

AMPT at the High Baryon Density Region

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RHIC-BES on-line seminar (Series II), June 29, 2021

Outline

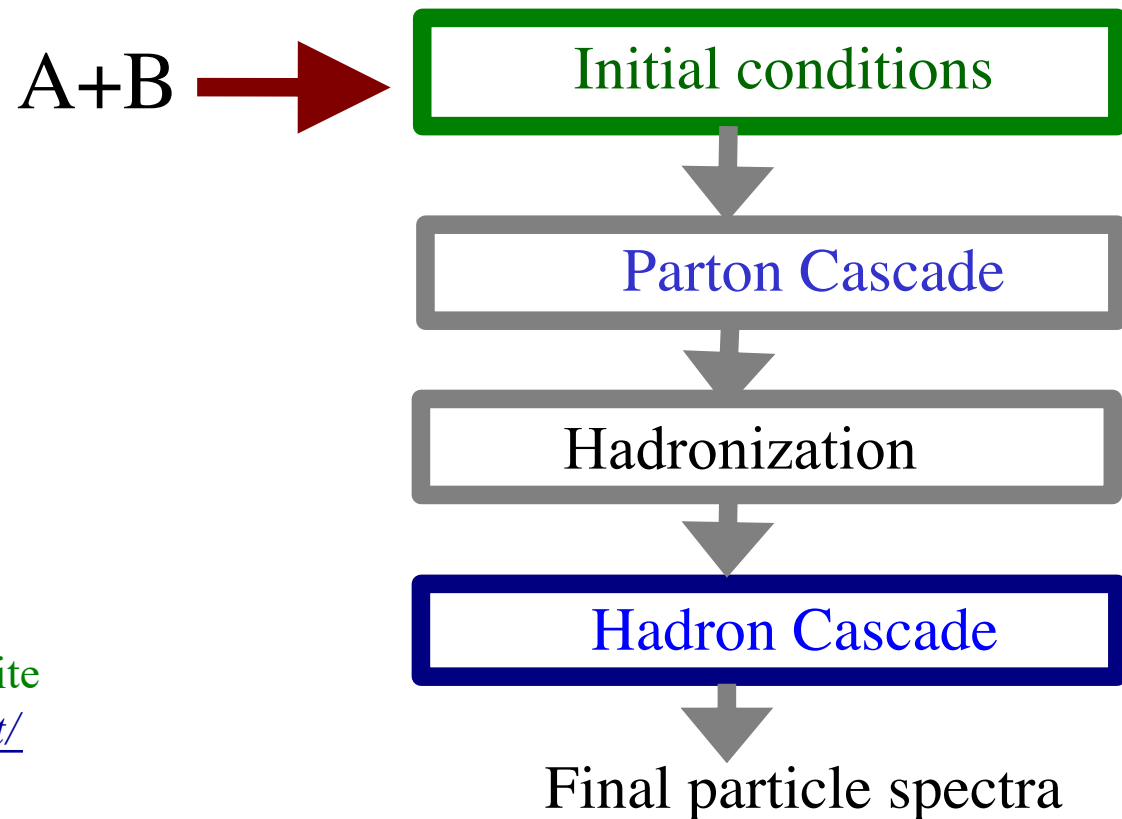
- Overview of the AMPT model
- Importance of nuclear thickness at lower energies
- Improved parton cascade and extraction of parton η/s
- Future developments for high baryon density physics
- Summary

Based on arXiv:1704.08418, 2001.10140,
2012.13825, 2102.06937, 2103.10815
*(in collaboration with Todd Mendenhall,
Xinli Zhao, Guo-Liang Ma, Yu-Gang Ma,
Han-Sheng Wang, Wei-jie Fu,
Chao Zhang, Liang Zheng, Shusu Shi)*

A Multi-Phase Transport (AMPT)

was constructed as a self-contained kinetic description of heavy ion collisions:

- evolves the system from initial condition to final observables;
- particle productions of different flavours at different P_T & y ;
- non-equilibrium initial condition & dynamics (e.g. fluctuations & correlations).



Source codes at the ECU website

<http://myweb.ecu.edu/linz/ampt/>

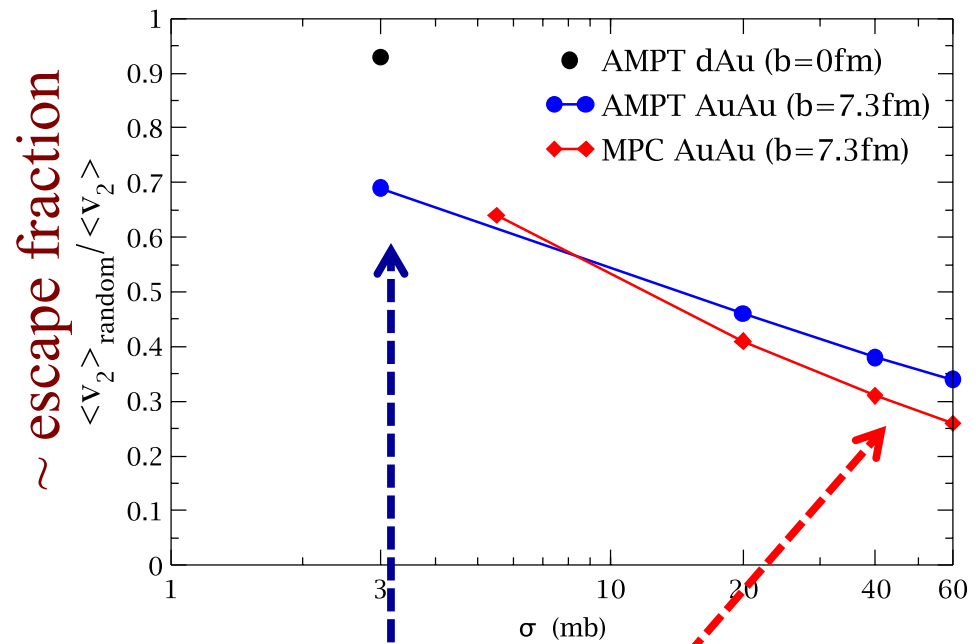
ZWL, Ko, Li, Zhang & Pal, PRC (2005);

ZWL & Zheng, NST (in press, 2021)

Transport versus hydrodynamics for finite systems

The escape mechanism:
interaction-induced response
to anisotropic geometry
from kinetic theory.

L He et al. PLB (2016),
ZWL et al. NPA (2016)



- Escape mechanism dominates v_2 for small systems & even semi-central AuAu@RHIC.
- At very large opacity (large system/energy/ σ), hydrodynamic collective flow will dominate v_2

Heiselberg and Levy, PRC (1999), Borghini, Feld and Kersting, EPJC (2018),
Kurkela, Wiedemann and Wu, PLB (2018) & EPJC (2019)

- It is important to develop transport model/kinetic theory & compare with hydrodynamics to understand physics/collectivity of finite size systems.
- Transport model (non-equilibrium, microscopic picture) & hydrodynamics (EoS, transport coefficients) nicely complement each other.

String Melting version of AMPT (AMPT-SM)

From the Bjorken formula for initial energy density in central AA collisions:

$$\epsilon(\tau) = \frac{1}{\tau A_T} \frac{dE_T(\tau)}{dy} \quad \sim \begin{matrix} 5 & 12 & 40 \\ \text{SPS} & \text{RHIC} & \text{LHC} \end{matrix} \text{ GeV/fm}^3$$

 Transverse area

Proper formation time,
taken as $0.5 \text{ fm}/c$

\gg critical energy density
for QCD phase transition:
 $\epsilon_c \sim O(1/2) \text{ GeV/fm}^3$

→ At high-enough energies,

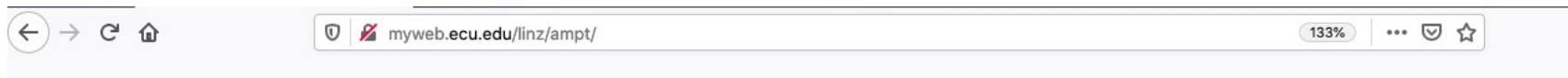
hadronic matter such as strings cannot exist at early times,

the initial matter should be represented by a high density partonic matter

→ the **string melting** version of AMPT

ZWL & Ko, PRC (2002)

AMPT codes are available online since 2004



AMPT source codes

(updated December 25, 2018):

A Multi-Phase Transport (AMPT) model is a Monte Carlo transport model for nuclear collisions at relativistic energies.

Each of the following versions contains:

the source codes, an example input file, a Makefile, a readme, a required subdirectory for storing output files, and a script to run the code.

1. [ampt-v1.11-v2.11.tgz](#) (11/2004)
2. [ampt-v1.21-v2.21.tgz](#) (10/2008)
3. *[Other older versions inbetween](#)*
4. [ampt-v1.26t5-v2.26t5.zip](#) (4/2015)
5. [ampt-v1.26t7-v2.26t7.zip](#) (10/2016)
6. [ampt-v1.26t7b-v2.26t7b.zip](#) (5/2018)
7. [ampt-v1.26t9-v2.26t9.zip](#) (9/2018)
8. [ampt-v1.26t9b-v2.26t9b.zip](#) (12/2018)

String Melting AMPT since 4/2015 can reasonably describe the bulk matter at high energies at RHIC and LHC.

This readme file lists the main changes up to version v1.26t9b-v2.26t9b ("t" means a version under test):

AMPT Users' Guide

```
*****
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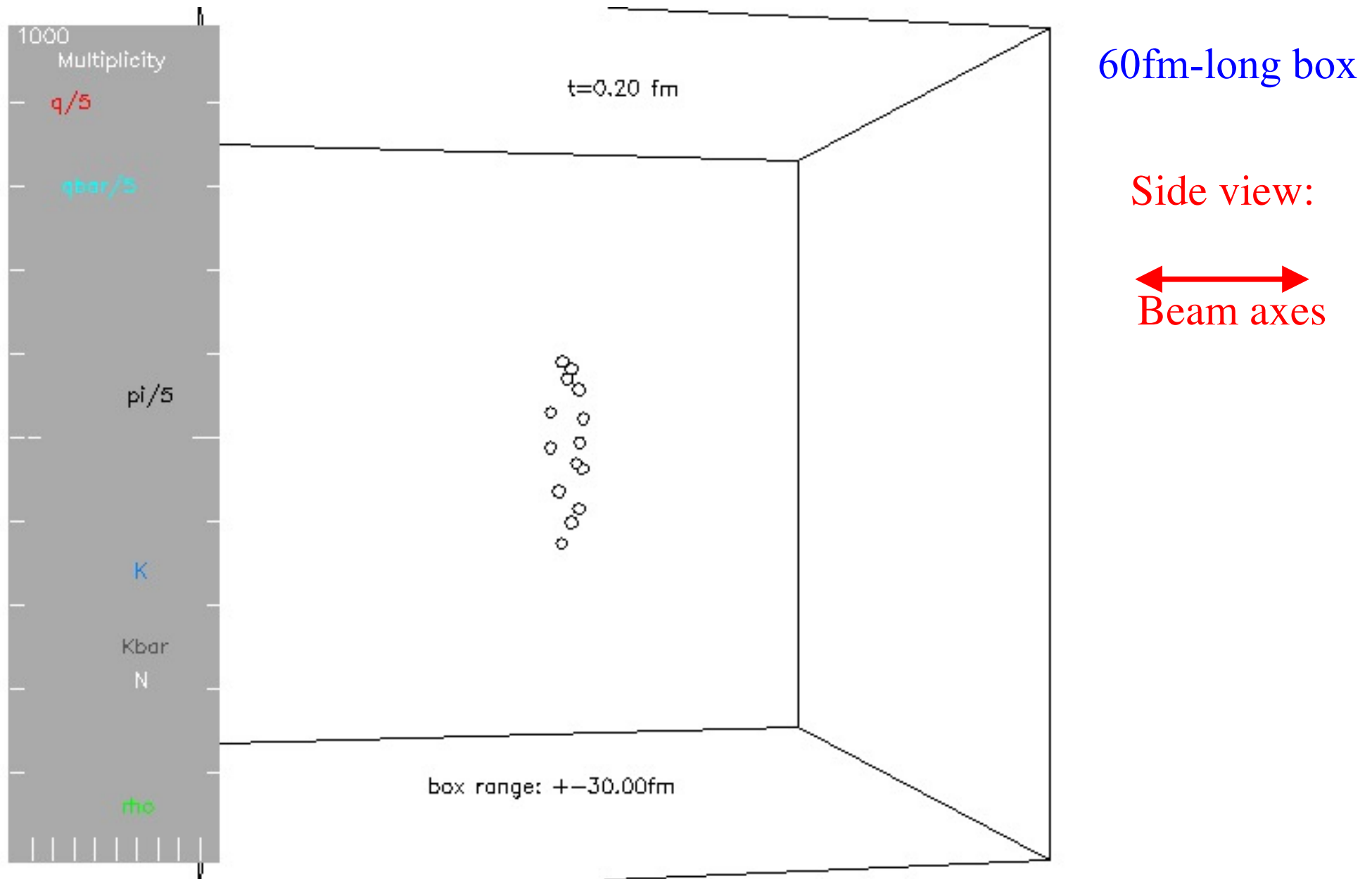
```
12/2018 test version v1.26t9b/v2.26t9b:
```

```
* Fixed bugs that can cause segmentation fault (especially for default AMPT at high energies):
```

```
exclude endpoints of 0. and 1. in random values from RANART()
  in amptsub.f in order to avoid crash in case of 0 branching ratio,
  resample random value from RLU(0) when TEVB in PYSSPA does not change
  (this may happen when RLU(0)>-0.9999 and thus
  ex==|LOG(RLU(0))*B0/(5.*WTSUM)|<-3e-8 and thus EXP(ex)=1.00000000,
  which may cause IS(2) to be undefined and be >10^8 (out of bound)).
```

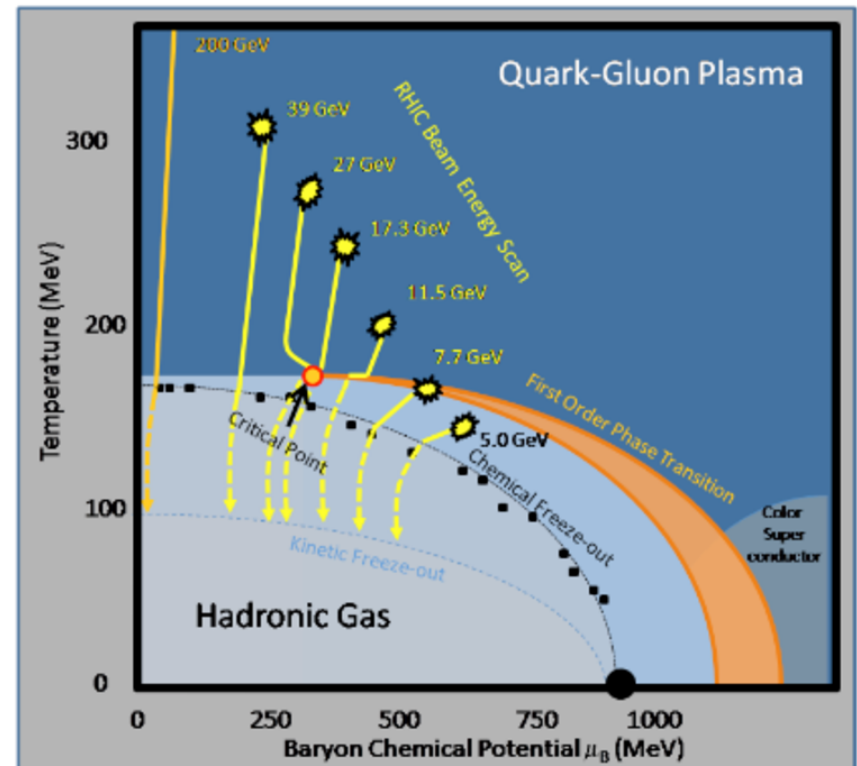
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The above modification can be found by searching "clin-12/2018".
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A central Au+Au event at 200A GeV from the AMPT-SM



Higher baryon densities / lower energies

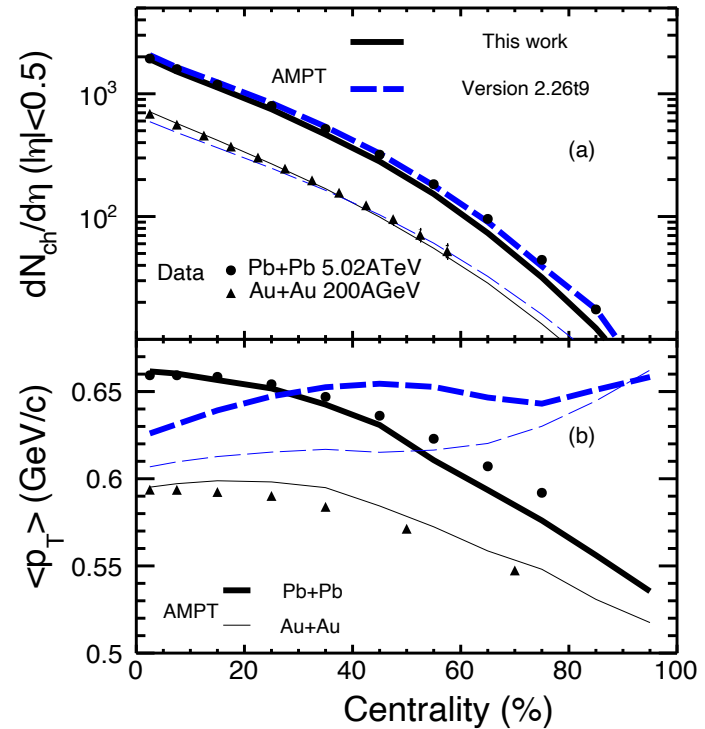
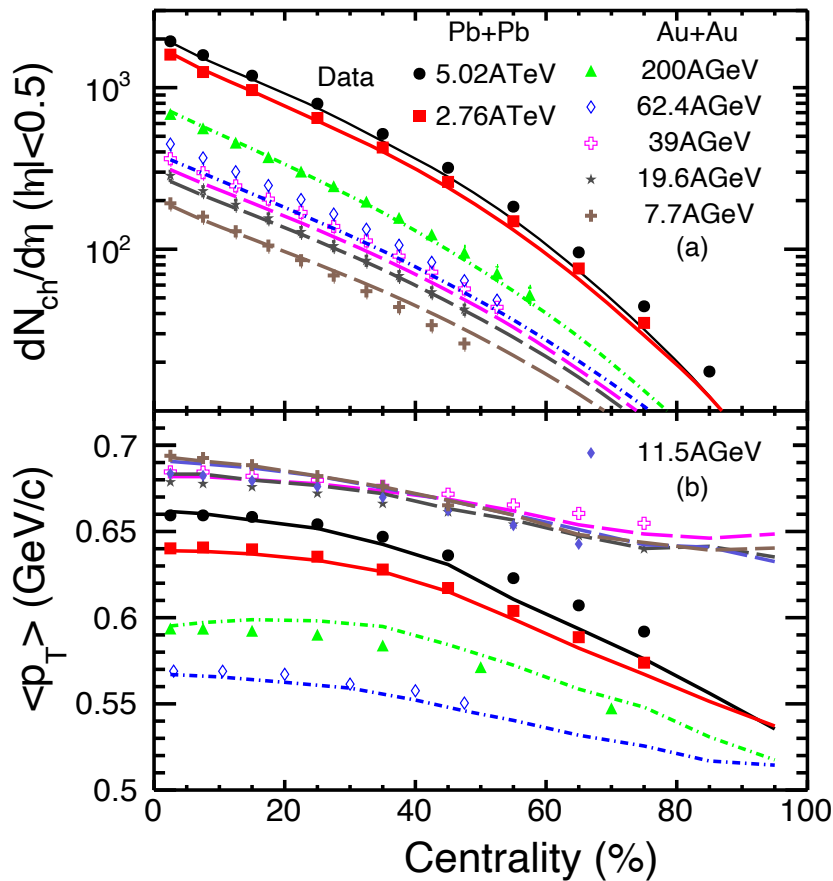
- For lower energies such as BES, particular interests are in high baryon density physics including the QCD critical end point (CEP).
- Before addressing possible effects of the critical point, one needs to know trajectory of nuclear collisions on the QCD phase diagram, including time evolutions of energy density ε & net-baryon density n_B (or T & μ_B)



from STAR arXiv:1007.2613

Results from AMPT-SM at different energies

String melting AMPT can now reasonably describe both large and small systems, including their centrality dependence after we apply local nuclear scaling on 2 key input parameters (*Lund b parameter* & *minijet pT cutoff p₀*):
 Chao Zhang et al. 2103.10815



Centrality dependence of $\langle p_T \rangle$
 is now much better than public AMPT

However, finite nuclear thickness is neglected in AMPT-SM as it was constructed for high energy nuclear collisions.

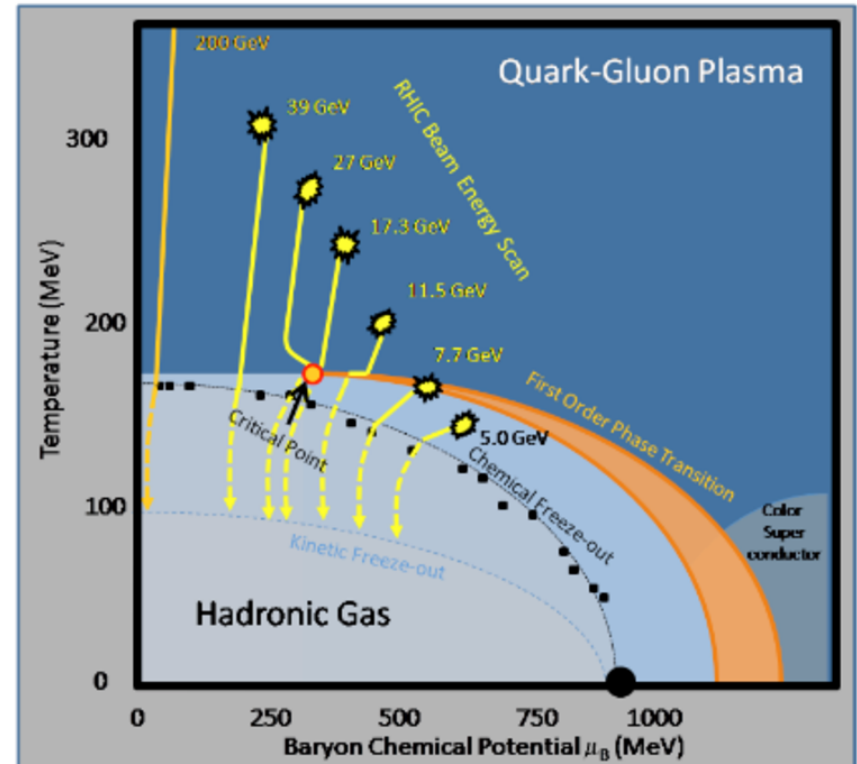
Importance of nuclear thickness at lower energies

- Effects of nuclear thickness at low energies are big on *energy density* ε & *net-baryon density* n_B (consequently on T & μ_B)

ZWL, 1704.08418,
Mendenhall & ZWL, 2012.13825,
H.S. Wang et al. 2102.06937

- For hydrodynamics models: dynamical initialization scheme is needed instead of initial condition at a fixed proper time.

Okai et al. 1702.07541,
Shen et al. 1704.04109



from STAR arXiv:1007.2613

Effect of nuclear thickness on initial energy density

For central A+A collisions in CM frame with the hard sphere model for nucleus: crossing time is

$$d_t = \frac{2R_A}{\sinh y_{CM}} = \frac{2R_A}{\gamma \beta}$$

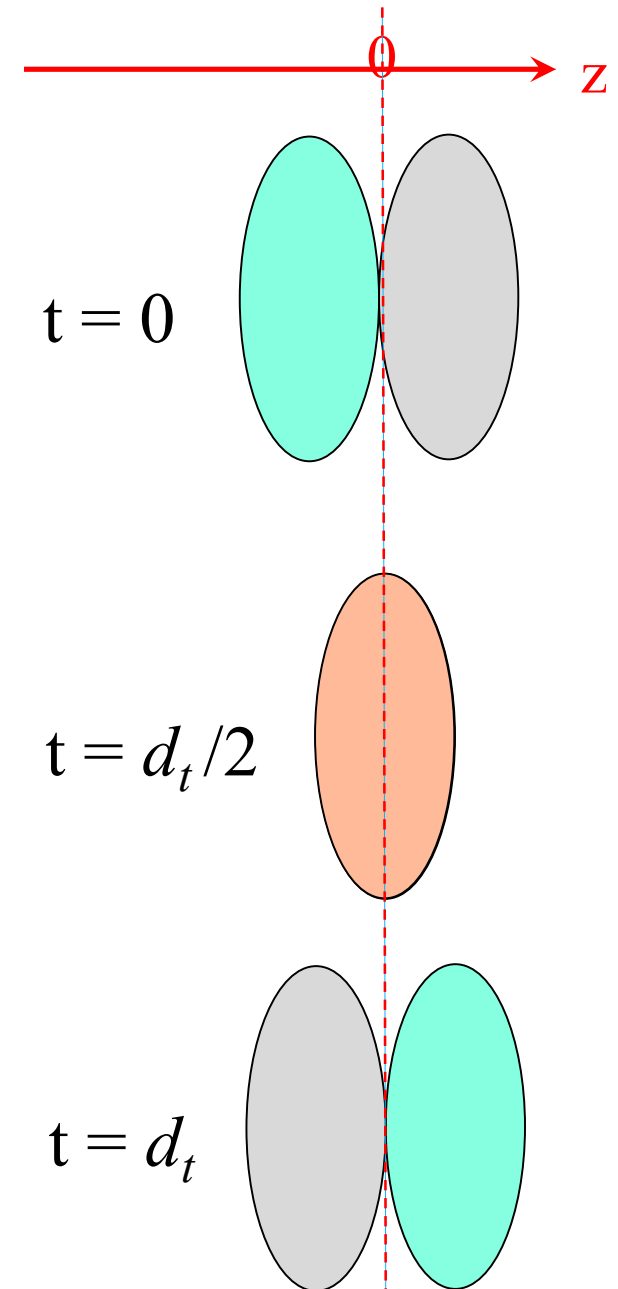
For central Au+Au collisions:

$\sqrt{s_{NN}}$ (GeV)	3	5	11.5	27	50	200
d_t (fm/c)	10.5	5.3	2.2	0.91	0.49	0.12

→ the Bjorken formula $\epsilon(\tau) = \frac{1}{\tau A_T} \frac{dE_T(\tau)}{dy}$

is only valid when $d_t \ll \tau_F$

or $\sqrt{s_{NN}} > \sim 50$ GeV for $\tau_F = 0.5$ fm/c



Effect of nuclear thickness on initial energy density

A schematic picture:

The shaded area

is the primary collision region that can contribute to $\epsilon(t)$, after considering formation time $t_F = \tau_F \cosh(y)$.

At late $t (> d_t + \tau_F)$, $\epsilon(t)$ comes from the full primary collision region (the big diamond area).

In these semi-analytical studies, we only consider the central region ($\eta_s \sim 0$) & neglect secondary scatterings or transverse expansion.

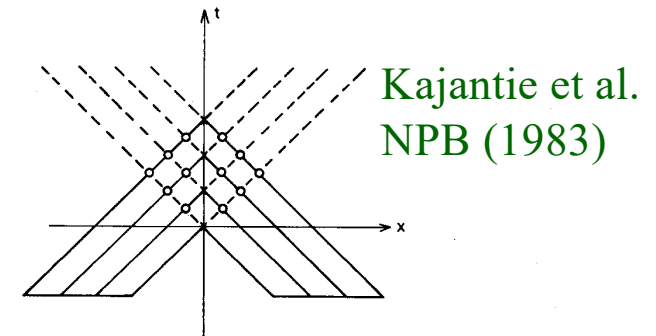
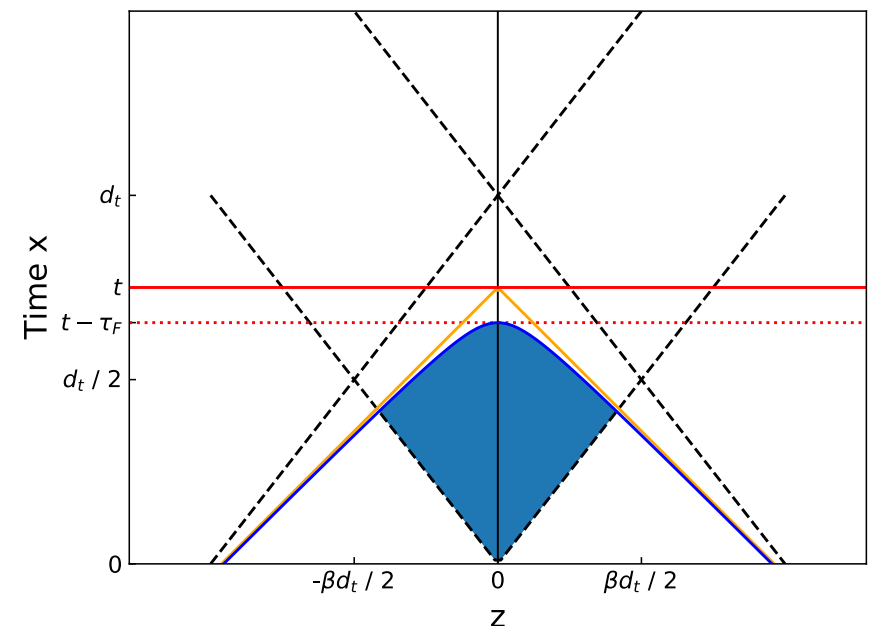


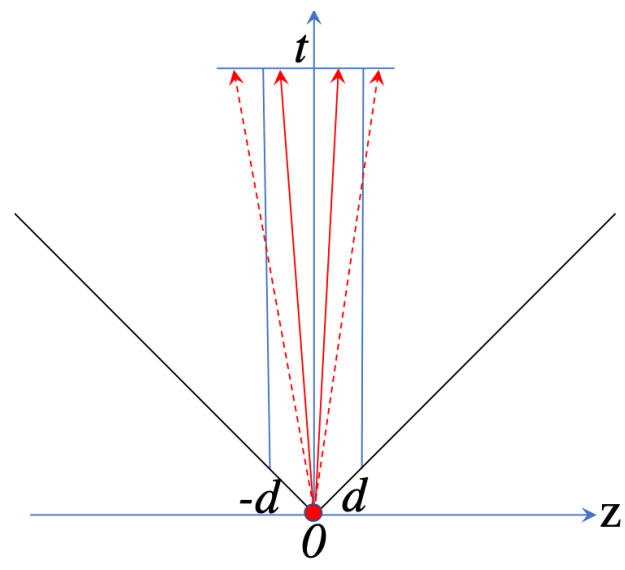
Fig. 5. An alternative description of the A + A collision. In addition to the pairwise N + N collisions on the time axis (crosses), the secondaries may further interact with the incoming nucleons (circles). This would enhance the energy density in the central region.



Mendenhall & ZWL 2012.13825

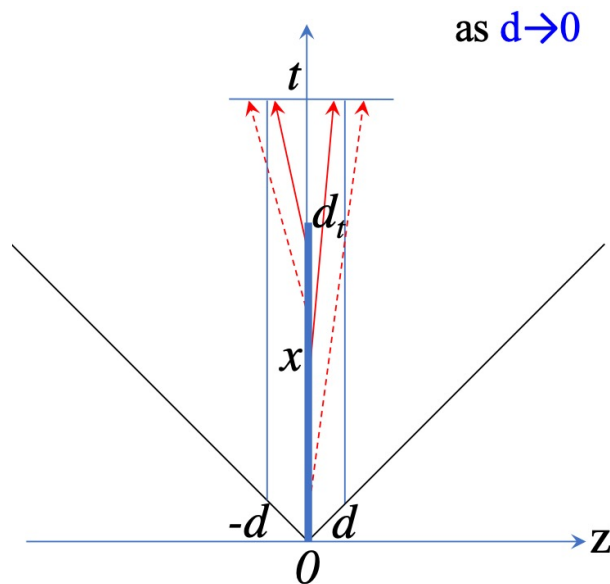
x : production time, $\in [0, d_t]$

Extension of Bjorken ε formula with nuclear thickness: 2)



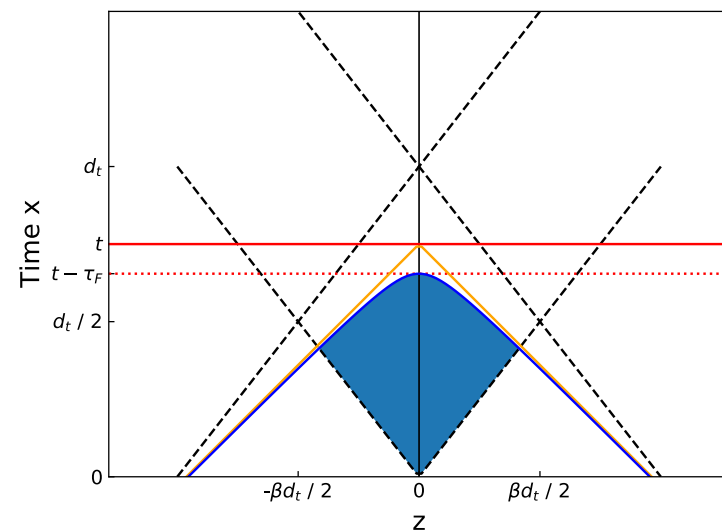
1) Without finite t or z :
the Bjorken ε formula

Bjorken, PRD (1983)



2) With finite t
(but not finite z -width)

ZWL 1704.08418



3) With both finite t & z
Mendenhall & ZWL 2012.13825

We first use this simpler method
to illustrate the qualitative effect
of nuclear thickness on initial $\varepsilon(t)$

Extension of the Bjorken ε formula: the uniform time profile

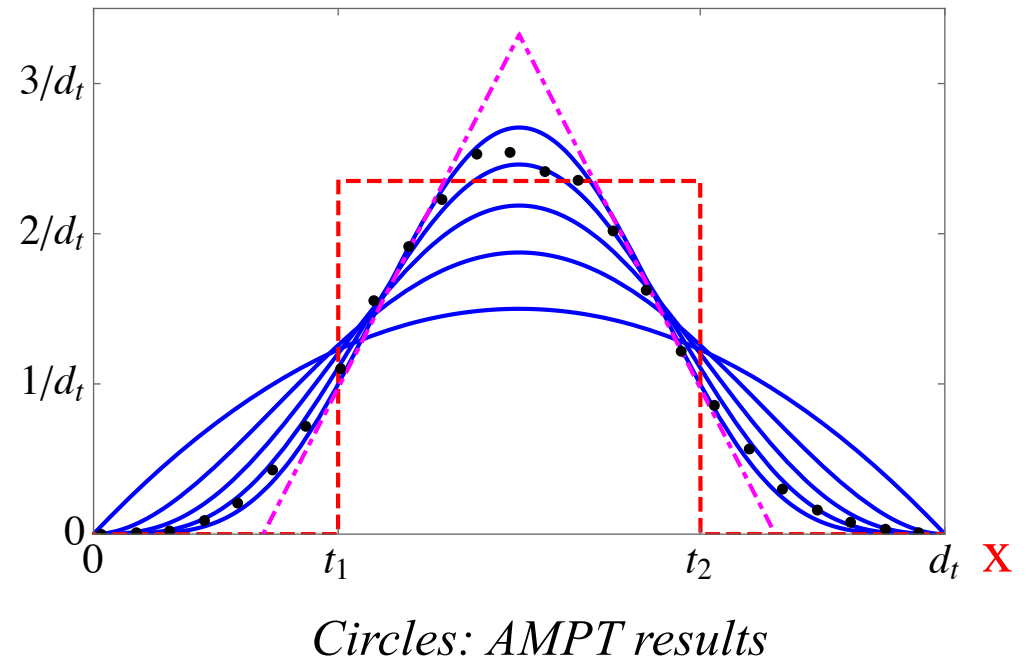
$$\varepsilon(t) = \frac{1}{A_T} \int_0^{t-\tau_F} \frac{dx}{(t-x)} \frac{d^2 E_T}{dy_0 dx}$$

For the simplest uniform profile,
initial energy (at $\eta_s \sim y_0 \sim 0$)
is produced uniformly
from time t_1 to t_2 :

$$\frac{d^2 E_T}{dy_0 dx} = \frac{1}{t_{21}} \frac{dE_T}{dy_0}$$

for $x \in [t_1, t_2]$,
with $t_{21} \equiv t_2 - t_1$

ZWL 1704.08418

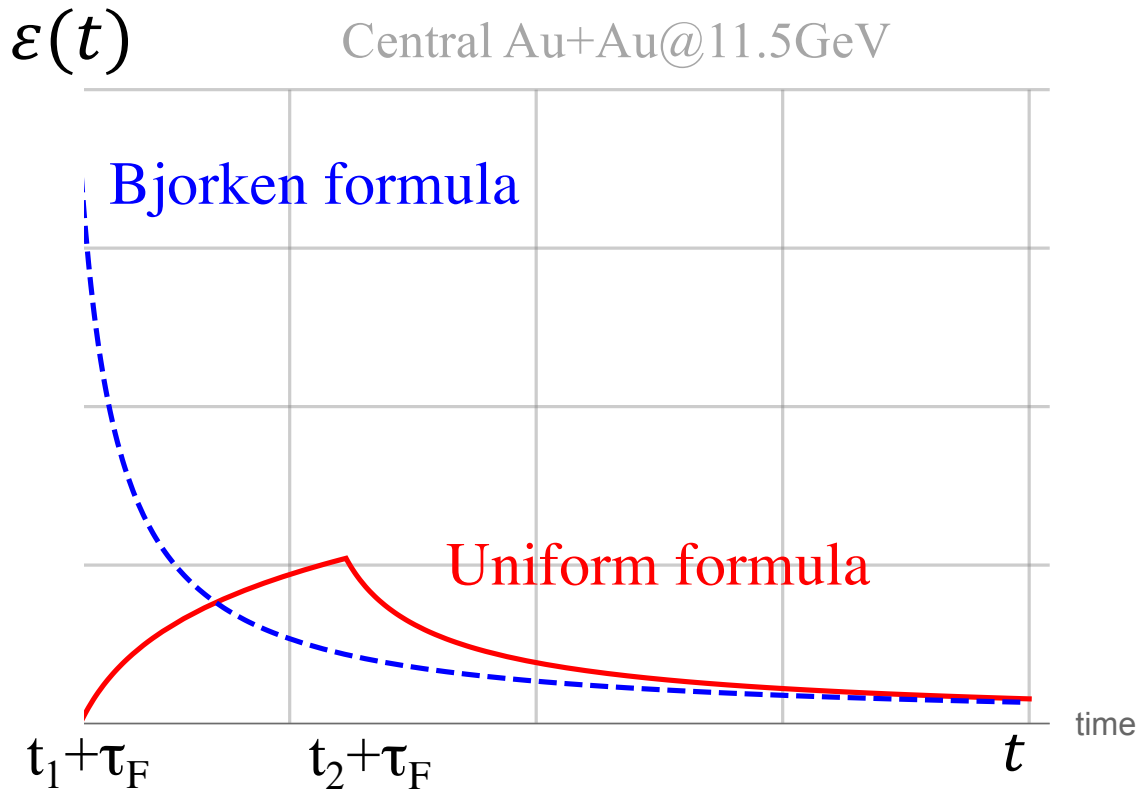


$dE_T/dy(y=0)$ parameterization from PHENIX PRC (2005)

Extension of the Bjorken ϵ formula: **the uniform time profile**

→ **solution:**
$$\epsilon_{\text{uni}}(t) = \frac{1}{A_{\text{T}} t_{21}} \frac{dE_{\text{T}}}{dy} \ln\left(\frac{t - t_1}{\tau_{\text{F}}}\right), \text{ if } t \in [t_1 + \tau_{\text{F}}, t_2 + \tau_{\text{F}}];$$

$$= \frac{1}{A_{\text{T}} t_{21}} \frac{dE_{\text{T}}}{dy} \ln\left(\frac{t - t_1}{t - t_2}\right), \text{ if } t \geq t_2 + \tau_{\text{F}}.$$



- *At high energies* (thin nuclei, or $t_{21}/\tau_{\text{F}} \rightarrow 0$):
 $\epsilon_{\text{uni}}(t) \rightarrow \epsilon_{\text{Bj}}(t)$
analytically.
- *At lower energies:*
very different from Bjorken.

Extension of the Bjorken ϵ formula: the uniform time profile

Peak energy density $\epsilon_{\text{uni}}^{\text{max}} = \epsilon_{\text{uni}}(t_2 + \tau_F) = \frac{1}{A_T t_{21}} \frac{dE_T}{dy} \ln\left(1 + \frac{t_{21}}{\tau_F}\right)$

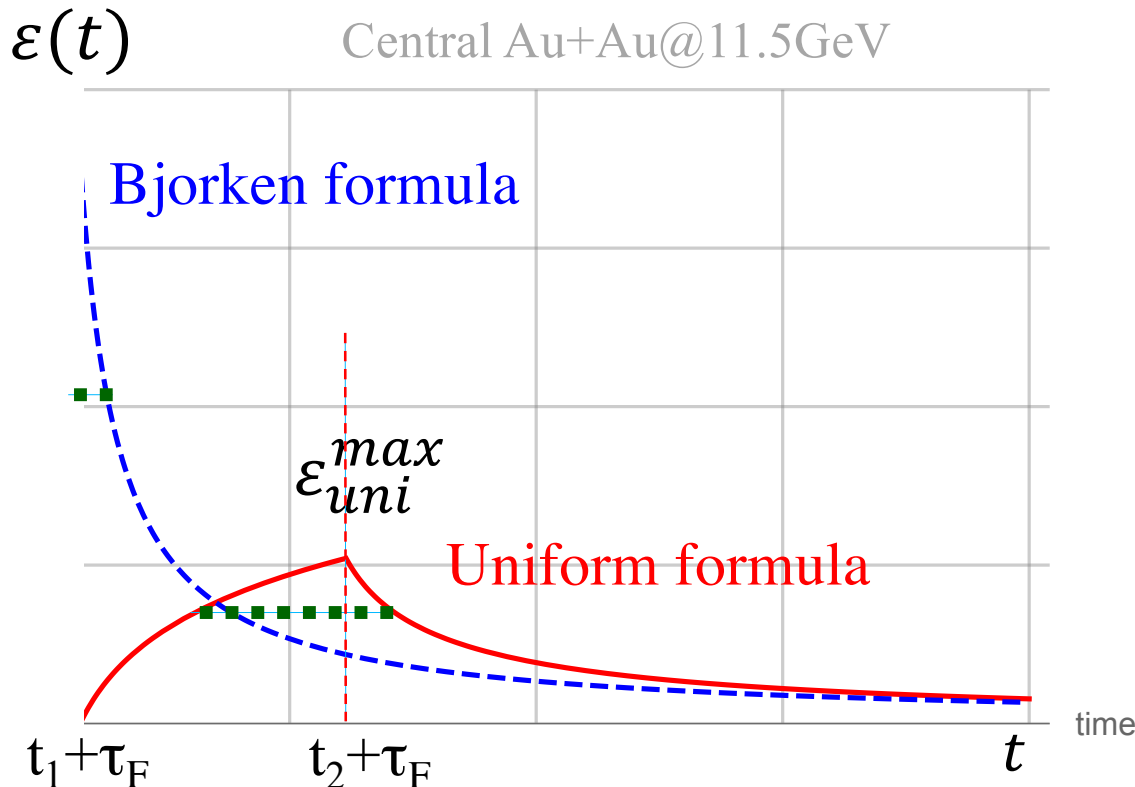
→ ratio over Bjorken: $\frac{\epsilon_{\text{uni}}^{\text{max}}}{\epsilon_{\text{Bj}}(\tau_F)} = \frac{\tau_F}{t_{21}} \ln\left(1 + \frac{t_{21}}{\tau_F}\right) \leq 1$ always.

At very low energies ($t_{21}/\tau_F \gg 1$):
ratio over Bjorken → 0;

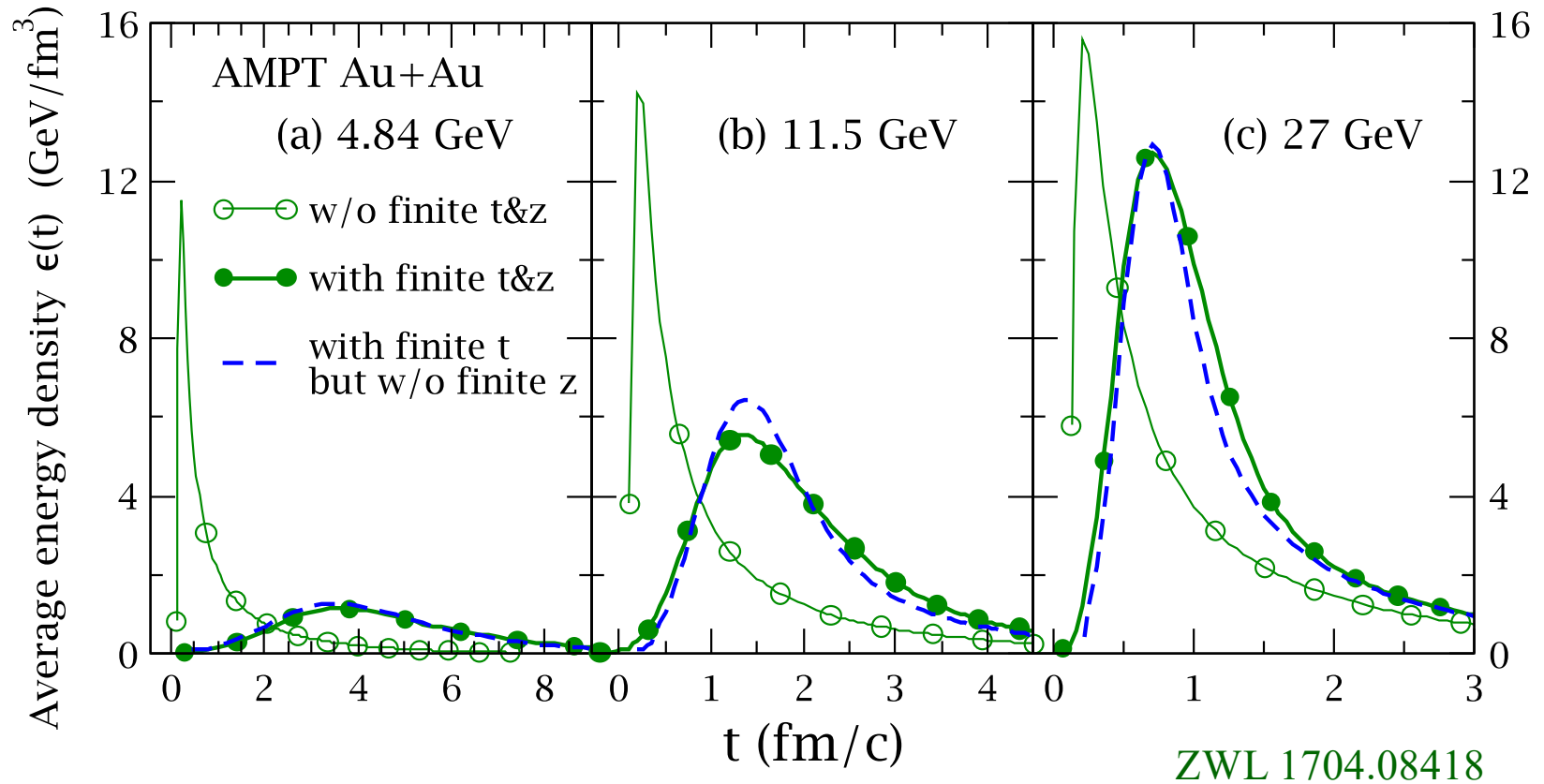
$$\epsilon_{\text{uni}}^{\text{max}} \propto \ln\left(\frac{1}{\tau_F}\right), \quad \text{not } \frac{1}{\tau_F}.$$

So the peak energy density

- \ll Bjorken value
- much less sensitive to τ_F
- FWHM width in $t \gg$ Bjorken



Nuclear thickness effects from AMPT-SM



We incorporated nuclear thickness in AMPT initial condition in a test version; the results show:

- Same qualitative features as our semi-analytical studies.
- Effect of nuclear thickness could be very important at low/BES energies.
- Peak energy density ϵ^{max} increases with $\sqrt{s_{NN}}$ much faster than Bjorken.

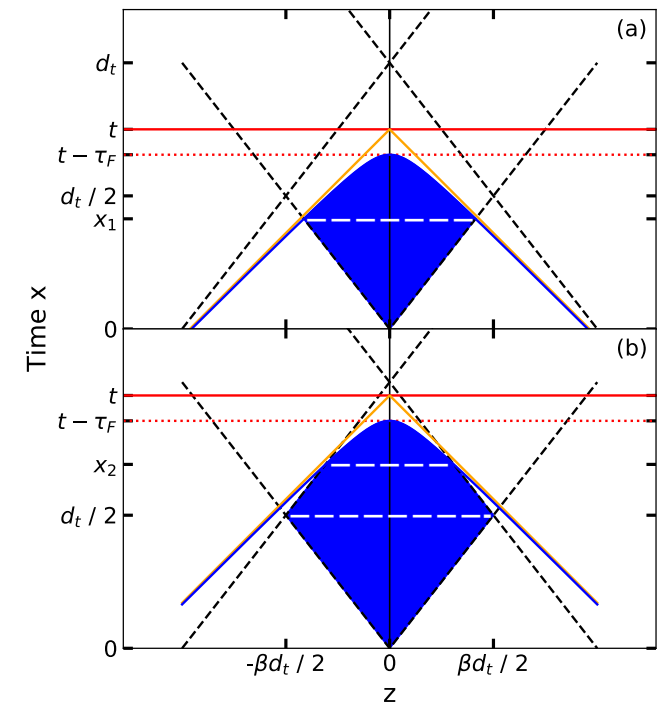
Extension of Bjorken ε formula with nuclear thickness: 3)

$$\varepsilon(t) = \frac{1}{A_T} \int_S \frac{dx dz}{(t-x)} \frac{d^3 m_T}{dy_0 dx dz} ch^3 y_0$$

S : integration area (shaded),
has 2 or 3 pieces depending on t :

3) With both finite t & z

Mendenhall & ZWL 2012.13825



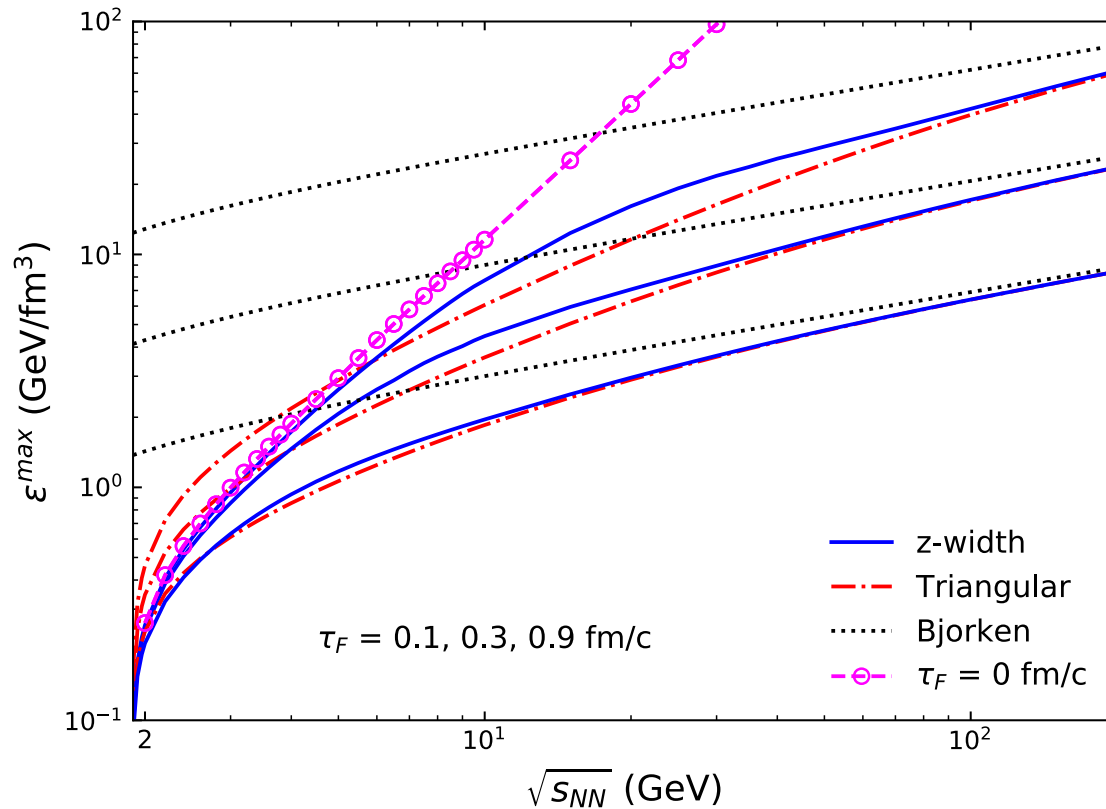
Depends on $\frac{d^3 m_T}{dy_0 dx dz}$: initial dm_T/dy production density in x - z plane:
assumed to be uniform in x - z plane
(similar to triangular time profile of earlier study ZWL 1704.08418).

dm_T/dy is based on improving $dE_T/dy(y=0)$ parameterization from PHENIX PRC (2005)
by including net-baryon contribution (important at low energies).

Extension of Bjorken ε formula with nuclear thickness: 3)

3) With both finite t & z

Mendenhall & ZWL 2012.13825



- Qualitatively similar to earlier study [ZWL 1704.08418](#)
 $\varepsilon^{max} \ll$ Bjorken value at low energies, \approx Bjorken value at high energies;
 ε^{max} & $\varepsilon(t)$ depend on τ_F more weakly than Bjorken at lower energies.
- **Surprise finding:** ε^{max} is **finite** at $\tau_F = 0$ at any colliding energy.

Extension of Bjorken ϵ formula with nuclear thickness: 3)

3) With both finite t & z

Mendenhall & ZWL 2012.13825

Todd Mendenhall has written a [web interface](#) to calculate *the initial energy density $\epsilon(t)$*

- Can be accessed at bottom of the [AMPT webpage](#)
- Takes input from user:

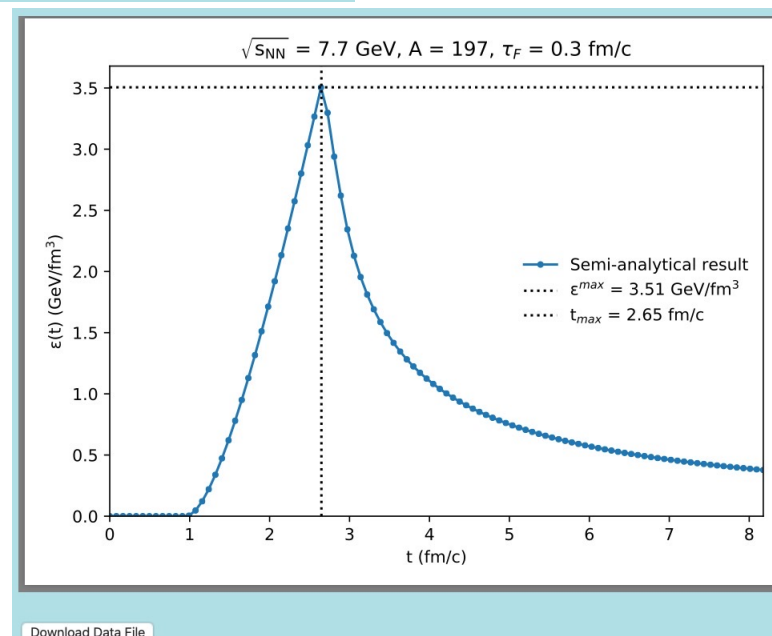
Atomic Mass Number (for projectile as well as target):

Center-of-Mass Energy per Nucleon Pair (AGeV):

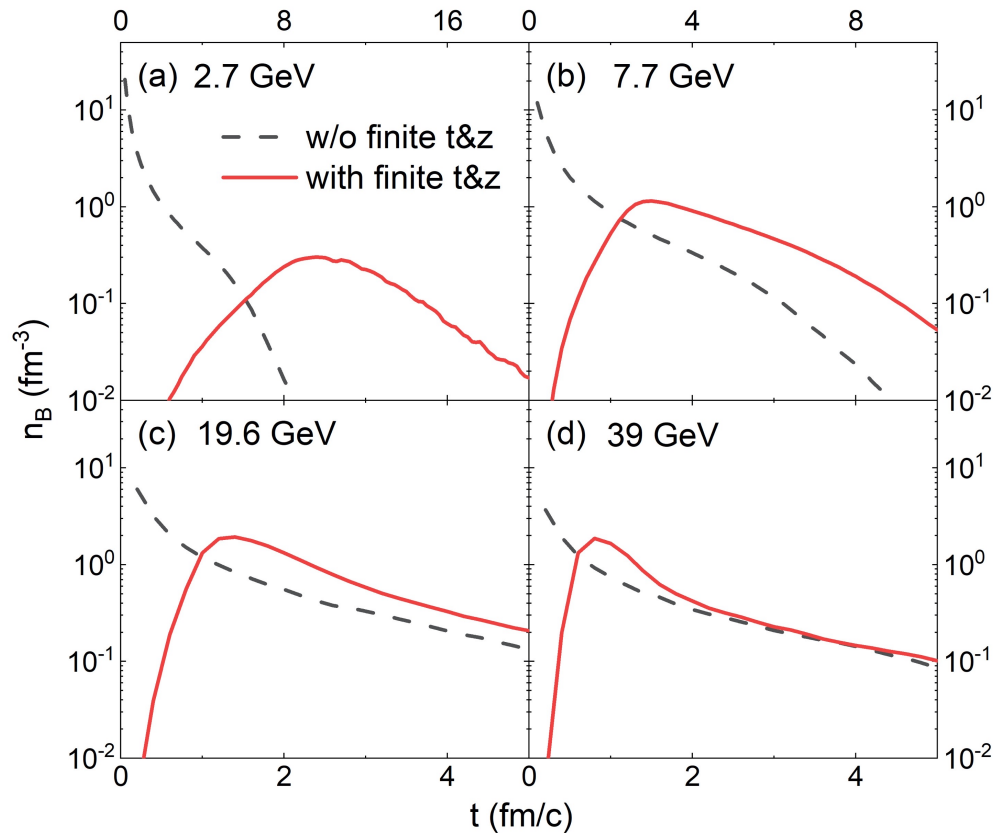
Formation time (fm/c):

Number of times to sample time evolution:

- Outputs $\epsilon(t)$ plot, user can download data file:



Nuclear thickness effects from AMPT-SM on μ_B

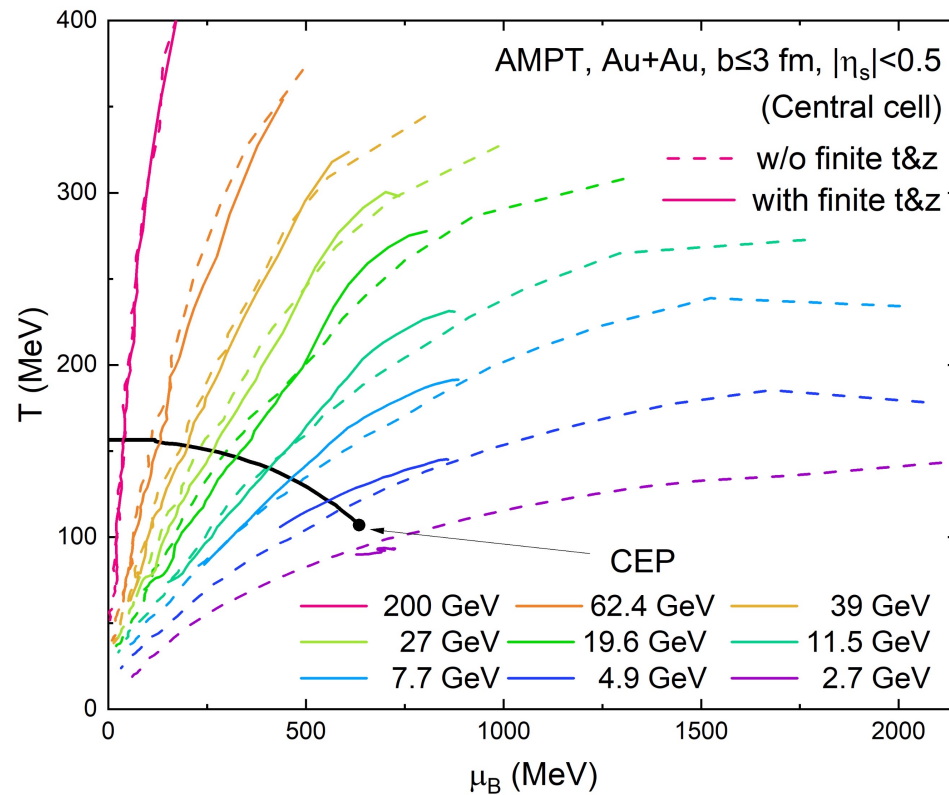


Han-Sheng Wang et al.
2102.06937,
preliminary

- Qualitatively similar to $\varepsilon(t)$ results ZWL 1704.08418, Mendenhall & ZWL 2012.13825:

n_B^{max} decreases drastically at very low energies (after including thickness); smaller change at higher energies.

Nuclear thickness effects from AMPT-SM on $T-\mu_B$



Han-Sheng Wang et al.
2102.06937,
preliminary

- Trajectory is plotted from t^{max} (time of reaching ε^{max}) to $t=10$ fm/c.
- T & μ_B both decrease significantly at very low energies (after including thickness); little change at high energies.

Improved parton cascade and extraction of parton η/s

Flows like v_2 & v_3 of large systems
mostly come from parton cascade in AMPT.

But ZPC/MPC cascade solution of the Boltzmann equation
is well known to suffer from **causality violation**.

Parton subdivision can resolve this problem:
but is very CPU-consuming
& alters e-by-e fluctuation/correlation.

We study

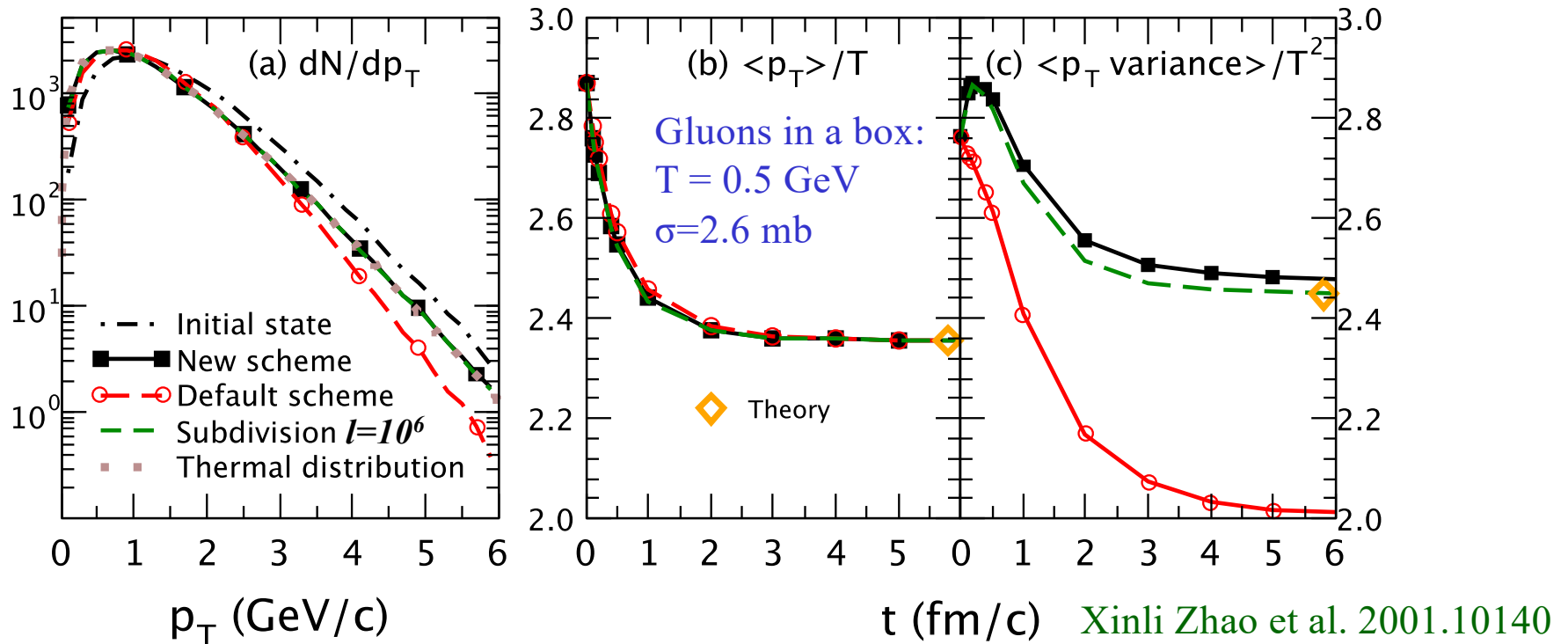
- how accurate ZPC is under expected densities/opacities
- how to accurately solve Boltzmann equation w/o subdivision
- extract parton η or η/s

Xinli Zhao et al. 2001.10140, Mendenhall & ZWL (ongoing)

Improved parton cascade and extraction of parton η/s

ZPC in AMPT numerically solves the Boltzmann equation for 2-body collisions:

$$p^\mu \partial_\mu f(\mathbf{r}, \mathbf{p}, t) = C[|\mathcal{M}|^2 f_1(\mathbf{r}_1, \mathbf{p}_1, t) f_2(\mathbf{r}_2, \mathbf{p}_2, t)]$$



We have found a **new scheme** to perform parton collisions:
 it almost eliminates **causality violation**
 & is more accurate than **default ZPC collision scheme**

Improved parton cascade and extraction of parton η/s

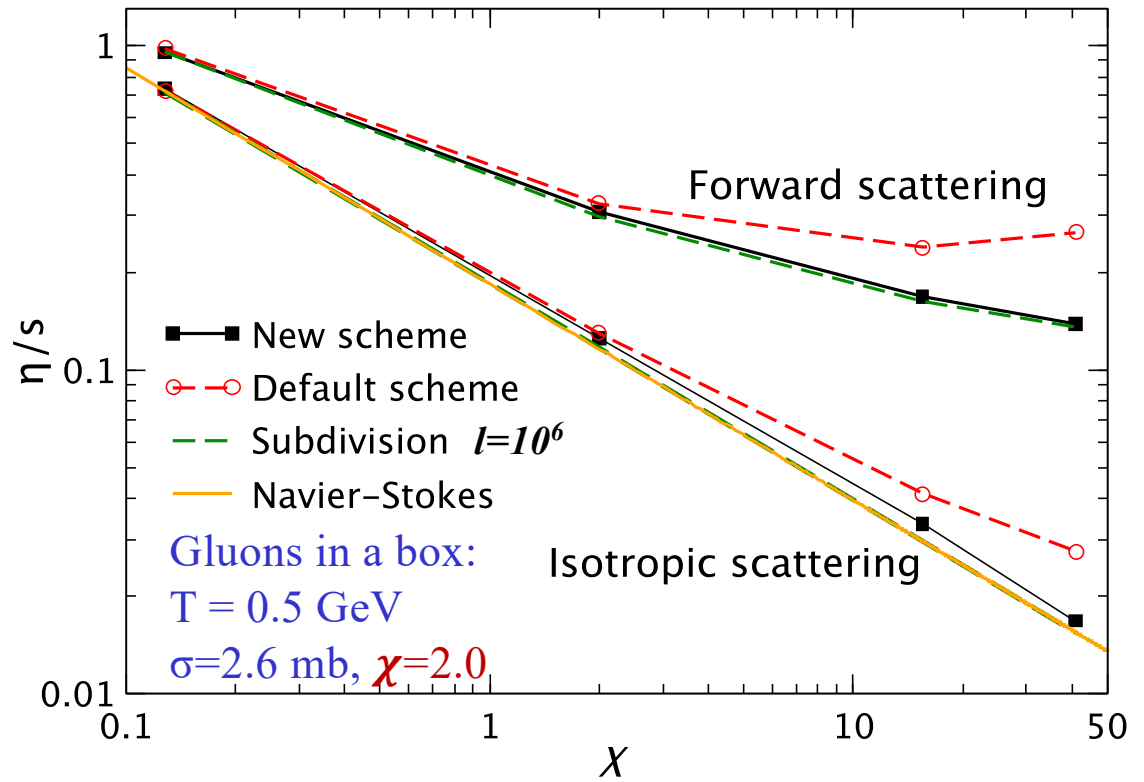
We calculate η with the Green-Kubo relation:

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

Xinli Zhao et al. 2001.10140

New scheme
 \approx Subdivision results
 $=$ Navier-Stokes values
 (for isotropic scatterings):

$$\left(\frac{\eta}{s}\right)^{NS} \approx \frac{0.1839}{\chi^{2/3}}$$



Opacity parameter $\chi = \sqrt{\frac{\sigma}{\pi}} / \lambda_{mfp} = n \sqrt{\frac{\sigma^3}{\pi}}$

Bin Zhang, Gyulassy and Pang, PRC (1998)

Next step: extract parton η/s
 of a 3-d expanding system

Future developments for high baryon density physics

Above developments of the AMPT model
lay a good foundation for studies of nuclear collisions
at low energies /high baryon densities.

Still, much important work needs to be done:

- 1) realistic Equation of State of the dense matter (*including QCD critical end point*)
- 2) inelastic parton reactions of different flavors (*QGP chemical composition*)
- 3) link parton cross sections with η/s and other transport coefficients
- 4) hadronization (*parton recombination/quark coalescence/fragmentation*)
- 5) potentials (*partonic and hadronic*)

+more

better hadron cascade

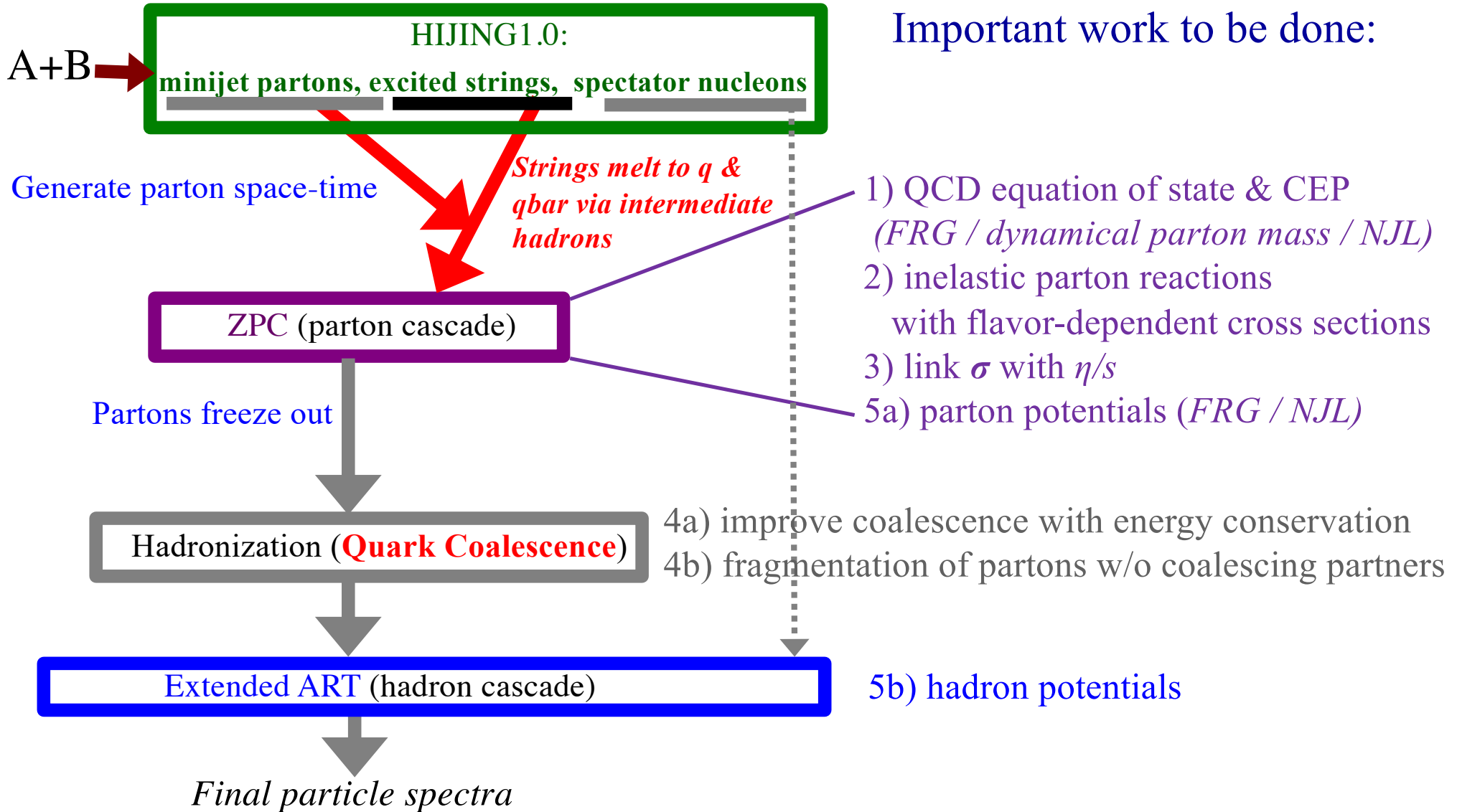
kinetics of spin polarizations

transport under electromagnetic fields

better model maintenance & integration to experiments

Future developments for high baryon density physics

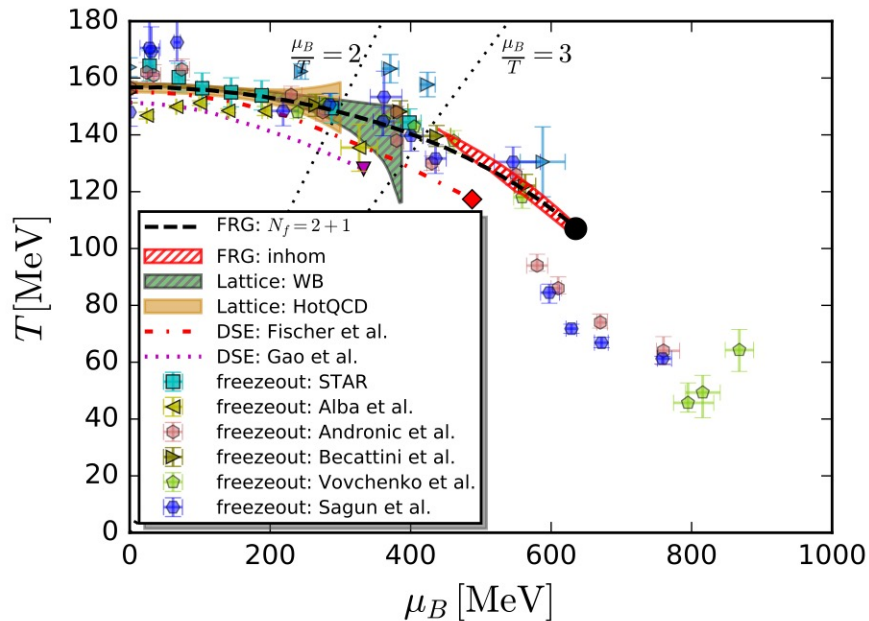
String Melting AMPT



Future developments for high baryon density physics

One direction is to couple AMPT with FRG since they complement each other:

From Wei-Jie Fu



FRG is applicable at high μ_B
and well suited for search of CEP.

Fu, Pawłowski and Rennecke, PRD (2020)

AMPT: a **dynamical** non-equilibrium model, can be directly compared with experimental data.

FRG (functional renormalization group): A non-perturbative but **static** QCD approach, consistent with lattice QCD.

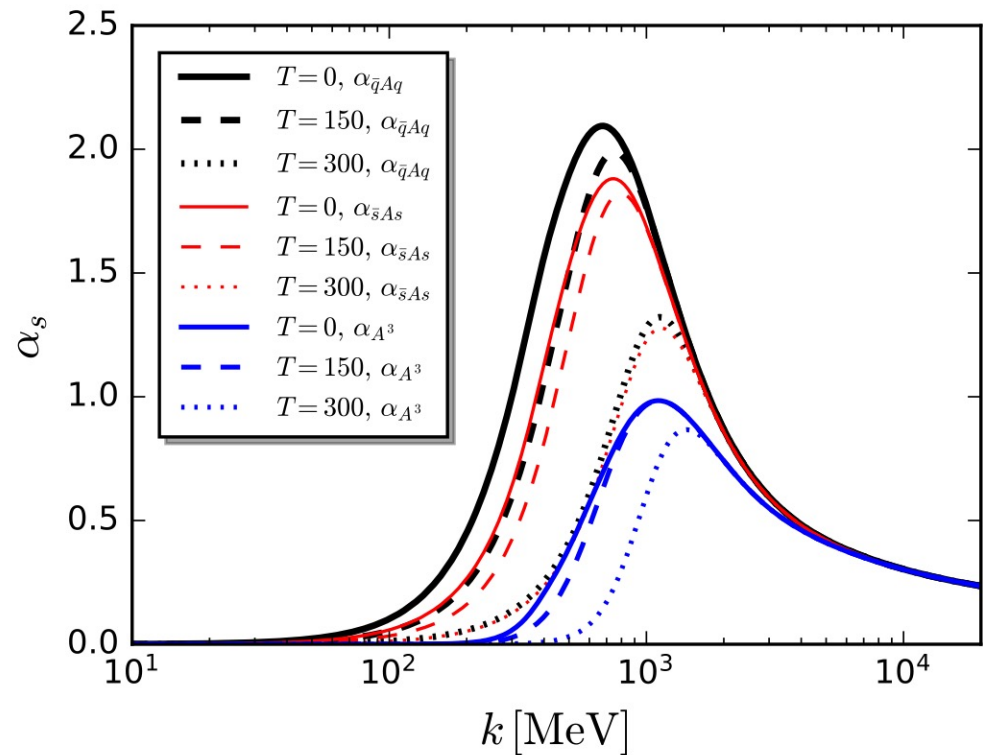
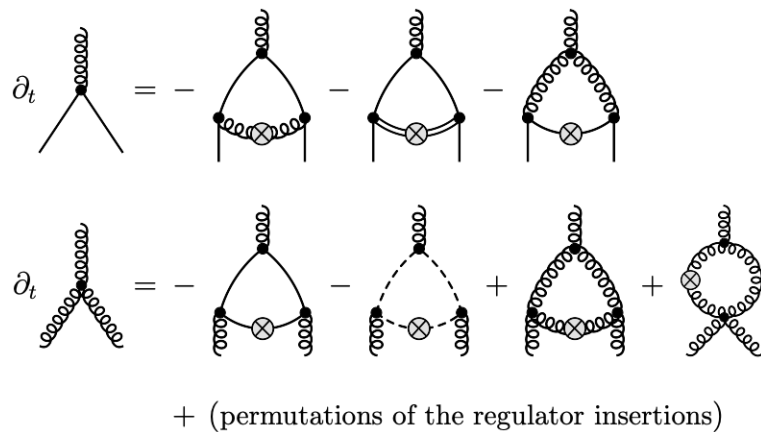
With the coupling of AMPT with FRG:

- first-principle QCD EoS from FRG provides information to improve EoS in AMPT.
- it enables calculation of dynamical evolution of fluctuations and possible effects of CEP.

Future developments for high baryon density physics

From Wei-Jie Fu

QCD strong couplings among quarks and gluons from FRG:



FRG could provide proper treatment of QCD interactions in AMPT

Fu, Pawłowski and Rennecke, PRD (2020)

Summary

- The string melting AMPT can now reasonably describe bulk observables of both large and small systems, including the centrality dependence
- At low energies like the BES, finite nuclear thickness has big effects on *energy density ε & net-baryon density n_B* , and consequently on event trajectories and relation to CEP
- We have incorporated finite nuclear thickness into AMPT to provide a better foundation for further studies at high baryon densities
- We have found a more accurate **new collision scheme** that almost eliminates causality violation in parton cascade and gives correct parton η values
- **Much important work needs to be done, especially on realistic Equation of State of the dense matter including the CEP**