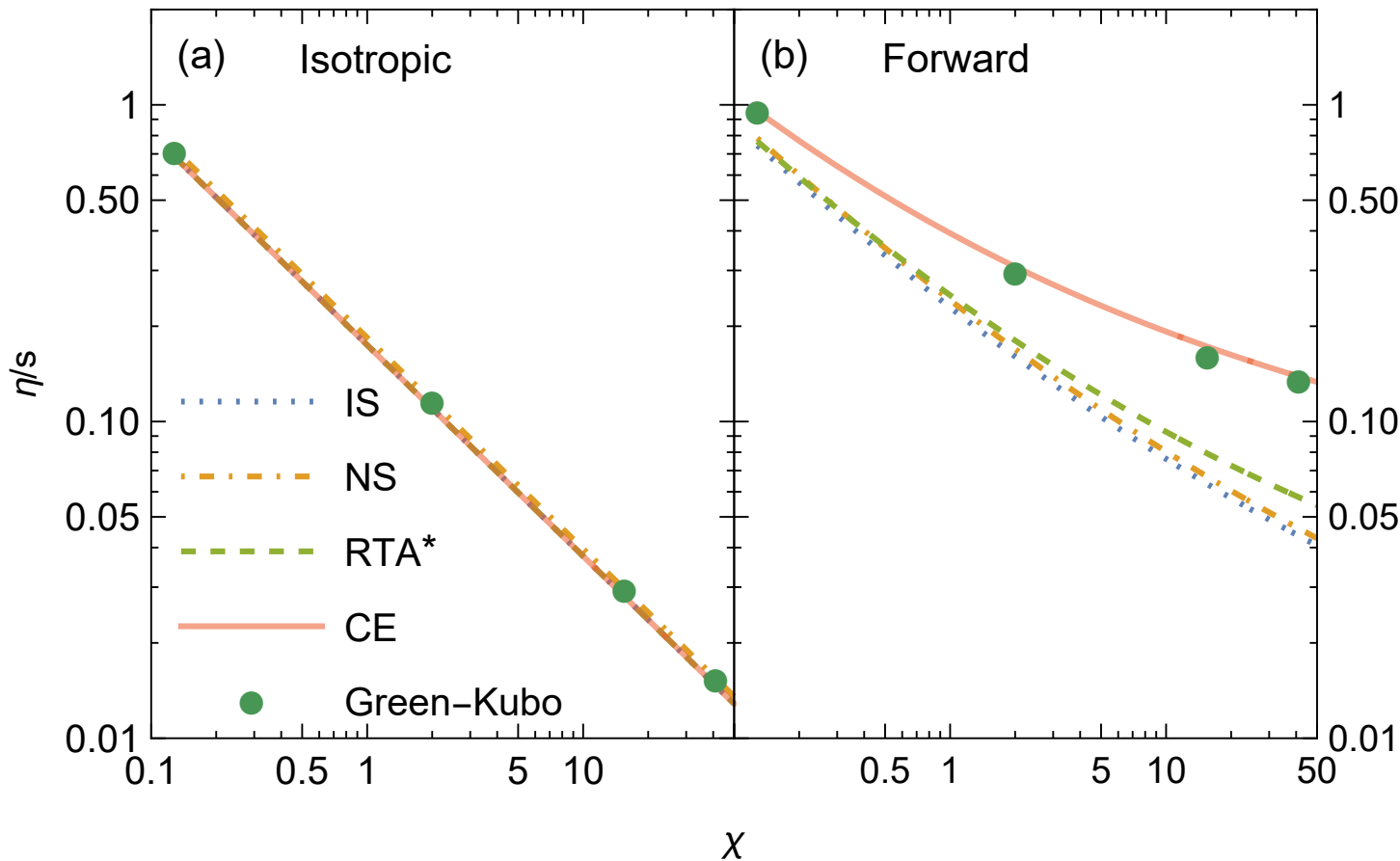


Zhao, Ma, Ma & ZWL, Phys Rev C (2020)

2) Comparison of η and η/s from different methods

Compare 4 analytical methods
with Green-Kubo results for η/s versus χ :

MacKay & ZWL, Eur Phys J C (2022)



χ (opacity parameter)
 \equiv interaction radius/mfp

$$\chi = \sqrt{\frac{\sigma}{\pi}} / \lambda = n \sqrt{\frac{\sigma^3}{\pi}}$$

Zhang, Gyulassy & Pang,
Phys Rev C (1998)

- For isotropic scatterings:
all methods agree well.

- For anisotropic scatterings:
CE agrees well with Green-Kubo;
but the other analytical methods
are not accurate.

2) Comparison of η and η/s from different methods

The fact that Green-Kubo agrees with CE (but not with RTA*) has been shown in Plumari, Puglisi, Scardina & Greco, Phys Rev C (2012) despite two typos in the eta formulae.

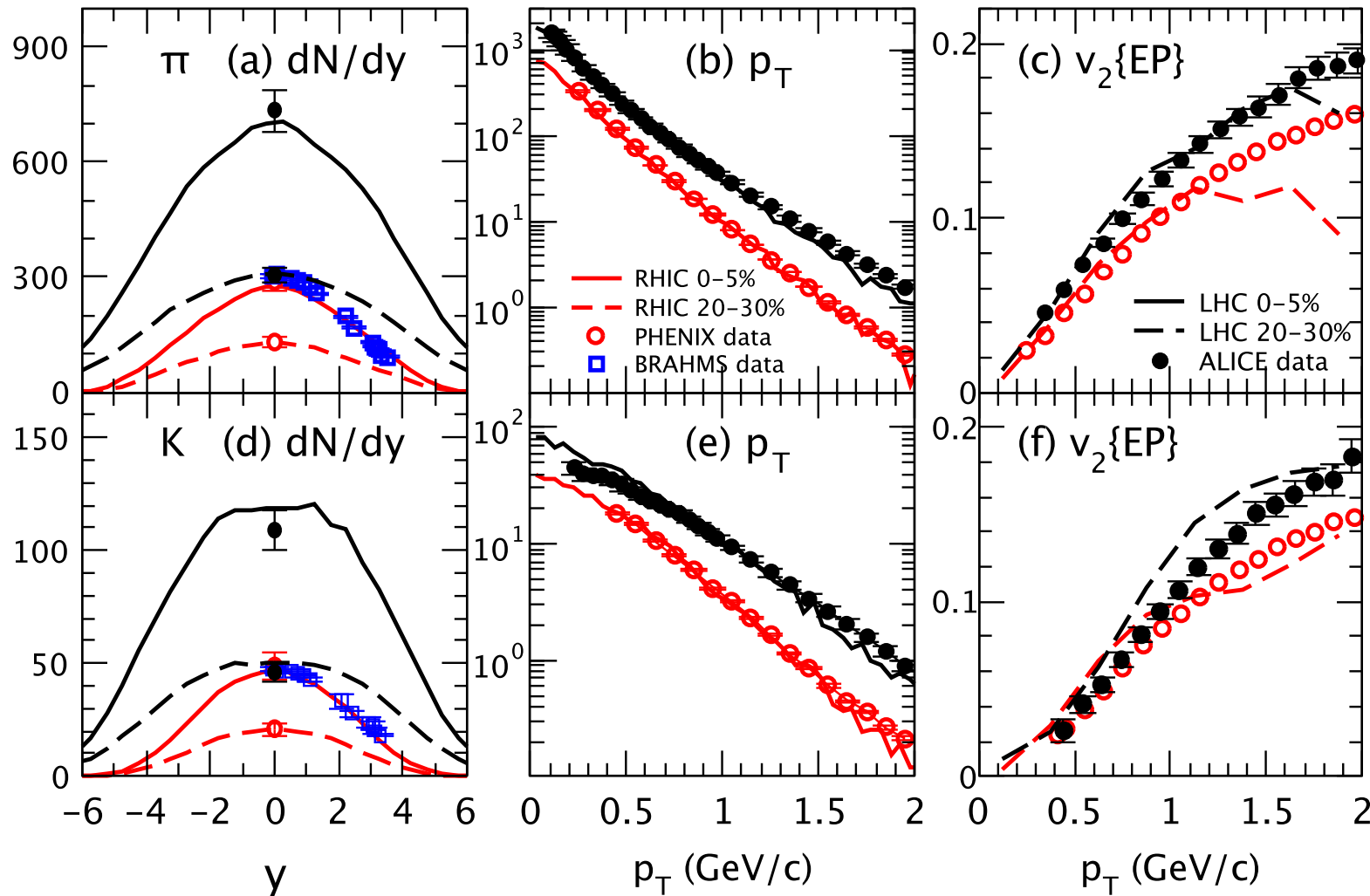
Wiranata, Koch, Prakash & Wang, Phys. Rev. C 88 (2013) extended the CE formula to a mixture of hadrons under elastic scatterings.

3) Application to parton matter in the AMPT model

We now apply the Chapman–Enskog (CE) method to study η and η/s of the parton matter in the string melting AMPT model for A+A.

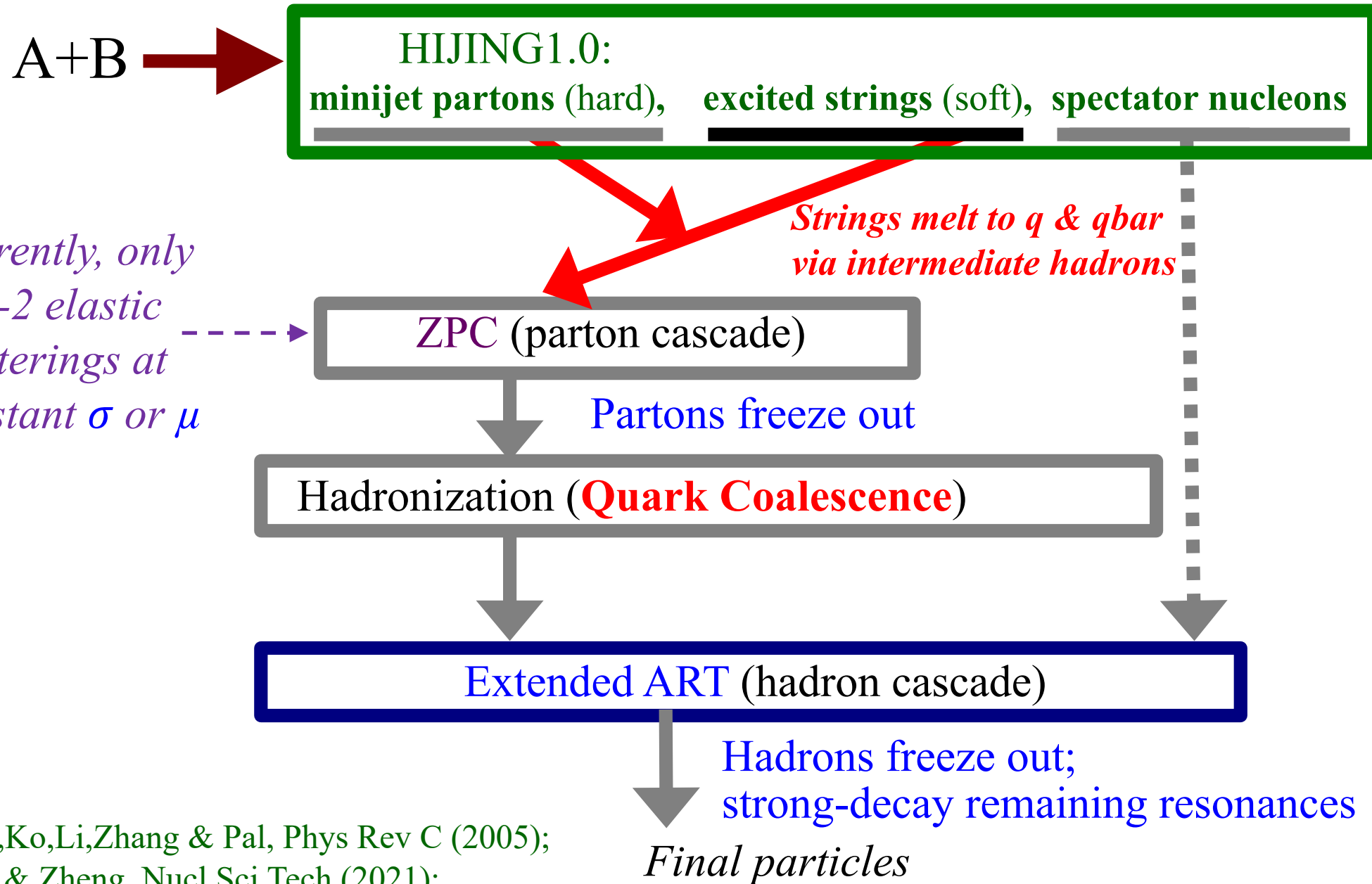
The AMPT model can reasonably describe the bulk matter observables at low p_T in A+A collisions:

ZWL, Phys Rev C (2014)



3) Application to parton matter in the AMPT model

Structure of the String Melting version of AMPT:



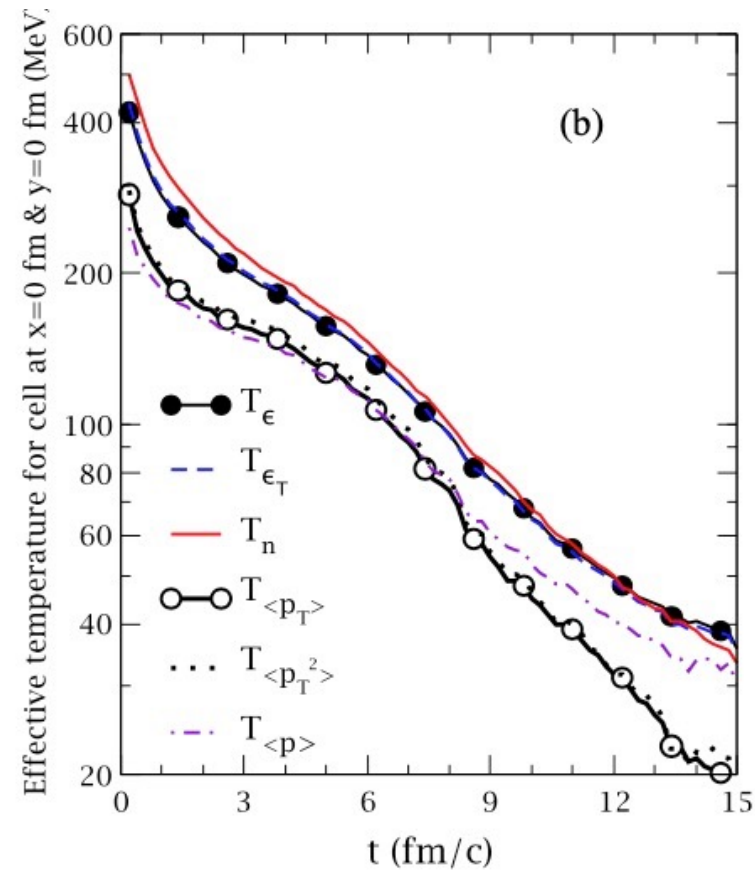
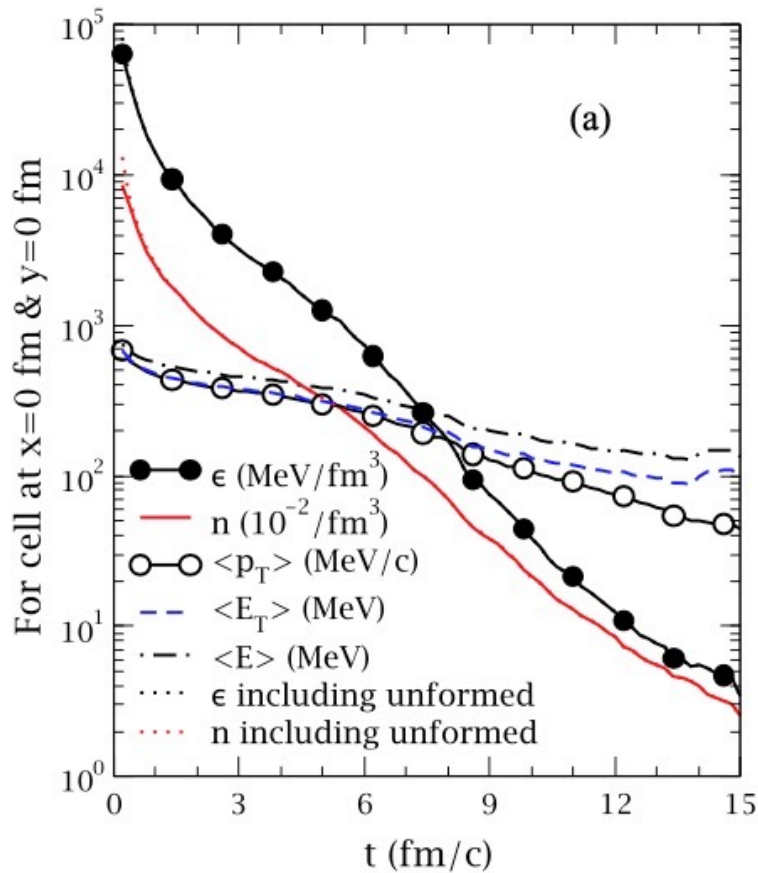
ZWL,Ko,Li,Zhang & Pal, Phys Rev C (2005);
ZWL & Zheng, Nucl Sci Tech (2021);
<https://myweb.ecu.edu/linz/ampt/>

3) Application to parton matter in the AMPT model

For parton matter in the center cell,
we have extracted the effective temperatures.

ZWL, Phys Rev C (2014)

For example, central Au+Au at 200A GeV:



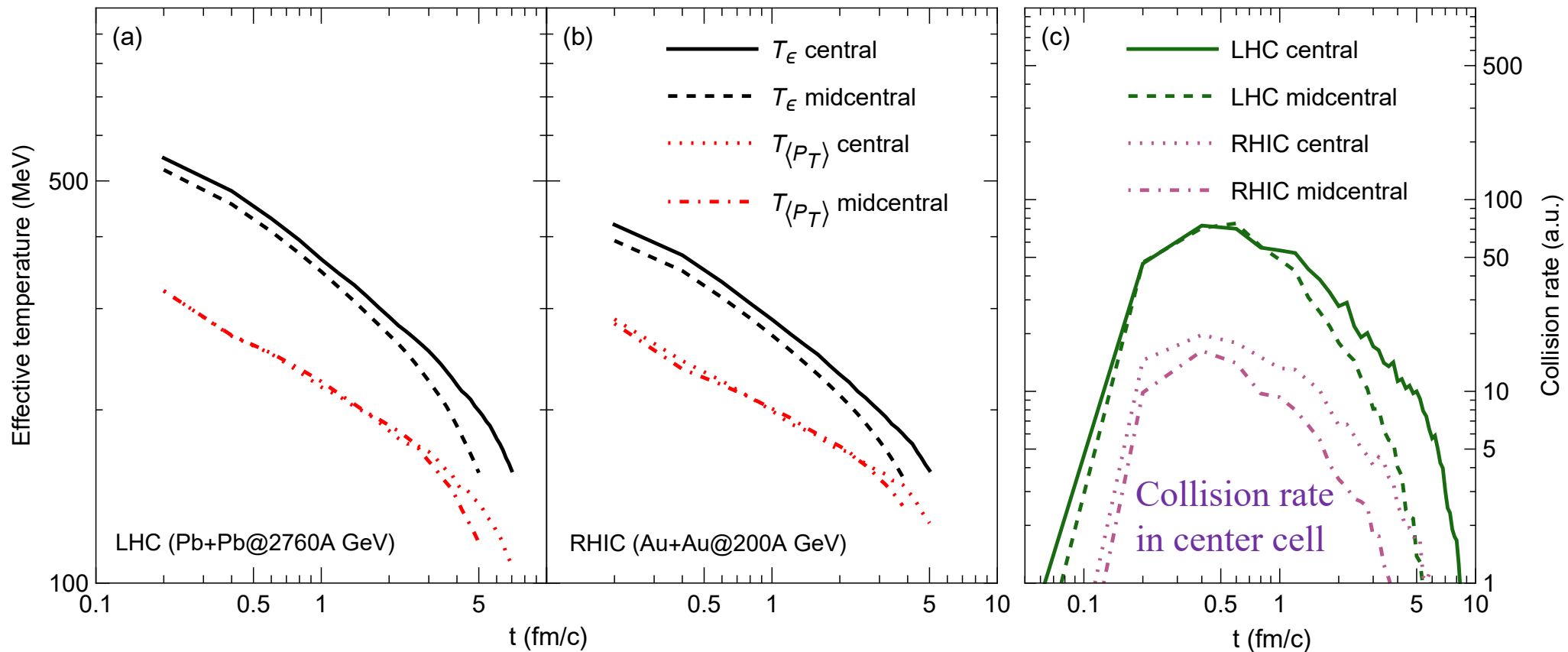
using $\epsilon = \frac{3g_B}{\pi^2} T_\epsilon^4$, $T_{\langle p_T \rangle} = \frac{4}{3\pi} \langle p_T \rangle$, ...

$T_{\langle p_T \rangle} < T_\epsilon \rightarrow$ the parton matter is not in chemical equilibrium.

3) Application to parton matter in the AMPT model

We have extracted effective temperatures $T_{\langle p_T \rangle}$ & T_ϵ of the center cell for 4 different collision systems:

ZWL, Phys Rev C (2014)



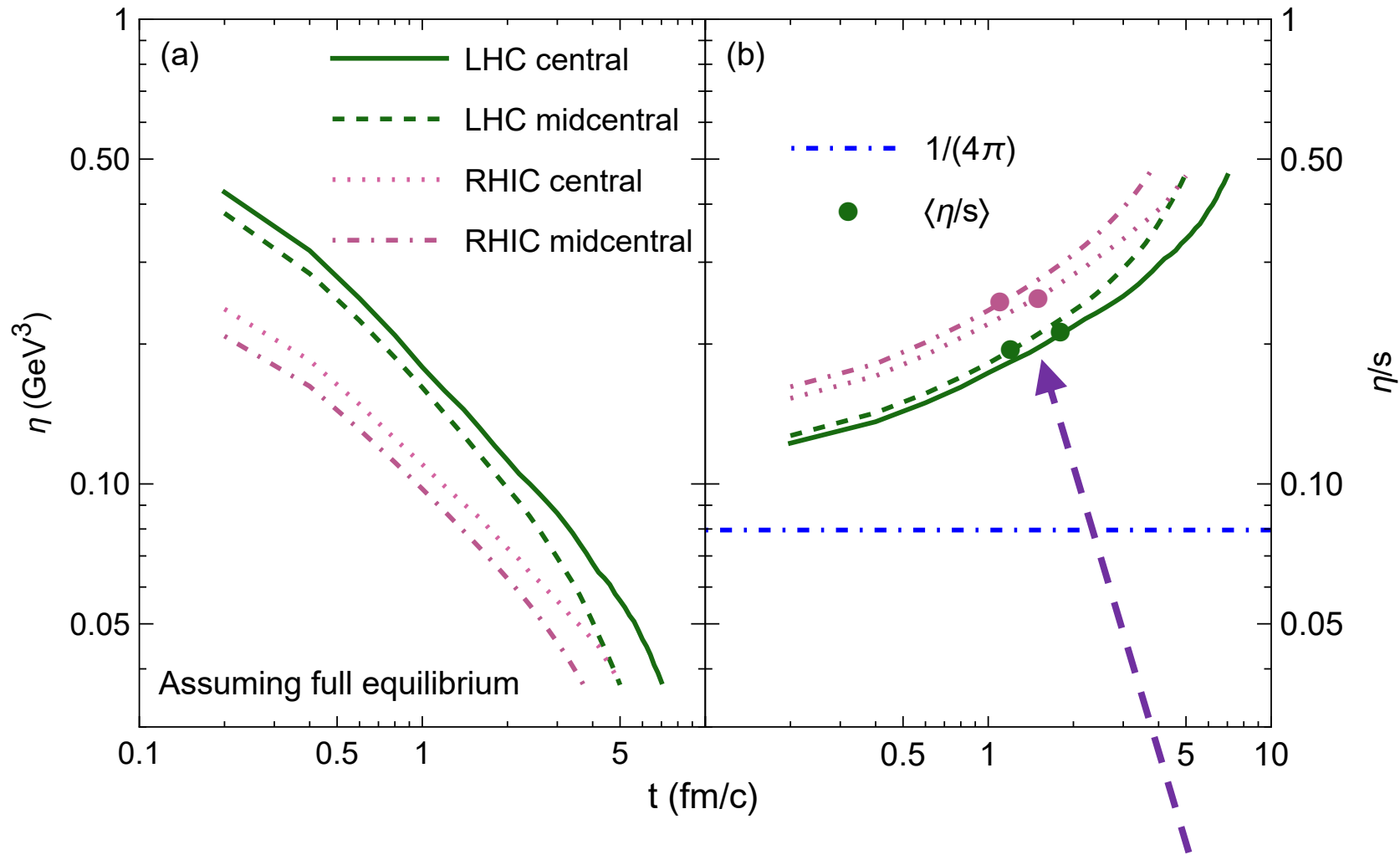
MacKay & ZWL, Eur Phys J C (2022)

We use these temperatures to calculate η and η/s of the center cell

(a volume around mid-pseudorapidity with 1 fm^2 in transverse area)

3) Application to parton matter in the AMPT model

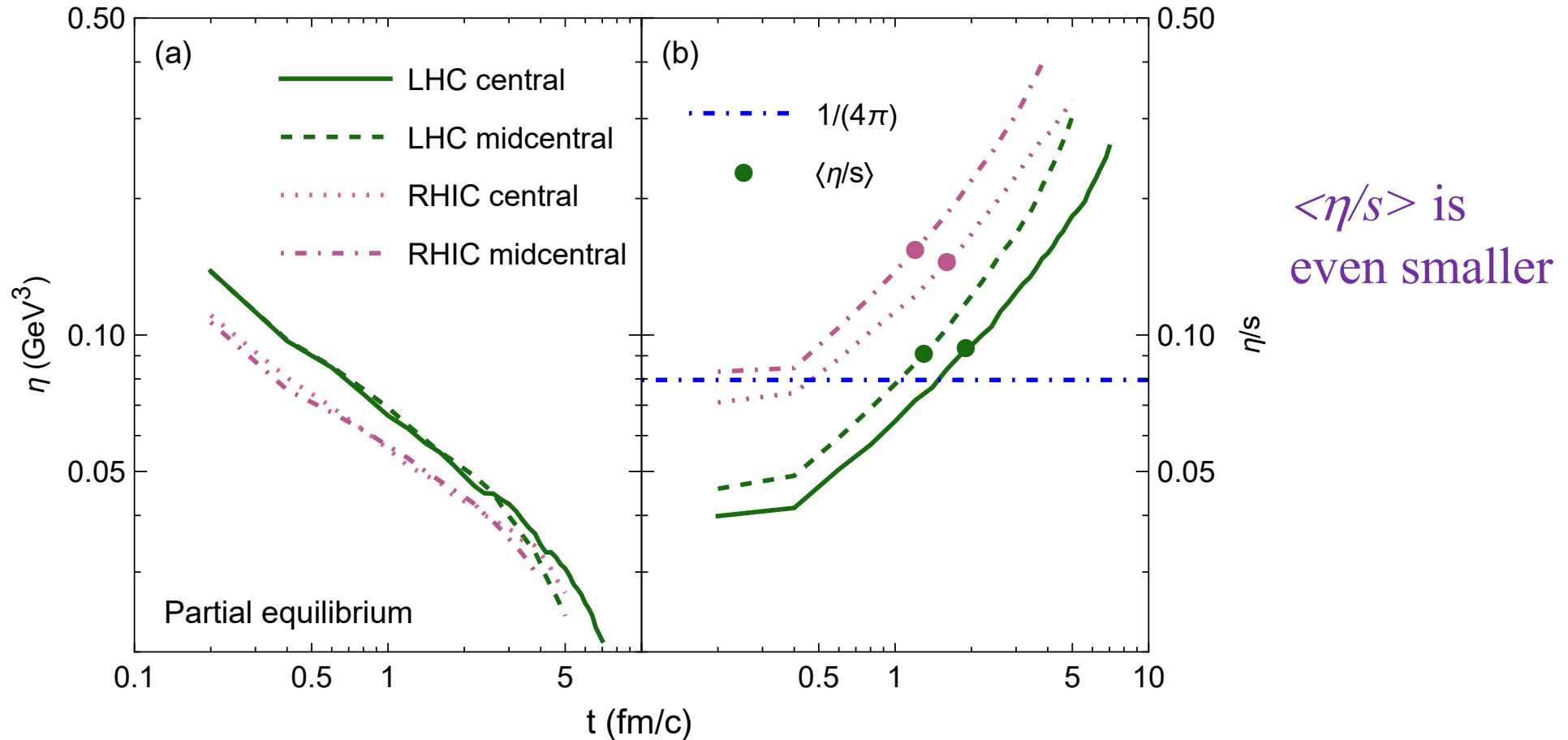
- When treating the matter as a QGP in full equilibrium ($N_f=3$), we use temperature T_ε to calculate both η and s :



- Time-averaged value weighted by collision rates: $\langle \eta/s \rangle$ is quite small.
- Temperature dependence of η/s is “wrong”, due to constant σ

3) Application to parton matter in the AMPT model

- When treating the matter as a QGP in partial chemical equilibrium, we use temperature $T_{\langle p_T \rangle}$ to calculate η but use T_ε to calculate s , since η is determined by momentum transfer but not density:



- η & η/s are lower in partial equilibrium due to $T_{\langle p_T \rangle} < T_\varepsilon$:
lower T (at constant μ) makes scattering more isotropic and effective

3) Application to parton matter in the AMPT model

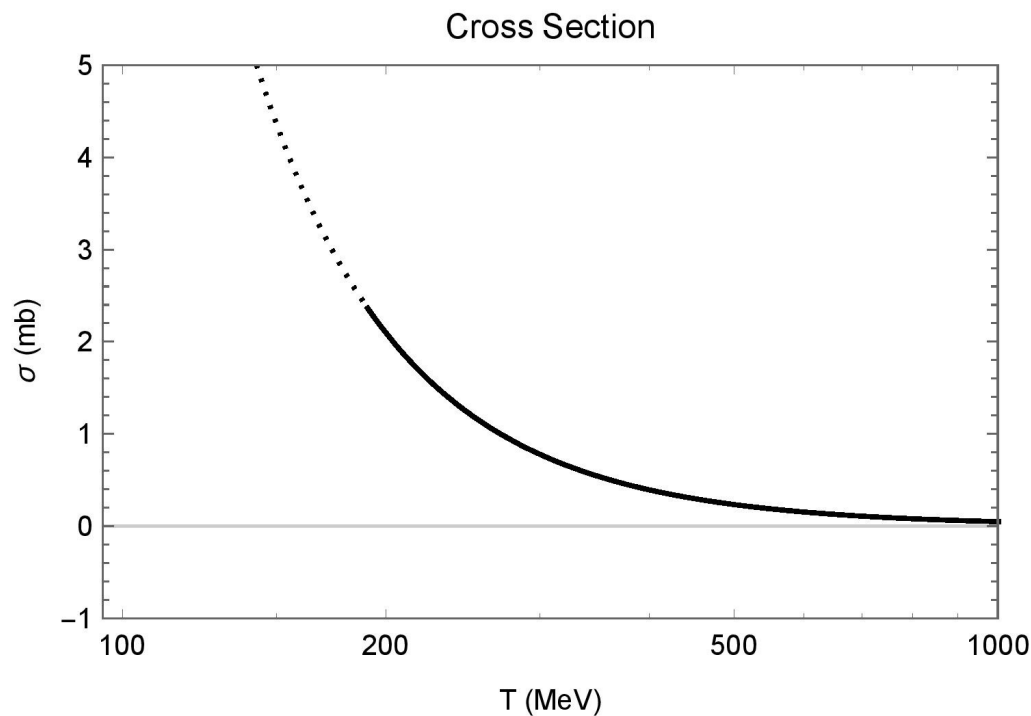
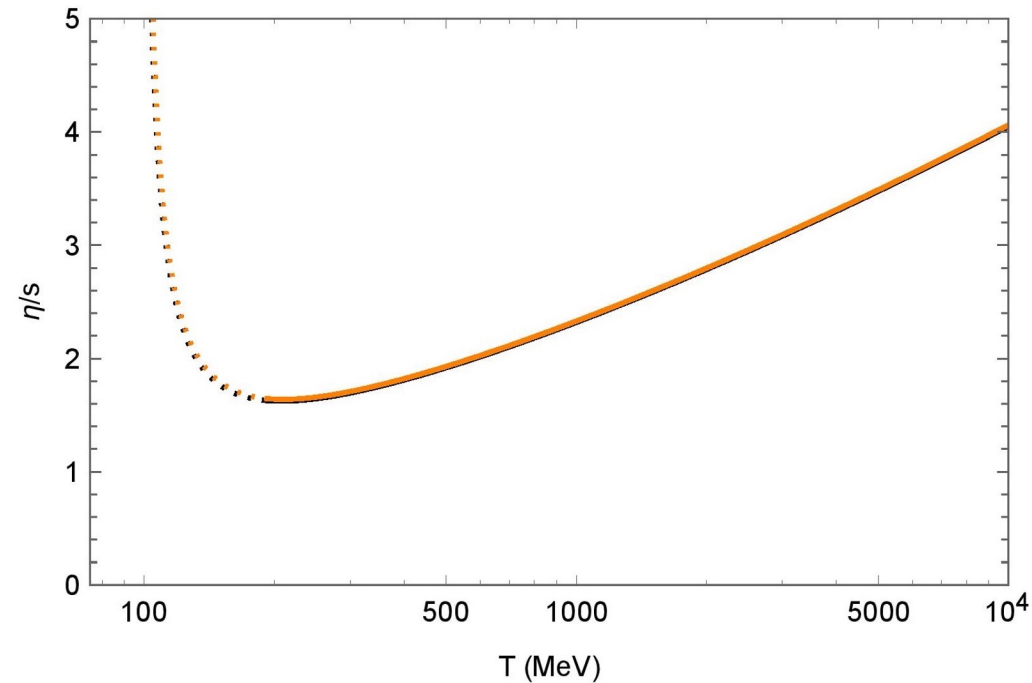
pQCD

with $\mu \propto gT$:

$$\frac{\eta}{s} \approx \frac{5.12}{g^4 \ln(2.42/g)}$$

Arnold, Moore & Yaffe, JHEP (2003);

Csernai, Kapusta & McLerran, Phys Rev Lett (2006)



If we use $\mu \propto gT$

→ $\sigma \propto 1/\mu^2$ will be larger at lower T

→ $\eta/s \propto T/\sigma$ will have the expected
 T - and t -dependences

→ a direction to improve ZPC/AMPT

Conclusions

- The Chapman–Enskog (CE) method gives accurate expression of η for parton matter under 2-to-2 scatterings
- Applying the CE method, $\langle \eta/s \rangle$ for parton matter in the center cell of high energy A+A collisions from the AMPT model is very small at $(1-3)/(4\pi)$
- T-dependence or time-dependence of η/s in AMPT is opposite to pQCD expectation, because of the constant σ or screening mass μ
- This problem can be resolved by adopting $\mu \propto gT$; this improvement will lead to a better ZPC/AMPT as a dynamical model for non-equilibrium studies

Thank you!