(Parton-Hadron-Quantum-Molecular-Dynamics)

- a novel microscopic transport approach to study heavy ion reactions

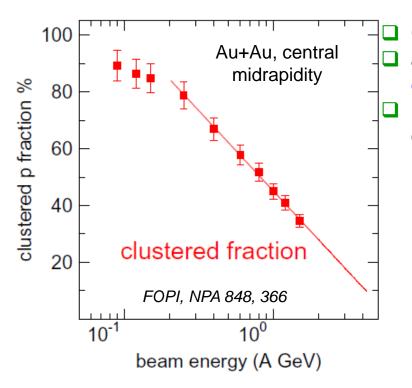
J. Aichelin

(E. Bratkovskaya, M. Winn, A. Le Fèvre, Y. Leifels, V. Kireyev, V. Kolesnikov, V. Voronyuk) arXiv:1907.03860

- ☐ Why a novel approach?
- ☐ Basics of the QMD Transport theory
- ☐ Fragment Formation
- Comparison with existing data
- ☐ Perspectives for BMN/NICA/FAIR/RHIC



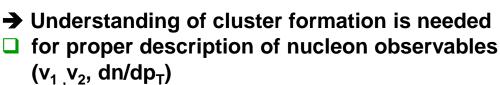
Clusters in HICs



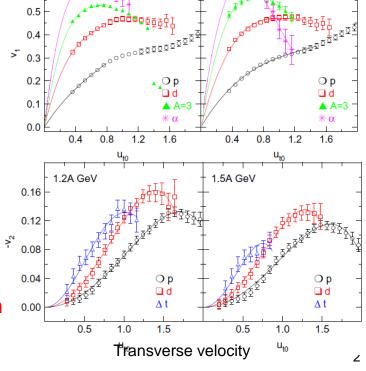
Clusters are very abundant at low energy; at 3 AGeV in central Au+Au collisions ~20% of the baryons are in clusters!

... and baryons in clusters have quite different properties $(v_1, v_2, dn/dp_T)$

Au+Au, semi-central

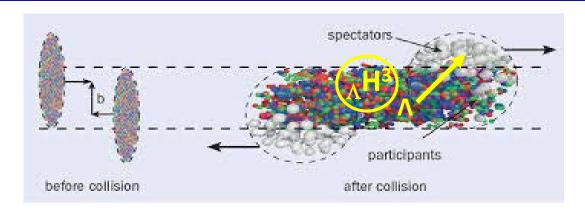


- ☐ to explore new physics opportunities like
- hyper-nucleus formation
- possible signals of the 1st order phase transition
 - cluster formation at midrapidity (RHIC, LHC)



FOPI, NPA 876,1

Why do we study especially hypermatter production?



Access to the nuclear dynamics:

different mechanisms for hypernucleus production vs. rapidity:

- at mid-rapidity : Λ -coalescence hypernuclei test the phase-space distribution of baryons in the expanding participant matter
- at target/projectile y: Λ -absorption by spectators elucidates the physics at the interface between spectator and projectile matter

Hypernuclei as bound objects:

batech

- give access to the third dimension of the nuclear chart (strangeness)
- give information on hyperon-nucleon and hyperon-hyperon interactions
- important e.g. for neutron stars (production of hypermatter at high density and low temperature)
- new field of hyperon spectroscopy

Modelling of fragment and hypernucleus formation

Present microscopic approaches:

- VUU(1985), BUU(1985), (P)HSD(96), SMASH(2016) solve the time evolution of the one-body phase-space density in a mean field → no dynamical fragments
- UrQMD is a n-body model but makes clusterization via coalescence and a statistical fragmentation model
- QMD is a n-body model but is limited to energies < 1.5 AGeV</p>
 - → describes fragments at SIS energies, but conceptually not adapted for NICA/FAIR energies and higher

In order to understand the microscopic origin of cluster formation one needs:

- a realistic model for the dynamical time evolution of HICs
- dynamical modelling of cluster formation based on interactions

Dynamical modelling of cluster formation is a complex task which involves: the fundamental nuclear properties, quantum effects, variable timescales



The goal: to develop a unified n-body microscopic transport approach for the description of heavy-ion dynamics and dynamical cluster formation from low to ultra-relativistic energies

Realization: combined model PHQMD = (PHSD & QMD) & SACA

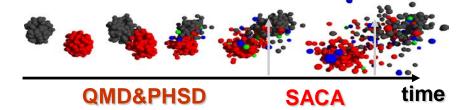
Parton-Hadron-Quantum-Molecular Dynamics

Initialization → propagation of baryons: QMD (Quantum-Molecular Dynamics)

Propagation of partons (quarks, gluons) and mesons
+ collision integral = interactions of hadrons and partons (QGP)
from PHSD (Parton-Hadron-String Dynamics)

Clusters recognition:

SACA (Simulated Annealing Clusterization Algorithm) vs. MST (Minimum Spanning Tree)





Transport eqs. for N-body theories like (PH)QMD,AMD,FMD

Roots in Quantum Mechanics

Remember QM cours when you faced the problem

- we have a Hamiltonian $\hat{H} = -\frac{\hbar^2 \nabla^2}{2m} + V$
- the Schrödinger eq.

$$\hat{H}|\psi_j>=E_j|\psi_j>$$

has no analytical solution

we look for the ground state energy



Walther Ritz

Ritz variational principle:

Assume a trial function $\psi(q,\alpha)$ which contains one adjustable parameter α , which is varied to find the lowest energy expectation value:

$$\frac{d}{d\alpha} < \psi |\hat{H}|\psi > = 0 \rightarrow \alpha_{min}$$

determines α for which $\psi(q,\alpha)$ is closest to the true ground state and $<\psi(\alpha_{min})|\hat{H}|\psi(\alpha_{min})>=E_0(\alpha_{min})$ closest to true ground state E



Extended Ritz variational principle (Koonin, TDHF)

Take trial wavefct with time dependent parameters and solve

$$\delta \int_{t_1}^{t_2} dt < \psi(t) |i\frac{d}{dt} - H|\psi(t) > = 0.$$
(1)

QMD trial wavefet for particle i with p_{oi} (t) and q_{oi} (t)

$$\psi_i(q_i, q_{0i}, p_{0i}) = Cexp[-(q_i - q_{0i} - \frac{p_{0i}}{m}t)^2/4L] \cdot exp[ip_{0i}(q_i - q_{0i}) - i\frac{p_{0i}^2}{2m}t]$$

For N particles:
$$\psi_N = \prod_{i=1}^N \psi_i(q_i, q_{0i}, p_{0i})$$
 QMD

$$\psi_N^F = Slaterdet[\prod^N \psi_i(q_i, q_{0i}, p_{0i})]$$
 AMD/FMD

For the QMD trial wavefct eq. (1) yields

$$\frac{dq}{dt} = \frac{\partial \langle H \rangle}{\partial p} \quad ; \quad \frac{dp}{dt} = -\frac{\partial \langle H \rangle}{\partial q}$$

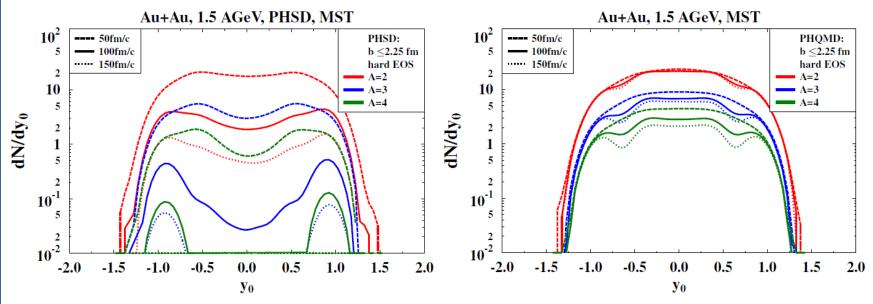
For Gaussian wavefct eq. of motion very similar to Hamilton's eqs. (but only for Gaussians !!)



Time dependence of cluster formation: QMD vs. MF

mean field propagation all two or more body correlation suppressed

QMD propagation correlations present



QMD propagation: number of clusters are stable vs. time (MST finds at 50 fm/c almost the same clusters as at 150fm/c)

MF propagation (per construction not suited for cluster studies):

- -- number of fragments strongly time dependent
- -- fragments disappear with time

ubatech

-- midrapidity fragments disappear early, projectile/target fragments later

→ no common time for coalescence

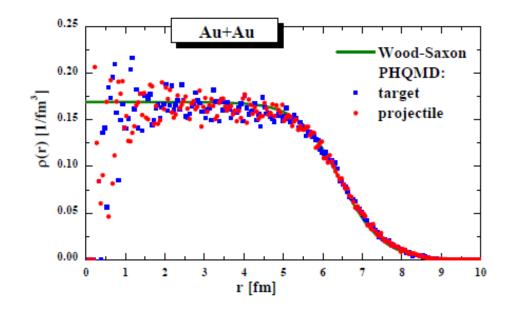


Initial condition:

to describe fragment formation and to guaranty the stability of nuclei

The initial distributions of nucleons in proj and targ has to be carefully modelled:

- Right density distribution
- Right binding energy



local Fermi gas model for the momentum distribution

Potential in PHQMD

above ε =0.5 GeV/fm³ transition to QGP like in PHSD

Below:

Relativistic molecular dynamics (PRC 87, 034912) too time consuming

The potential interaction is most important in two rapidity intervals:

- □ at beam and target rapidity where the fragments are initial final state correlations and created from spectator matter
- at midrapidity where at the late stage the phase space density is sufficiently high that small fragments are formed

In both situations we profit from the fact that the relative momentum between neighboring nucleons is small and therefore nonrelativistic kinematics can be applied. Potential interaction between nucleons

$$V_{i,j} = V(\mathbf{r_i}, \mathbf{r_j}, \mathbf{r_{i0}}, \mathbf{r_{j0}}, t) = V_{\text{Skyrme}} + V_{\text{Coul}}(+V_{mom})$$

$$= \frac{1}{2}t_1\delta(\mathbf{r_i} - \mathbf{r_j}) + \frac{1}{\gamma + 1}t_2\delta(\mathbf{r_i} - \mathbf{r_j})\rho^{\gamma - 1}(\mathbf{r_i}, \mathbf{r_j}, \mathbf{r_{i0}}, \mathbf{r_{j0}}, t)$$

$$+ \frac{1}{2}\frac{Z_iZ_je^2}{|\mathbf{r_i} - \mathbf{r_j}|}(+V_{mom}),$$

$$\langle V(\mathbf{r_i}, t) \rangle = \sum_{j \neq i} \int d^3r d^3r' d^3p d^3p'$$

$$V(\mathbf{r}, \mathbf{r'}, \mathbf{r_i}, \mathbf{r_j}) f(\mathbf{r}, \mathbf{p}, \mathbf{r_i}, \mathbf{p_i}, t) f(\mathbf{r'}, \mathbf{p'}, \mathbf{r_j}, \mathbf{p_j}, t)$$

$$\langle V_i^{Skyrme}(\mathbf{r_i}, t) \rangle = \alpha \left(\frac{\rho_{int}(\mathbf{r_i}, t)}{\rho_0} \right) + \beta \left(\frac{\rho_{int}(\mathbf{r_i}, t)}{\rho_0} \right)^{\gamma}$$

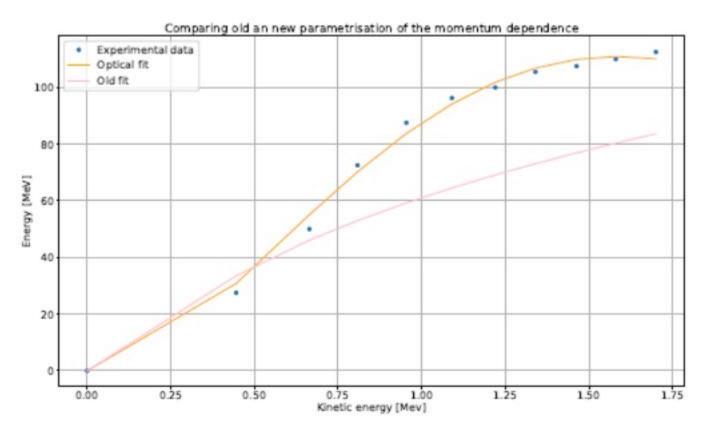
To describe the potential interactions in the spectator matter we transfer the Lorentz-contracted nuclei back into the projectile and target rest frame, neglecting the small time differences

$$\rho_{int}(\mathbf{r_i}, t) \rightarrow C \sum_{j} \left(\frac{4}{\pi L}\right)^{3/2} e^{-\frac{4}{L}(\mathbf{r_i^T}(t) - \mathbf{r_j^T}(t))^2} \cdot e^{-\frac{4\gamma_{cm}^2}{L}(\mathbf{r_i^L}(t) - \mathbf{r_j^L}(t))^2}.$$

For the midrapidity region $\gamma \rightarrow 1$. and we can apply nonrelativisitic kinematics as well

All elastic and inelastic cross sections from PHSD - therefore at high energy the spectra of produced particles are similar to PHSD results (however initial distribution is different in PHSD and PHQMD)

pA experiments show that the NN potential is not static but becomes increasingly repulsive with beam energy



$$V_{mom}(\mathbf{p_{i0}}, \mathbf{p_{j0}}) = \exp\left(-c\sqrt{(\mathbf{p_{i0}} - \mathbf{p_{j0}})^2}\right) \left(a(\mathbf{p_{i0}} - \mathbf{p_{j0}})^2 + b(\mathbf{p_{i0}} - \mathbf{p_{j0}})^4\right) \frac{\rho}{\rho_0}$$

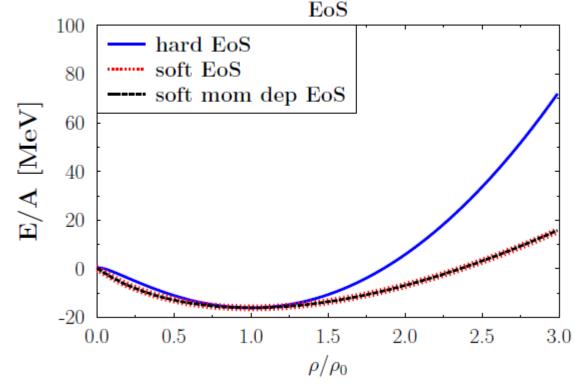
with the parameters

- $a = 236.326(MeV)^{-1}$ $b = -20.7304(MeV)^{-2}$
- $c = 0.901519 \text{MeV}^{-1}$



How to fix the strength of the potential?

In infinite matter a potential corresponds to an equation of state (EoS)



Equation of state cannot be calculated:

Brückner G-matrix is a low density expansion:

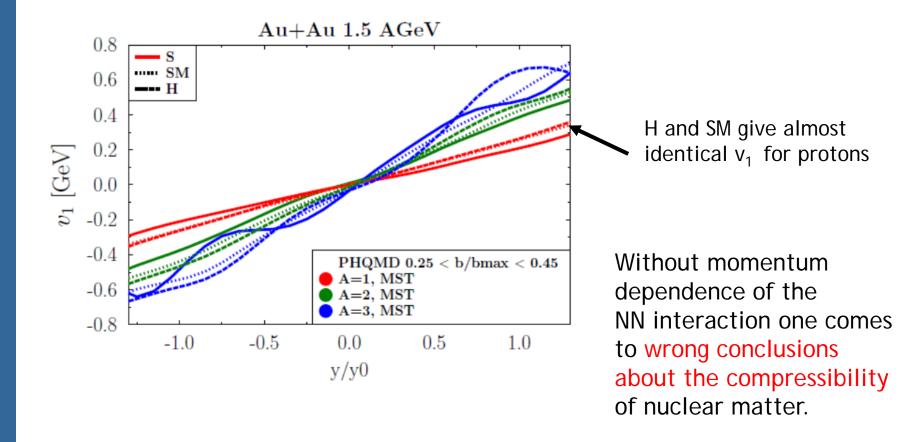
Expansion parameter: $a \cdot k_F$ a=hard core range (.6 fm)

$$k_F = p_F / hbar = 1.28 (\rho/\rho_0)^{1/3} \frac{1}{fm}$$



Why is the momentum dependence important?

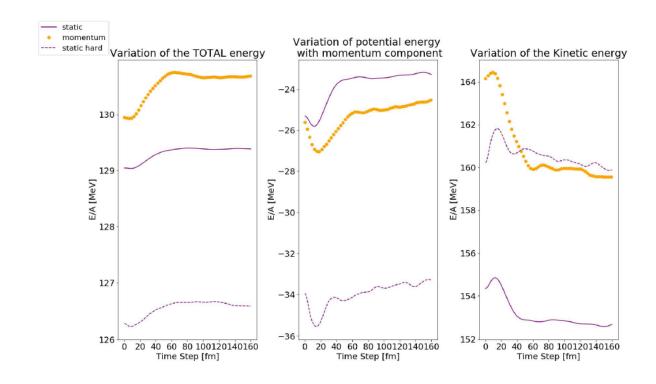
SM and H may give almost identical results. Example: v₁



Also the centrality dependence of v₁ is different for S and SM and SM is closer to H than to S



Energy conservation of the numcerical realization Au+Au 600 AMeV



омона 17

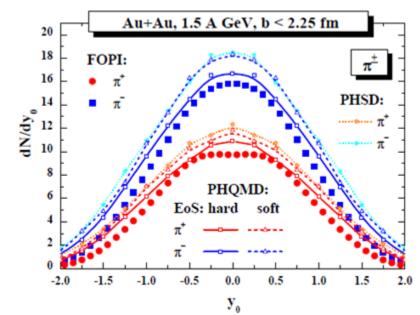
Results

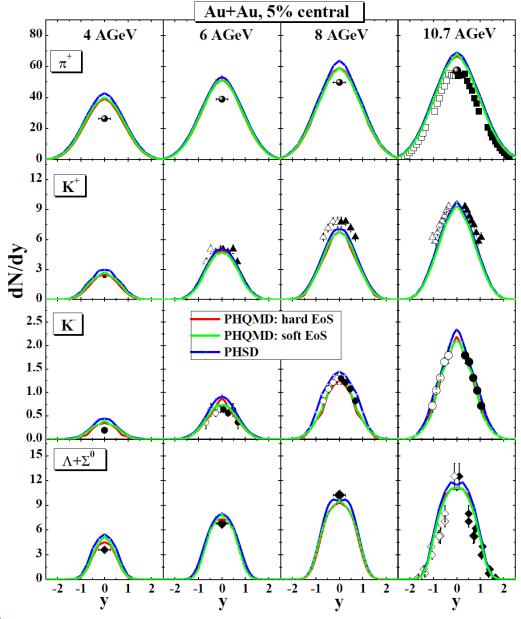
First Results of PHQMD

Produced particles

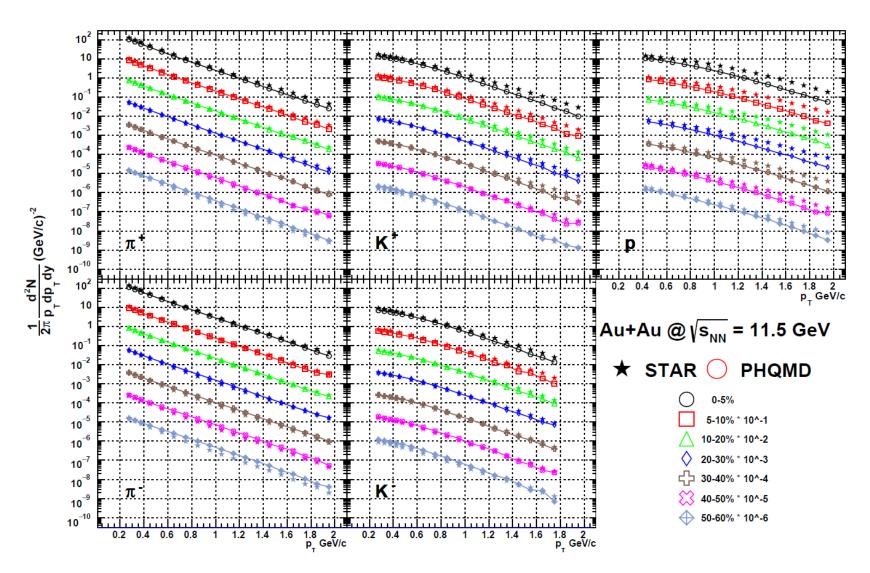
are well reproduced at SIS/NICA/FAIR energies

(dominated by collisions)





.. And also the most recent STAR data at 11.5 AGeV



PHOMD

Methods to identify fragments in theories which propagate nucleons:

Static approaches:

means fragment multiplicity determined at a fixed time point

- -- coalescence (early, assumption: no coll. later)
- -- statistical model (V,T,N) very late $\rho << \rho_0$

Dynamical approaches:

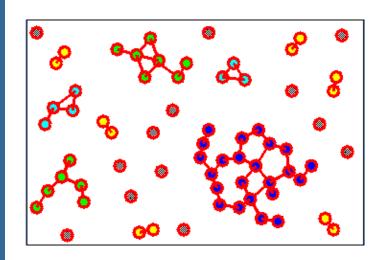
means fragment multiplicity is fct. of time

- -- minimum spanning tree (correlation in coord. space)
- -- simulated annealing (correlation in mom and coord. space)
- -- time dep. perturbation theory using Wigner densities

I. Minimum Spanning Tree (MST) is a cluster recognition method applicable for the (asymptotic) final state where coordinate space correlations may only survive for bound states.

The MST algorithm searches for accumulations of particles in coordinate space:

- 1. Two particles are bound if their distance in coordinate space fulfills $\left| \vec{r}_i \vec{r}_j \right| \leq 2.5 \, fm$
- 2. A particle is bound to a cluster if it is bound with at least one particle of the cluster.



Additional momentum cuts (coalescence) change little: large relative momentum -> finally not at the same position

If we want to identify fragments earlier one has to use momentum space info as well as coordinate space info

Idea by Dorso et al. (Phys.Lett.B301:328,1993):

- a) Take the positions and momenta of all nucleons at time t.
- b) Combine them in all possible ways into all kinds of fragments or leave them as single nucleons
- c) Neglect the interaction among clusters
- d) Choose that configuration which has the highest binding energy

Simulations show: Clusters chosen that way at early times are the preclusters of the final state clusters.

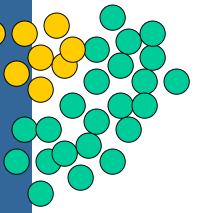
(large persistent coefficient)

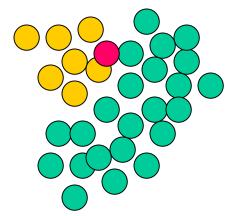
How does this work?

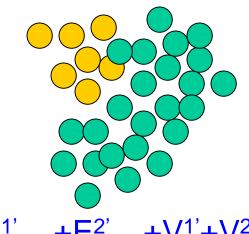
Simulated Annealing Procedure: PLB301:328,1993 later SACA, now FRIGA: Nuovo Cim. C39 (2017) 399

Take randomly 1 nucleon out of a fragment

Add it randomly to another fragment







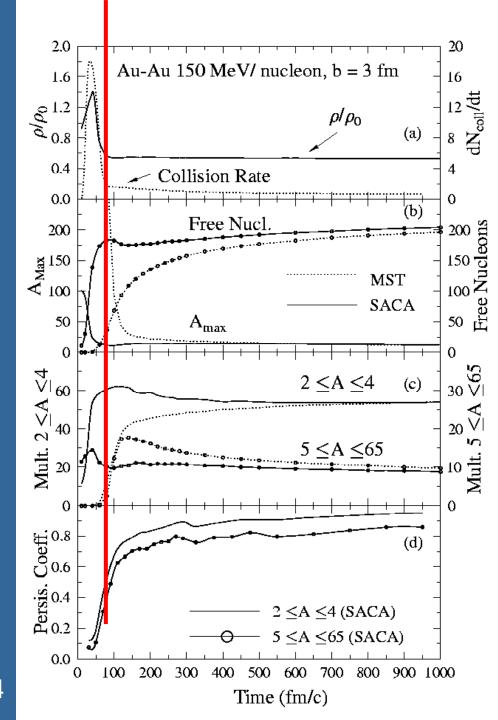
$$E = E_{kin}^{1} + E_{kin}^{2} + V^{1} + V^{2}$$

$$E'=E_{kin}^{1'}+E_{kin}^{2'}+V_{kin}^{1'}+V_{kin}^{2'}$$

If E' < E take the new configuration

If E' > E take the old with a probability depending on E'-E Repeat this procedure very many times

→ Leads automatically to the most bound configuration

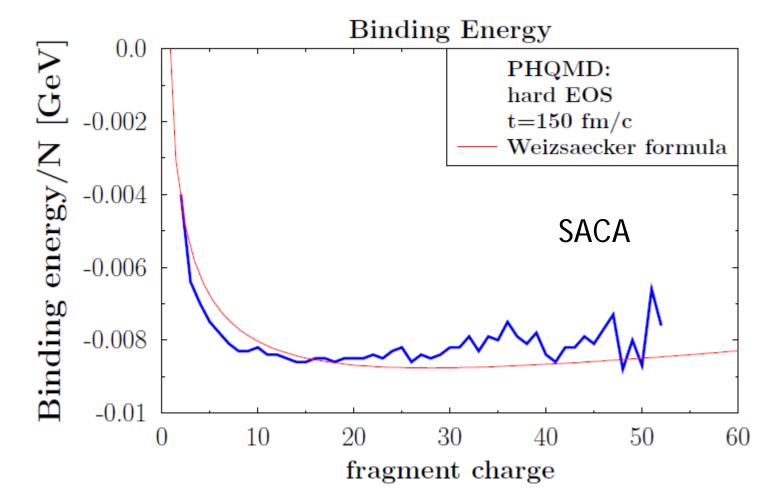


SACA can really identify the fragment pattern very early as compared to the Minimum Spanning Tree (MST) which requires a maximal distance in coordinate space between two nucleons to form a fragment

At1.5t_{pass} Amax and multiplicities of intermediate mass fragments are determined

Fragment formation in PHQMD





There are two kinds of fragments

formed from spectator matter
close to beam and target rapidity
initial-final state correlations
HI reaction makes spectator matter unstable

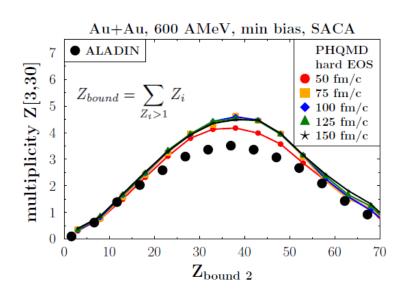
formed from participant matter created during the expansion of the fireball "ice" (E_{bind} ≈8 MeV/N) in "fire"(T≥ 100 MeV) origin not known yet seen from SIS to RHIC (quantum effects may be important)



First Results of PHQMD

Spectator Fragments

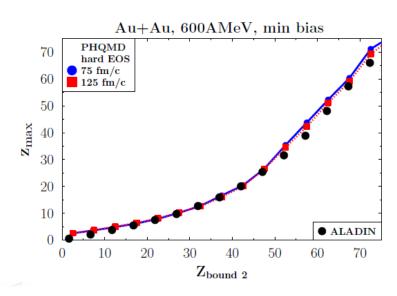
experm. measured up to E_{beam} =1 AGeV (ALADIN)

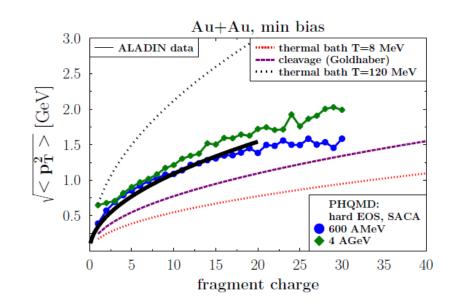


agreement for very complex fragment observables like the

- energy independent "rise and fall"
- ☐ largest fragment (Z_{bound})

rms(p_t) shows \sqrt{Z} dependence

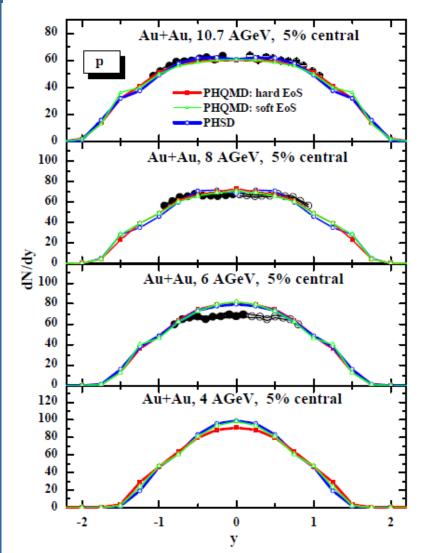


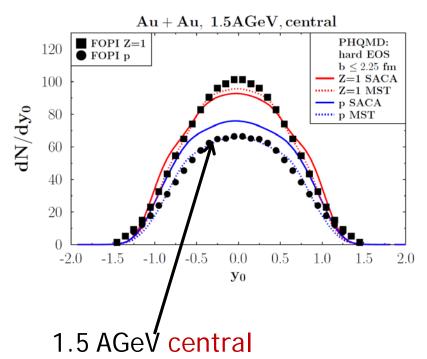


First Results of **PHQMD**

Protons at midrapidity well described

midrapidity fragment production increases with decreasing energy

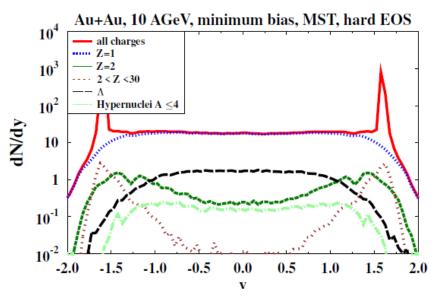


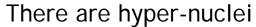


> 30% of protons bound in cluster To improve: better potential for small clusters

.. and what about hyper-nuclei?

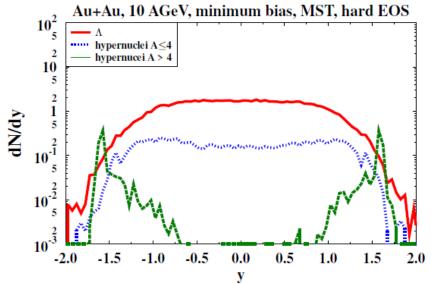
First Results of PHQMD

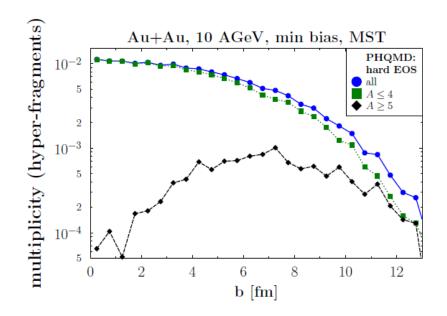




- at midrapidity (A small)
- at beam rapidity (A large)
 few in number but
 more than in other reactions
 to create hyper-nuclei

Central collisions → light hyper-nuclei
Peripheral collisions → heavy hyper-nuclei







Conclusions

We presented a new model, PHQMD, for the NICA/CBM energies which allows - in contrast to all other models - to predict the

dynamical formation of fragments

- allows to understand the proton spectra and the properties of light fragments (dn/dp_Tdy, v₁,v₂, fluctuations)
- allows to understand fragment formation in participant and spectator region
- allows to understand the formation of hypernuclei
- should allow to understand fragment formation at RHIC/LHC

Very good agreement with the presently available fragment data as well as with the AGS/SPS single particle spectra

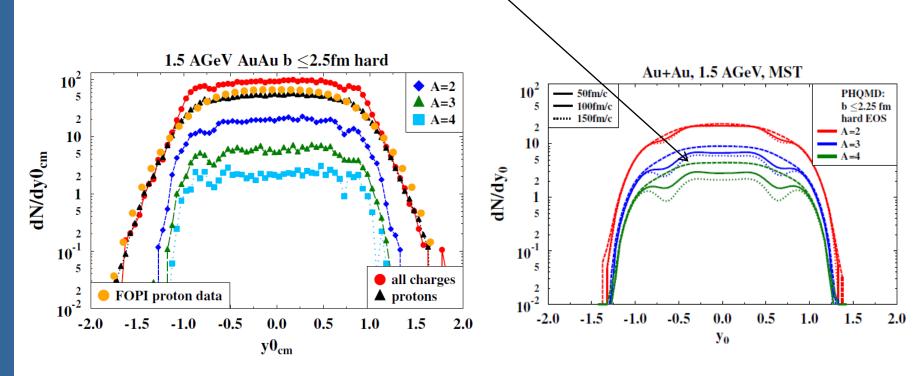
But a lot has still to be done!!



First Results of PHQMD

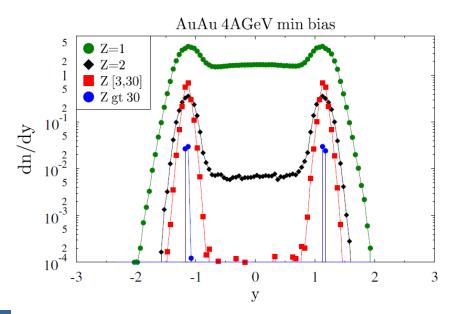
There are all kinds of dynamically produced fragments at midrapidity and they are stable

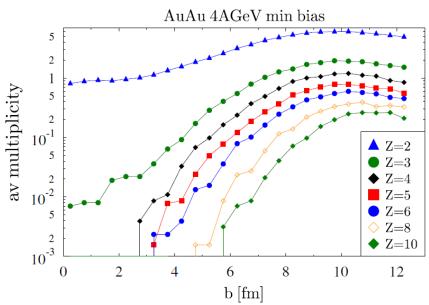
(MST finds at 50fm/c almost the same fragments as at 150fm/c)





First Results of PHQMD





- ☐ Only for most central events fragments do not play a big role
- ☐ Heavy fragments appear only in the residue rapidity range
- Complicated fragment pattern for larger impact parameters
- \square M₇ (b) is different for each fragment charge



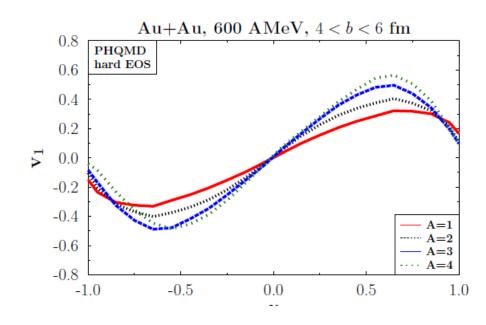
Dynamical variables - v₁

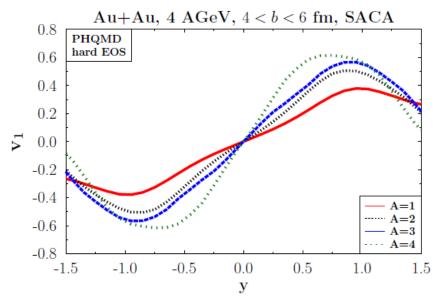
v₁ quite different for nucleons and fragments(as seen in experiments)

nucleons come from participant regions(-> small density gradient)

fragments from interface spectator-participant (strong density gradient)

 v_1 increases with E_{beam} larger density gradient $\rightarrow F_T t_p = p_T$ larger



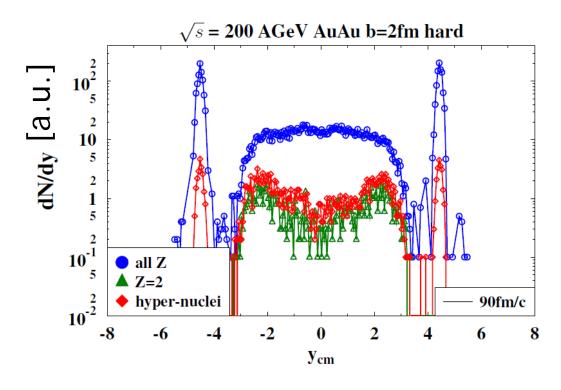


First Results of

PHQMD

At RHIC

hyper-nuclei also from spectator matter Z=2 fragments at midrapidity very preliminary



Back up

Proposals have been made to form clusters in the mean field approaches (which uses test particle method)

using a coalescence description for test particles

$$P_d(\mathbf{r_1},\mathbf{r_2},\mathbf{p_1},\mathbf{p_2},\mathbf{t}) = \underbrace{\rho_d^W(\mathbf{p_1}-\mathbf{p_2},\mathbf{r_1}-\mathbf{r_2})}_{\text{deuteron Wigner density}}$$

On can argue that this is theoretically not consistent because 1 and 2 are test particles, no nucleons. In addition:

- result depends on the number of test particles
- result depends on time t when nucleons coalesce
- □ time is different for different particles: PRC56,2109
- no information about the formation process



Importance of correlations and fluctuations

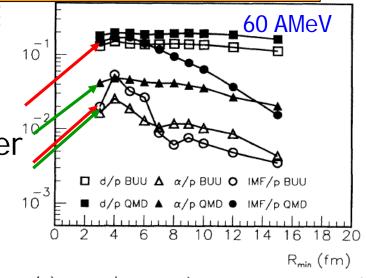
Ratio

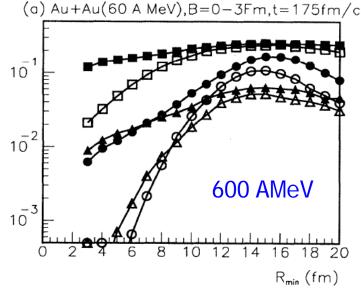
MST analysis with variable R_{min}

Two particles i,j are bound together

if
$$|\vec{r}_i - \vec{r}_j| \le R_{min}$$

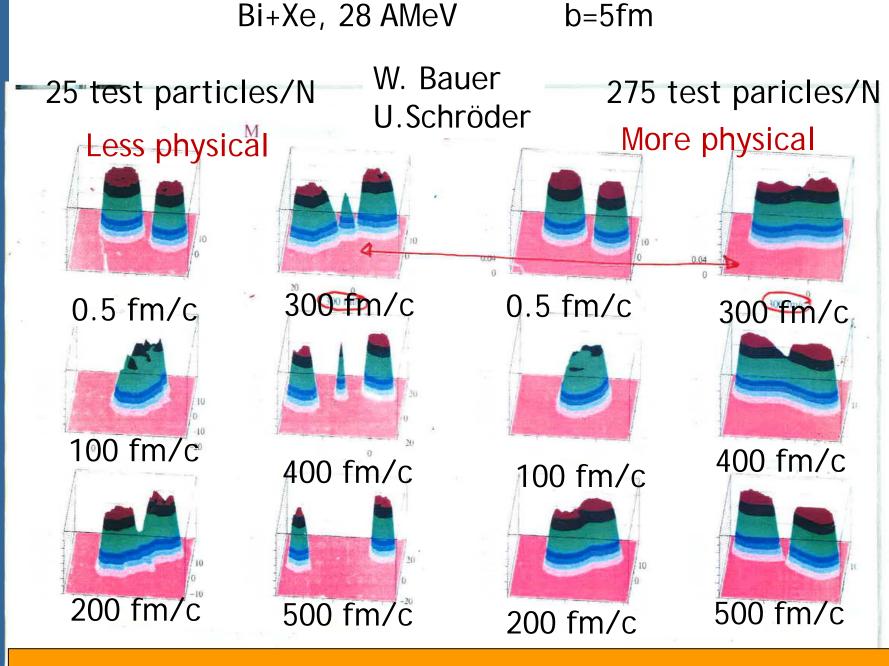
BUU obtained by event mixing of QMD events





(b) Au + Au(600 A MeV), B = 0 - 3Fm, t = 175fm/cGossiaux, Keane (EOS coll) et al

PRC51 (1995) 3357

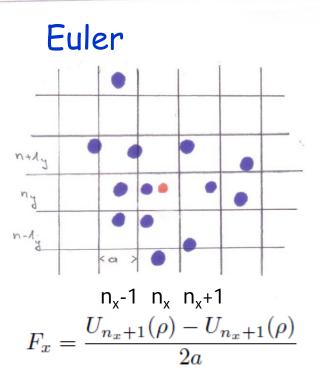


38

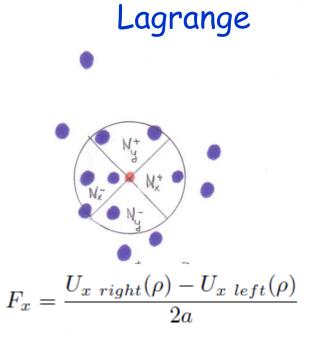
PHQMD

When is N sufficiently large?

One uses delta like forces: $F(r) = \delta(r)$ (Skyrme) but then point-like test particles $f = \sum \delta(r-r_i(t))$ do almost never interact. Solution: one uses grids (and introduces the grid size **a** which plays a similar role as the width in QMD).



Result different if number of test particles is finite (usually N=100)



Average distance between nucleons 2fm. Grid size ≈ 1fm (surface). Therefore very many test particles necessary to avoid numerical fluctuations: 100tp->12 in a cell->30% fluctuation

VUU, BUU, HSD, SMASH solve a Boltzmann type eq.

$$\frac{\partial f}{\partial t} + \frac{\mathbf{p}}{m} \cdot \nabla f + \mathbf{F} \cdot \frac{\partial f}{\partial \mathbf{p}} = \left(\frac{\partial f}{\partial t}\right)_{\text{coll}}$$

Same interaction, not possible classically

$$\left(\frac{\partial f}{\partial t}\right)_{\text{coll}} = \iint gI(g,\Omega)[f(\mathbf{p'}_A,t)f(\mathbf{p'}_B,t) - f(\mathbf{p}_A,t)f(\mathbf{p}_B,t)] d\Omega d^3\mathbf{p}_A d^3\mathbf{p}_B.$$

v · differential cross section

Only the test particle method made it possible to solve the BUU equations in complex situations

Test particle method -> replace integrals by sums (MC) integration

$$f(\mathbf{r}, \mathbf{p}, t) = \sum_{i=1}^{N \to \infty} \delta(\mathbf{r} - \mathbf{r}_i(t)) \ \delta(\mathbf{p} - \mathbf{p}_i(t))$$
 test particle \neq nucleon

If N small unphysical fluctuations

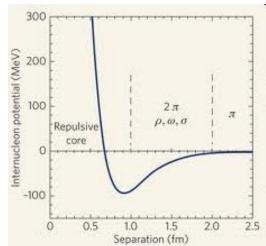
What means N ->∞ in reality?

PHOMD

How does a collision term appear?

The Hamiltonian (Schrödinger and Boltzmann eq.) contains V = NN potential

The NN potential has a hard core, would make transport calculations very unrealistic (Bodmer 75) (independent of the beam energy the participants would thermalize like in a cascade calculation with



would thermalize like In a cascade calculation without Pauli blocking)

Solution (taken over from TDHF):

Replace the NN potential V_{NN} by the solution of the Bethe-Salpeter eq. in T-matrix approach (Brueckner)

$$T = V + V T$$

$$T_{\alpha}(E;q,q') = V_{\alpha}(q,q') + \int k^2 dk \ V_{\alpha}(q,k) \ G_{Q\overline{Q}}^0(E,k) \ T_{\alpha}(E;k,q')$$



$$T_{\alpha}(E;q,q') = V_{\alpha}(q,q') + \int k^2 dk \ V_{\alpha}(q,k) \ G_{Q\overline{Q}}^0(E,k) \ T_{\alpha}(E;k,q')$$

Consequences:

 V_{NN} is real \Rightarrow T is complex = ReT + i Im T

corresponds to V_{NN} σ_{elast} in Hamiltonian collisions
(Skyrme) done identically
BUU (test-particles)
and QMD (particles)

To this one adds inelastic collisions (BUU, HSD, SMASH and QMD - the same way)!

→ Therefore in BUU and QMD the spectra of produced particles are (almost) identical (intensively checked in the past)



- take a small number of test particles (N₁):
 - mathematically this is then not a correct solution of the differential (BUU) equation
 - in practise problems with energy and momentum conserv.
 - assumes, relations between physical (σ, T, ρ) and mathematical fluctuations $(1=\frac{\Gamma}{N})$ which are difficult to justify
- add a fluctuating force to the BUU equation Colonna, Suraud, Ayik......
 - mathematically correct
 - difficult to determine these fluctuations size in Δr and Δp , dependence of T, ρ , (as effectively in QMD)..???
- move in BUU several testparticles simultaneously (Bertsch..)
 - how many and which ones?
 - in which way?

Question: Why not start directly from a N-body theory where fluctuations are (better) under control? (Width L fixed by nucl. density profile etc.)



How to determine the width L?

- surface of the nucleus -> L not too large
- correlations of the relative 2-part. wavefct in a nucleus (healing distance) ≈ 2fm
- range of nuclear potential ≈ 2 fm

$$L = 4.33 \text{ fm}^2$$

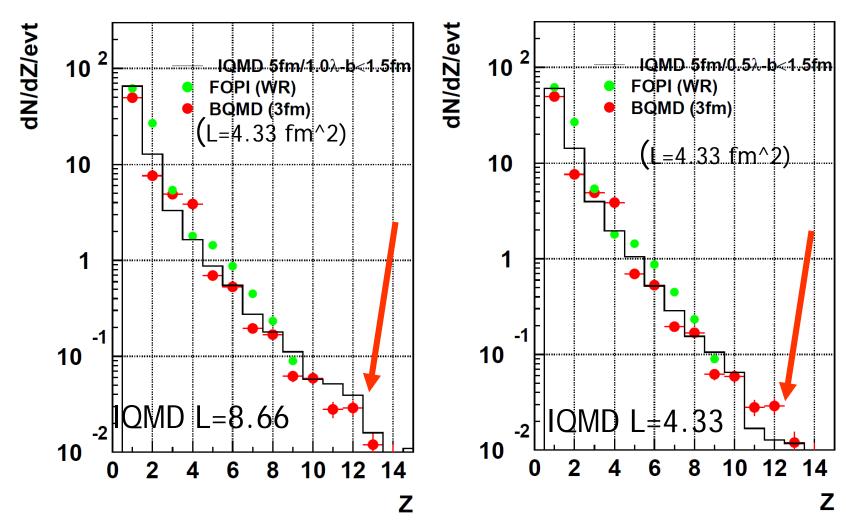
Where L shows up in the observables?

- initially the average over many simulations gives the same $\rho(r)$ as BUU ''d³pf (r;p;t) but the density in each simulation fluctuates around $\rho(r)$ Initial state fluctuations depend on L
- L determines the local density change if a nucleons is kicked out by a hard collision (spectator fragmentation)
 L influences spectator fragmentation
- L plays also a role when fragments are formed from prefr.

 in participant fragmentation (via binding energies)

Influence of L on fragment yield (Y. Leifels)





There are differences but they are modest



Modeling of fragment and hypernucleus formation

The goal: Dynamical modeling of cluster formation by a combined model PHQMD = (QMD & PHSD) & SACA (FRIGA)

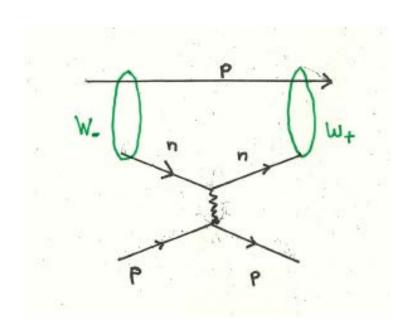
- Parton-Hadron-Quantum-Molecular-Dynamics a non-equilibrium microscopic transport model which describes n-body dynamics based on QMD propagation with collision integrals from PHSD (Parton-Hadron-String Dynamics) and cluster formation by the SACA model or by the Minimum Spanning Tree model (MST).
- MST can determine clusters only at the end of the reaction.
- Simulated Annealing Clusterization Algorithm cluster selection according to the largest binding energy (extension of the SACA model -> FRIGA which includes hypernuclei). FRIGA allows to identity fragments very early during the reaction.



time

PHOMD

III. Wigner density formalism (Remler (NPA 402, 596))



d-wave function

$$\Psi_d(\mathbf{r},\mathbf{R}) \propto exp^{-(\mathbf{r}-\mathbf{r_0})^2L} exp^{-(\mathbf{R}-\mathbf{R_0})^2L/4}$$

d-Wigner density

$$\rho_d^W(\mathbf{r},\mathbf{p}) \propto exp^{-(\mathbf{r}-\mathbf{r_0})^2L} exp^{-(\mathbf{p}-\mathbf{p_0})^2/L\hbar}$$

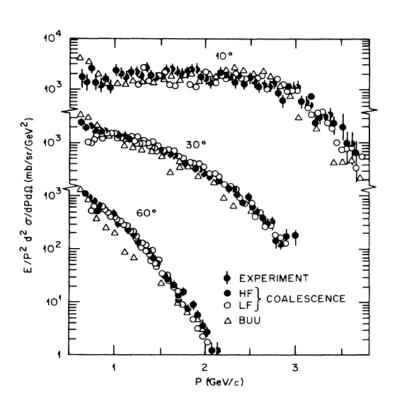
Yields for the rate

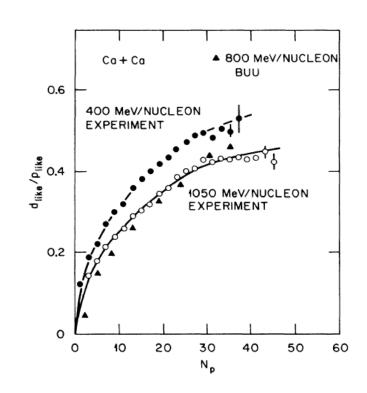
$$\Gamma(t) \ = \underbrace{\sum_{i=1,2} \sum_{j \geq 3}}_{coll \ between \ n \ or \ p \ and \ rest} \delta(t-t_{ij}(\nu)) \overbrace{\int \prod_{i} \frac{d^{3} p_{i} d^{3} x_{i}}{h^{3}} \underbrace{\rho_{d}^{W}(\mathbf{p_{1}, x_{1}, p_{2}, x_{2}})}_{Wigner \ density \ of \ d} [\rho_{N}^{W}(t+\epsilon) - \rho_{N}^{W}(t-\epsilon)] }$$

$$= \underbrace{\sum_{i=1,2} \sum_{j \geq 3}}_{coll \ between \ n \ or \ p \ and \ rest} \delta(t-t_{ij}(\nu)) [\rho_{d}^{W}(t+\epsilon) - \rho_{d}^{W}(t-\epsilon)] }$$

Easy to apply at lower energies

Ca+Ca 800 AMeV (PRC35,1291)





At higher energies: role of resonances? In PHQMD under construction



QMD (like AMD and FMD) are true N-body theories.

N-body theory: Describe the exact time evolution of a system of N particles. All correlations of the system are correctly described and fluctuations correctly propagated.

Roots in classical physics:

A look into textbooks on classical mechanics: If one has a given Hamiltonian

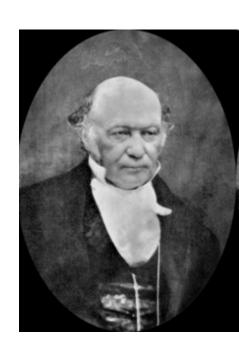
$$H(\mathbf{r}_1,..,\mathbf{r}_N,..,\mathbf{p}_1,..,\mathbf{p}_N,t)$$

$$\frac{d\mathbf{r}_i}{dt} = \frac{\partial H}{\partial \mathbf{p}_i}; \quad \frac{d\mathbf{p}_i}{dt} = -\frac{\partial H}{\partial \mathbf{r}_i}$$

For a given initial condition

$$\mathbf{r}_1(t=0), ..., \mathbf{r}_N(t=0), \mathbf{p}_1(t=0), ..., \mathbf{p}_N(t=0)$$

the positions and momenta of all particles are predictible for all times.



William Hamilton

$m_{\rm r}^{-1} dN/(dm_{\rm r} dy) [({ m GeV})^{-2}]$ 10^2 10^{1} 10° 10⁻¹ K-*0.1 10-2 10⁴ 10^3 10^2 10¹ 10^{0} 10-1 K*0.1 10⁻² 10⁴ 10^3 10^2 10^{1} 10^{0} 10⁻¹ 10-2 K-*0.1 0.2 0.0

 10^4

 10^3 10^2 10^{1}

 10^{0}

10-1

10⁻²

10⁴

 10^3

PHQMD

Au+Au, 11 AGeV, 5%, midrapidity

Au+Au, 8 AGeV, 5%, midrapidity

Au+Au, 6 AGeV, 5%, midrapidity

Au+Au, 4 AGeV, 5%, midrapidity

0.6

 $m_{T}^{-}m_{0}^{-}$ [GeV]

0.4

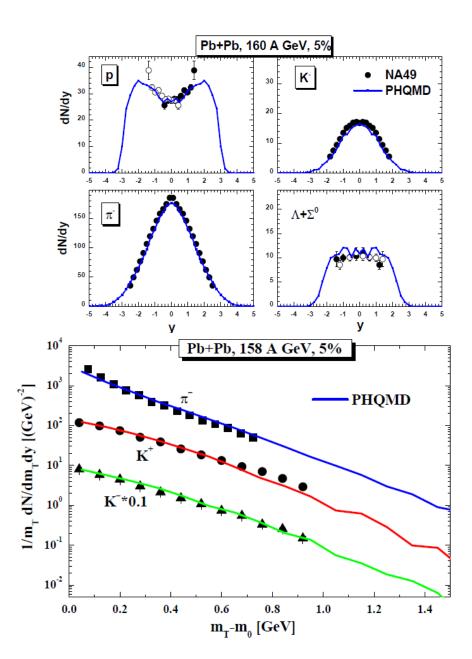
▲ E866/E917

E895

0.8

1.0

As well as at SPS energies





Parton-Hadron-String-Dynamics (PHSD)



PHSD is a non-equilibrium microscopic transport approach for the description of strongly-interacting hadronic and partonic matter created in heavy-ion collisions

Initial A+A collision

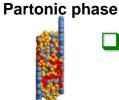
Dynamics: based on the solution of generalized off-shell transport equations derived from Kadanoff-Baym many-body theory



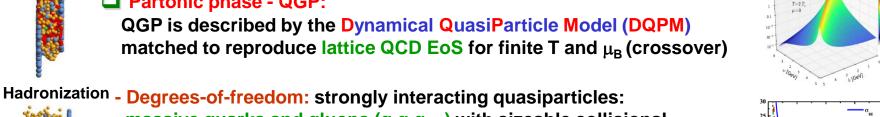
Initial A+A collisions:

N+N → string formation → decay to pre-hadrons + leading hadrons

 \square Formation of QGP stage if local ε > ε_{critical}: dissolution of pre-hadrons → partons



Partonic phase - QGP:





massive quarks and gluons (g,q,q_{bar}) with sizeable collisional widths in a self-generated mean-field potential

- Interactions: (quasi-)elastic and inelastic collisions of partons



Hadronization to colorless off-shell mesons and baryons: Strict 4-momentum and quantum number conservation



