







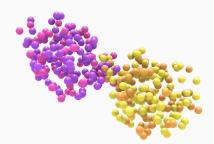


# PHQMD: Effects of EOS: Flow, Clustering

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#### **PHQMD**

PHQMD(Parton Hadron Quantum Molecular Dynamics)<sup>1</sup> is a simulation code that is based on PHSD(Parton Hadron String Dynamics)<sup>2</sup> (quasi particle description of QGP of PHSD)but uses N-body dynamics for propagation of nucleons

### **Quantum Molecular Dynamics**

Using n-body theory to track nucleon by nucleon the collisions, and calculate the two body interactions between all of the nucleons.

- IQMD<sup>3</sup>(Limited to lower energies 1.5 GeV/A)
- UrQMD<sup>4</sup>(Relativistic, used for coalescence, no potential)

<sup>&</sup>lt;sup>1</sup>J. Aichelin and, E. Bratkovskaya et al. *PR C101, 044905* 

<sup>&</sup>lt;sup>2</sup>W. Cassing and E. Bratkovskaya PR C78,034919

<sup>&</sup>lt;sup>3</sup>C. Hartnack et al. Eur. Phys. J. A1:151

<sup>&</sup>lt;sup>4</sup>S. A. Bass et al. NP A41:225

#### Why is it needed?

Current models are not well adapted to address cluster formation, and current QMD models are not usable at relativistic energies

#### **Challenges**

Production of clusters with binding energy of  $B_E \approx -8[MeV]$  in environment of the fireball at  $E \approx 100[MeV]$ 

Hyper-clusters production  $^5$ ,  $^6$  is an area of renewed interest, and which is not currently addressed

Transition form low energy, hadron dominated, interactions and high energy, quark and gluon dominated, reactions

Transition regime is characterised by a finite chemical potential thus a finite net baryonic density

<sup>&</sup>lt;sup>5</sup>C. Rappold et al. *PL B747,129* 

<sup>&</sup>lt;sup>6</sup>J. Adam et al *PR C93, 024917* 

#### Future experimental data

New energy range from 2 GeV/A - 100 GeV/A, will be used to investigate the first order phase transition from hadronic to QGP matter, and degrees of freedom of hadronic matter (strangeness) .

- FAIR
- NICA

#### Experimental results so far

- 0.6 [GeV/A] ALADIN <sup>7</sup>
- 1.23 [GeV/A] HADES 8
- 1.5 [GeV/A] FOPI 9

<sup>&</sup>lt;sup>7</sup>A. Schüttauf et al *NP A607,457* 

<sup>&</sup>lt;sup>8</sup>J. Adamczewski-Musch et al NP A982

<sup>&</sup>lt;sup>9</sup>W. Reisdorf et al. NP A876

### Brief PHQMD model overview

Generalised Ritz variational principle

$$\delta \int_{t_1}^{t_2} dt \langle \Psi(t) | i \frac{d}{dt} - H | \Psi(t) \rangle = 0$$
 (1)

Gaussian test wave function which yields the Wigner density:

$$f(\vec{r_i}, \vec{p_i}, \vec{r_{i0}}, \vec{p_{i0}}, t) = \frac{1}{\pi^3 \hbar^3} \exp\left[-\frac{2}{L} (\vec{r_i} - \vec{r_{i0}})^2\right] \exp\left[-2L(\vec{p_i} - \vec{p_{i0}})^2\right]$$
(2)

Classical type equations of motion for expectation value of the Hamiltonian for Gaussian wavefunctions

$$\dot{r}_{i0} = \frac{\partial \langle H \rangle}{\partial p_{i0}} \qquad \dot{p}_{i0} = -\frac{\partial \langle H \rangle}{\partial r_{i0}} \tag{3}$$

### Brief PHQMD model overview

Hamiltonian composed of Kinetic term and Two body potentials

$$\langle H \rangle = \sum_{i}^{N} \langle H_{i} \rangle = \sum_{i}^{N} \left( \langle T_{i} \rangle + \sum_{i \neq j}^{N} \langle V_{i,j} \rangle \right)$$
 (4)

Potentials ( Coulomb , Asymmetry , Skyrme and Momentum dependent Skyrme)

$$\langle V \rangle = \langle V_c \rangle (\vec{r}) + \langle V_a \rangle (\vec{r}) \begin{cases} \langle V_s \rangle (\rho_{int}) \text{ Static} \\ \langle V_s \rangle (\rho_{int}) + \langle U_{opt} \rangle (\Delta \vec{p}^2) \text{ Momentum dependent} \end{cases}$$
(5)

Interaction density

$$\rho_{int}(i,t) = C \sum_{i,i \neq i}^{N} \exp\left[-\frac{1}{L}(\vec{r_i} - \vec{r_j})^2\right]$$
 (6)

## E.o.S parametrisation

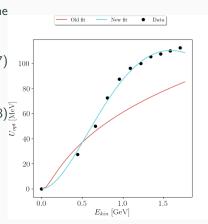
Momentum dependent Skyrme parametrisation

$$\langle V_{tot} \rangle = \langle V_s \rangle \left( \rho_{int} \right) \qquad (7) \qquad 100 - 4$$

$$+ \langle U_{opt} \rangle \left( \Delta p^2 \right) \qquad 80 - 4$$

$$U_{opt}(\Delta p^2) = \exp \left[ -c\sqrt{\Delta p} \right] \qquad (8) \stackrel{\text{S}}{\underset{\text{S}}{\overset{\text{S}}{=}}} \qquad 60 - 4$$

$$\left( a\Delta p + b\Delta p^2 \right) \rho(\vec{r}) \qquad 20 - 4$$



<sup>&</sup>lt;sup>1</sup>S. Hama, et al., Phys. Rev. C 41, 2737

Optical potential fit to expermimental  ${\rm data}^{10}$ 

## E.o.S parametrisation

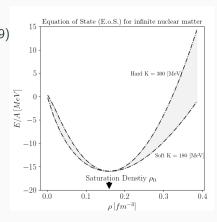
### Static Skyrme parametrisation

$$\left\langle V_{s}\right
angle \left( 
ho_{int}
ight) \propto lpha 
ho + eta 
ho^{\gamma}$$

Constraints for the potentials

$$\begin{cases} \rho_0 = 0.1695 [fm^{-3}] \\ E_0 = -16 [MeV] \\ K_0 = 200 \text{ or } 380 [MeV] \\ (\text{Soft/Hard}) \end{cases}$$

$$K_0 = 9 \frac{\partial^2 E/A}{\partial \rho^2} \Big|_{\rho = \rho_0}$$



Equations of state parametrised for soft and hard nuclear matter compressibility

### E.o.S parametrisation

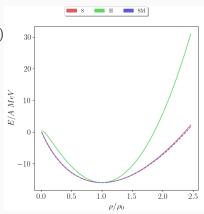
### Static Skyrme parametrisation

$$\langle V_s \rangle (\rho_{int}) \propto \alpha \rho + \beta \rho^{\gamma}$$
 (10)

Constraints for the potentials

$$\begin{cases} \rho_0 = 0.1695 [fm^{-3}] \\ E_0 = -16 [MeV] \\ K_0 = 200 \text{ or } 380 [MeV] \\ (\text{Soft/Hard}) \end{cases}$$

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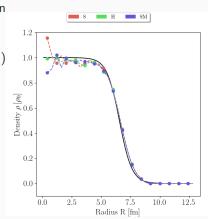
Equations of state parametrised for soft and hard nuclear matter compressibility

### **Nucleon distribution**

Wood-Saxon distribution in position space of nucleons

$$\rho^{WS}(r) = \frac{\rho_0}{1 + \exp\left[\frac{r - R_A}{a}\right]}$$
 (11)

$$\begin{cases} R_A = r_0 A^{1/3} \\ r_0 = 1.125[fm] \\ a = 0.535[fm] \end{cases}$$



Initial Wood-Saxon position distribution for the three E.o.S.

## **Binding energy**

Total binding energy of our nucleus  $( < B_F > = -10[MeV] )$ 

$$B_{E} > = -10[\textit{IMeV}]$$

$$B_{E} = E_{kin}(\vec{p}) + E_{c}(\vec{r}) + E_{a}(\vec{r})$$

$$\begin{cases} E_{s}(\vec{r}) \\ E_{s}(\vec{r}) + E_{m}(\vec{r}, \Delta p) \end{cases}$$

$$(12) \begin{cases} E_{s}(\vec{r}) \\ E_{s}(\vec{r}) \\ E_{s}(\vec{r}) \end{cases}$$

$$(13) \begin{cases} E_{s}(\vec{r}) \\ E_{s}(\vec{r}) \\ E_{s}(\vec{r}) \end{cases}$$

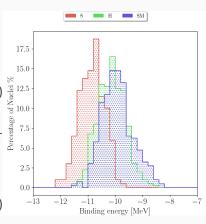
$$(14) \begin{cases} E_{s}(\vec{r}) \\ E_{s}(\vec{r}) \\ E_{s}(\vec{r}) \end{cases}$$

$$(15) \begin{cases} E_{s}(\vec{r}) \\ E_{s}(\vec{r}) \\ E_{s}(\vec{r}) \end{cases}$$

$$(16) \begin{cases} E_{s}(\vec{r}) \\ E_{s}(\vec{r}) \\ E_{s}(\vec{r}) \\ E_{s}(\vec{r}) \end{cases}$$

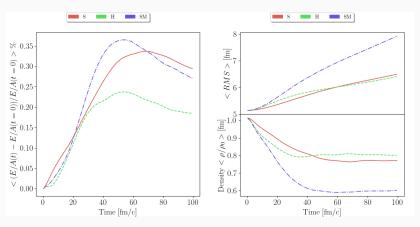
Constraint on momentum distribution

$$0 \le \sqrt{m^2 + \vec{p}_i^2} - m \le -V(\rho_{int})$$
(13)



Initial binding energy of the nuclei

## **Propagation**



Conservation of the total system energyEvolution of the average radial distance of the nucleons, and the average density of the nuclei

#### Cluster formation

#### SACA<sup>11</sup> and MST<sup>12</sup>

To form the cluster we use two methods

- MST: simple form of spanning tree to create clusters based on distance from other nucleons
- SACA: complex simulated annealing of all possible cluster patterns of the nucleons to find the lowest sum of binding energies of all clusters

All the following results are for the SACA method

<sup>&</sup>lt;sup>10</sup>R. K. Puri, C. Hartnack and J. Aichelin *PR C54, R28* 

<sup>&</sup>lt;sup>11</sup> J. Aichelin, Phys. Rept 202, 233

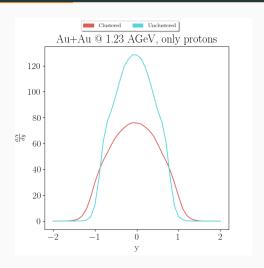
#### Cluster formation

### SACA<sup>13</sup>and MST<sup>14</sup>

To form the cluster we use two methods

- MST :
- SACA:

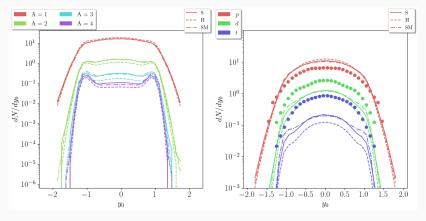
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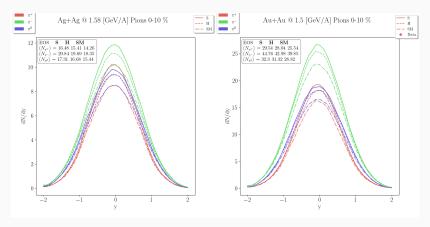
## FOPI results for AuAu E = 1.5 [GeV/A] Multiplicity



Fragment multiplicity

Z=1 isotopes multiplicities for central collisions ( b < 2 [fm])

## Effect of EOS on pion Multiplicity

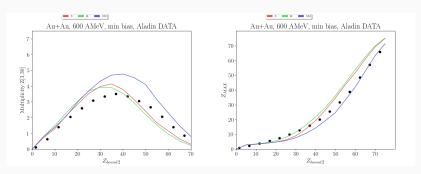


Pion multiplicity for Ag Ag 0-10 % @ 1.58

Pion multiplicity for Au Au 0-10 % @ 1.5

## Aladin results for AuAu E = 0.6 [GeV/A]

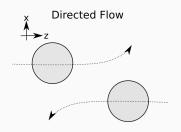
 $Z_{bound2}$  the sum of the charges all fragments with  $Z \geq 2$ , and  $Z_{MAX}$  the charge of the largest cluster

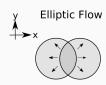


Rise and Fall curve for all three equations of state

Zmax curve for all three equations of state

# **Directed and Elliptic Flow**



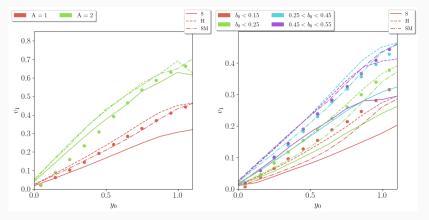


$$E\frac{\mathrm{d}^{3}N}{\mathrm{d}^{3}p} = \frac{1}{2\pi} \frac{\mathrm{d}^{2}N}{p_{T}\mathrm{d}p_{T}\mathrm{d}y}$$
$$\left(1 + 2\sum_{n=1}^{\infty} v_{n} \cos\left[n(\varphi - \Psi_{RP})\right]\right) \quad (14)$$

A Fourier expansions allows us to obtain the coefficients:

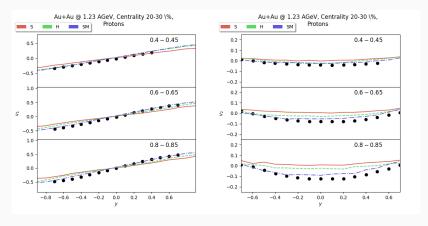
- Directed Flow  $v_1 = \left\langle \frac{p_x}{p_T} \right\rangle$
- Elliptic Flow  $v_2 = \left\langle \left(\frac{p_x}{p_T}\right)^2 \left(\frac{p_y}{p_T}\right)^2 \right\rangle$

## FOPI results for AuAu E = 1.5 [GeV/A] Flow

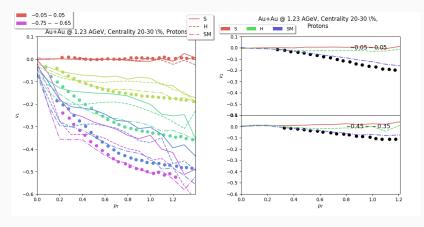


Directed flow for proton  $(v_1)$  and momentum

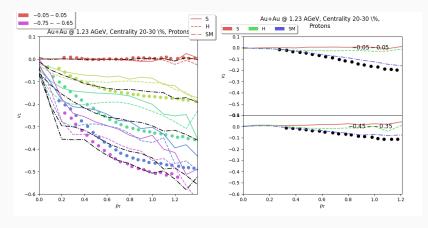
Directed proton flow  $(v_1)$  for 4 different deuteron with a threshold on transverseimpact parameter classes for all protons with a threshold on transverse momentum



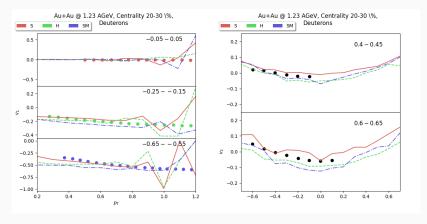
Directed proton flow  $(v_1)$  as a function Elliptical proton flow  $(v_2)$  as a function of rapidity for three  $p_T$  classes compared with the HADES data compared with the HADES data



Directed proton flow  $(v_1)$  as a function Elliptical deuteron flow  $(v_2)$  as a of  $p_T$  for various rapidity classes function of rapidity for two  $p_T$  classes compared with the HADES data



Directed proton flow  $(v_1)$  as a function Elliptical deuteron flow  $(v_2)$  as a of  $p_T$  for various rapidity classes function of rapidity for two  $p_T$  classes compared with the HADES data



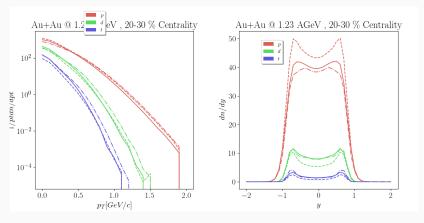
Directed deuteron flow  $(v_1)$  as a Elliptical deuteron flow  $(v_2)$  as a function of  $p_T$  for three various rapidity function of rapidity for two  $p_T$  classes classes compared with the HADES datacompared with the HADES data

#### **Conclusions**

- PHQMD allows use to study cluster formation in a unique and novel way
- It permits us to study the effect of nuclear EOS on a host of observables
- In most cases for the FOPI data the SM EOS give better agreement with the data In some cases the HADES data is better replicated for the SM EOS

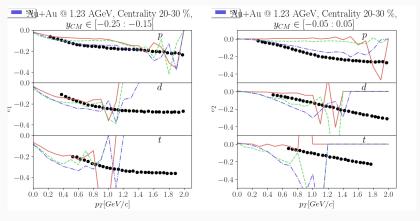
Thank you

# AuAu E = 1.23 [GeV/A] Multiplicity



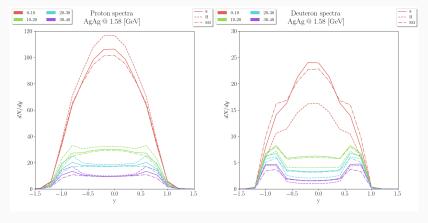
Multiplicity of Z=1 isotopes function of Multiplicity of Z=1 isotopes function of rapidity for Au+AU @ 1.23 AGeV  $p_T$  for Au+AU @ 1.23 AGeV

# AuAu E = 1.23 [GeV/A] Flow



Directed isotope flow of Z=1 isotopes function of  $p_T$  for Au+AU @ 1.23 AGeV

Elliptical Z=1 isotopes flow function of  $p_T$  for Au+AU @ 1.23 AGeV



Proton multiplicity Ag Ag @ 1.58 AGeVDeuteron multiplicity Ag Ag @ 1.58 function rapidity AGeV function rapidity