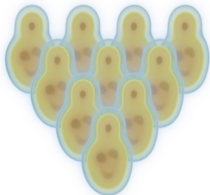


The intersection of heavy ions and nuclear structure

from the neutron skin of ^{208}Pb to the unexpected uses of a nuclear bowling pin

Govert Nijs

October 1, 2024

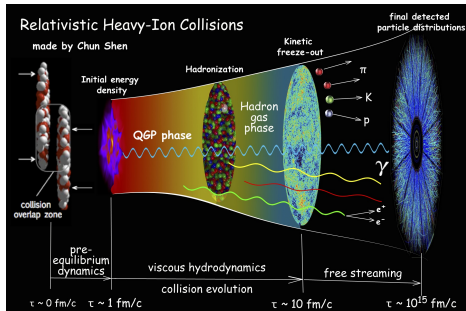


Based on:

- GN, van der Schee, 2312.04623
- Giacalone, GN, van der Schee, 2305.00015
- Giacalone, Bally, GN, Shen, Duguet, Ebran, Elhatisari, Frosini, Lähde, Lee, Lu, Ma, Meißner, Noronha-Hostler, Plumberg, Rodríguez, Roth, van der Schee, Somà, 2402.05995

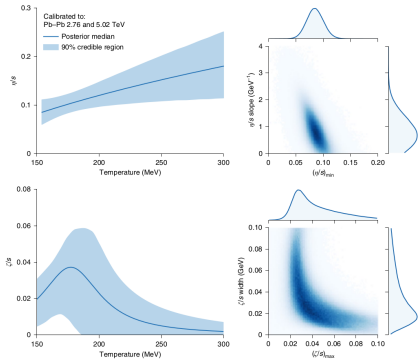
The status of the field

- The general picture of the stages of a heavy ion collision is known.
- Theoretical modelling follows these stages:
 - T_{RENTo} or IP-Glasma for the initial state.
 - Free streaming for the pre-hydrodynamic stage.
 - Viscous hydrodynamics with temperature dependent shear and bulk viscosity.
 - SMASH or UrQMD as a hadronic afterburner.
- Bayesian analysis gives a data-driven approach to understand each stage in more detail.

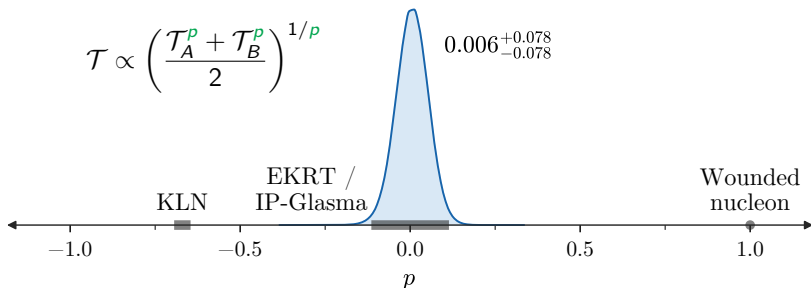


Uses of Bayesian analysis: viscosities

- We know the QGP phase is described by viscous hydrodynamics.
 - We know exactly what the free parameters are, i.e. η/s , ζ/s , ...
- We can use Bayesian analysis to find data-preferred values for these parameters.
- The values of the parameters provide an interface with microscopic theories of the QGP.



Uses of Bayesian analysis: parameterized phenomenology

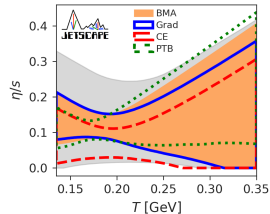
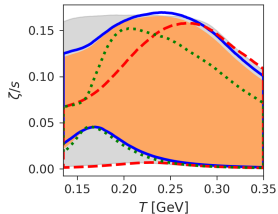


- For the initial state, there is no single widely accepted model.
- With a phenomenological model such as $\mathcal{T}_{\text{RENT}}^o$, aspects of microscopic models can be tested, such as the scaling shown here, parameterized by p .
 - IP-Glasma and EKRT are ruled in.
 - KLN and wounded nucleon are ruled out.



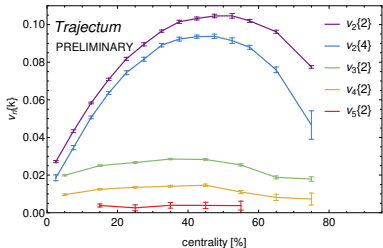
Uses of Bayesian analysis: deciding between models

- One can take this idea a step further, and actually compare different models.
- Here shown are different particlization schemes.
- By taking into account how well each model fits, one can even take a weighted average over models, known as *Bayesian model averaging*.

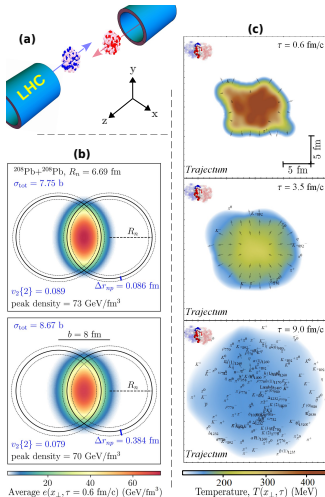


Model used: *Trajectum*

- New heavy ion code developed in Utrecht/MIT/CERN.
 - *Trajectum* is the old Roman name for Utrecht.
- Contains initial stage, hydrodynamics and freeze-out, as well as an analysis suite.
- Easy to use, example parameter files distributed alongside the source code.
- Fast, fully parallelized.
 - Figure (20k oversampled PbPb events at 2.76 TeV) computes on a laptop in 21h.
 - Bayesian analysis requires $\mathcal{O}(1000)$ similar calculations to this one.
- Publicly available at sites.google.com/view/governnijs/trajectum/.



Some simple intuition

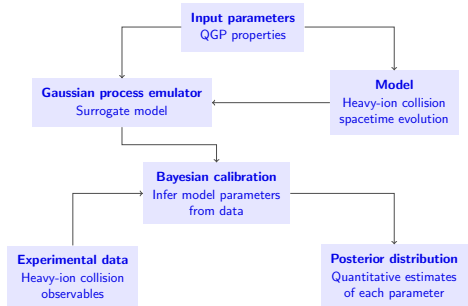


- Model details are not necessary to understand the contents of this talk.
 - Some details are available in the backup.
- Hydrodynamics can be intuitively understood:
 - Pressure gradients drive expansion.
 - Hotter systems expand faster, resulting in more transverse momentum.
 - Spatially anisotropic systems expand preferentially along the short axis, resulting in momentum anisotropy in the final state.



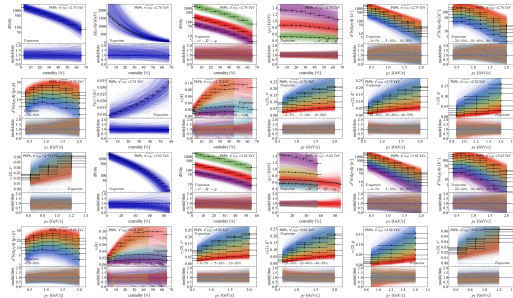
Bayesian analysis workflow

- In principle, Bayesian analysis is simply a fit to data.
- In practice the process is more complicated:
 - Generate a large number of randomly chosen parameter sets called *design points*.
 - Run the model for each one to obtain the prior.
 - Train the emulator.
 - Run the MCMC to obtain the posterior.
- The posterior then is a list of likely parameter sets.



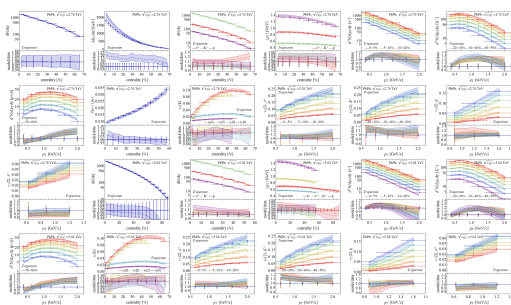
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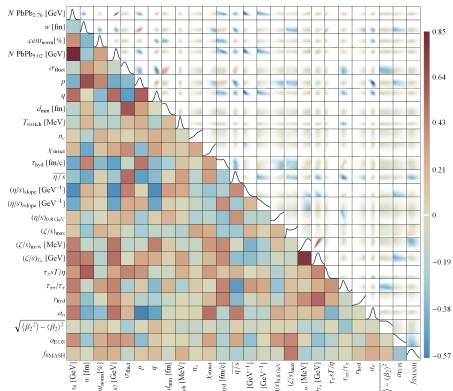
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Data used: 670 individual data points

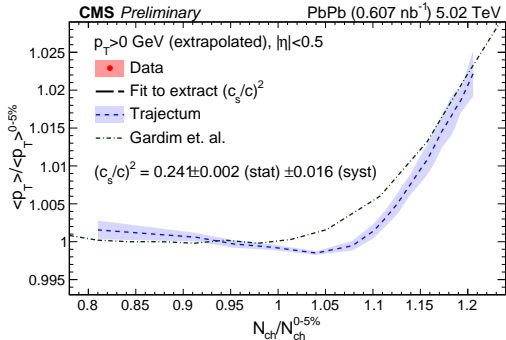
✓: data used
 ⌚: data available
 ✗: data unavailable

	PbPb 2.76 TeV				PbPb 5.02 TeV				pPb 5.02 TeV
	incl.	π^\pm	K^\pm	p	incl.	π^\pm	K^\pm	p	incl.
σ	✗	✗	✗	✗	✓	✗	✗	✗	✓
dN/dy	✓	✓	✓	✓	✓	✓	✓	✓	⌚
$\langle p_T \rangle$	✗	✓	✓	✓	✓	✓	✓	✓	⌚
$dE_T/d\eta$	✓	✗	✗	✗	✗	✗	✗	✗	✗
$\delta p_T / \langle p_T \rangle$	✓	✗	✗	✗	✗	✗	✗	✗	✗
$v_{2,3,4}\{2\}$	✓	⌚	⌚	⌚	✓	⌚	⌚	⌚	⌚
$v_2\{4\}$	✓	✗	✗	✗	✓	⌚	⌚	⌚	⌚
$d^2N/dp_T dy$	✗	✓	✓	✓	✗	✓	✓	✓	✗
$v_2\{2\}(p_T)$	✗	✓	✓	✓	✗	✓	✓	✓	⌚
$v_3\{2\}(p_T)$	✗	✓	⌚	⌚	✗	✓	⌚	⌚	⌚
$NSC(2, 3)$	⌚	✗	✗	✗	✓	✗	✗	✗	⌚
$NSC(2, 4)$	⌚	✗	✗	✗	✓	✗	✗	✗	⌚
$\rho(v_2\{2\}^2, \langle p_T \rangle)$	✗	✗	✗	✗	✓	✗	✗	✗	✗



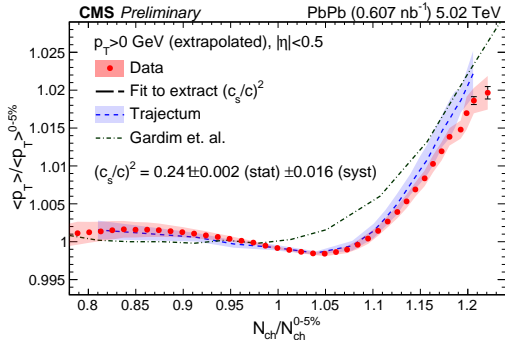
Using the posterior parameter values to make predictions

- The posterior parameter values can be used to make predictions for new observables.
 - When using multiple samples from the posterior, this includes systematic uncertainty from the parameter estimation.
- Here shown is the prediction for ultracentral $\langle p_T \rangle$.



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 - When using multiple samples from the posterior, this includes systematic uncertainty from the parameter estimation.
- Here shown is the prediction for ultracentral $\langle p_T \rangle$.
- Precise agreement between theory and experiment.



Neutron skin

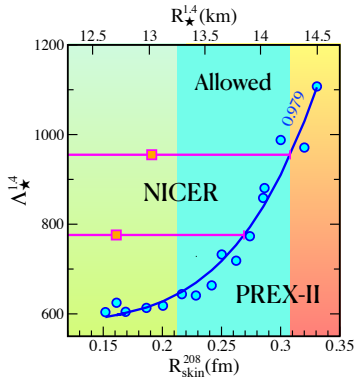
- In a ^{208}Pb nucleus, neutrons sit further from the center than protons.

- This is quantified by the *neutron skin*:

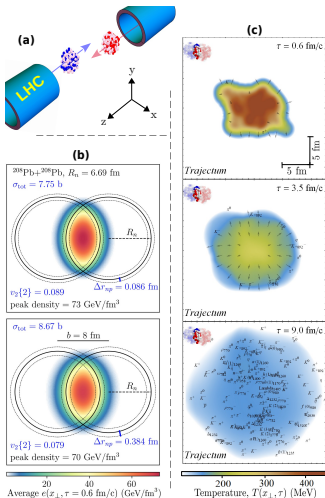
$$\Delta r_{np} = \langle r^2 \rangle_n^{1/2} - \langle r^2 \rangle_p^{1/2},$$

i.e. the *difference* in RMS radii of the neutron and proton distributions.

- Heavy nuclei and neutron stars are sensitive to the same nuclear interactions.
 - A constraint on Δr_{np} translates directly into a constraint on the radius of a $1.4M_{\odot}$ neutron star.
 - We can learn something about the low T , high μ_B region even at LHC energies!



How to measure neutron skin?



- To measure the neutron skin, we need the distributions of protons and neutrons inside the nucleus.
 - The proton distribution distribution is well-known from electron scattering.
- Several different methods are in use for the neutron distribution:
 - Polarized electron scattering off ^{208}Pb (PREX).
 - Photon tomography of ^{197}Au (STAR).
- Heavy ion collisions provide a completely orthogonal method.
 - Sensitive to the total matter distribution inside the nucleus.
 - Purely gluonic measurement.



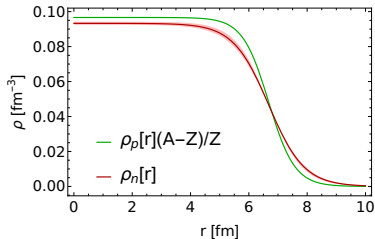
The Woods-Saxon distribution

- Nucleon positions are drawn from a Woods-Saxon distribution:

$$\rho_{WS}(r) \propto \frac{1}{1 + \exp\left(\frac{r-R}{a}\right)}.$$

- We fix R for both protons and neutrons.
- We fix a for protons, while varying a_n as a parameter.
- Neutron skin $\Delta r_{np} = \langle r^2 \rangle_n^{1/2} - \langle r^2 \rangle_p^{1/2}$ strongly depends on a_n :

$$\langle r^2 \rangle_{WS} = \frac{12a^2 \text{Li}_5(-e^{R/a})}{\text{Li}_3(-e^{R/a})}.$$



	proton	neutron
R [fm]	6.68	6.69
a [fm]	0.447	a_n



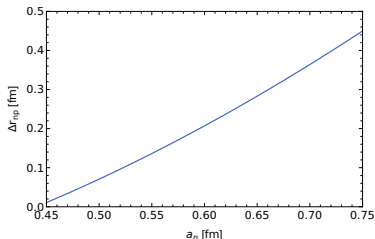
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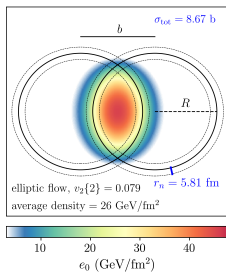
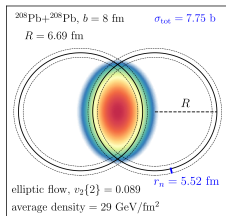


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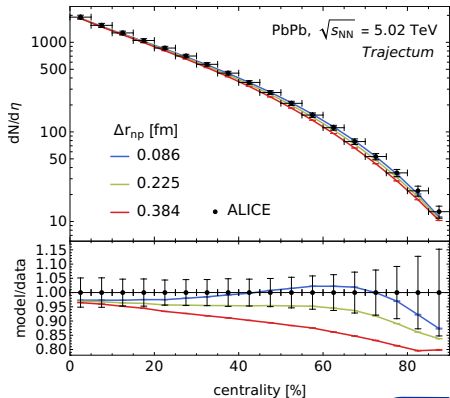
Do we have observables sensitive to a_n ?

- Initial geometry is sensitive to a_n .
Larger nuclei lead to:
 - Larger hadronic PbPb cross-section,
 - Larger initial QGP size,
 - Smaller initial QGP eccentricity.
- Final state observables are in turn sensitive to initial geometry. Larger Δr_{np} leads to:
 - Larger hadronic PbPb cross-section,
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 - Smaller mean transverse momentum,
 - Smaller elliptic flow.



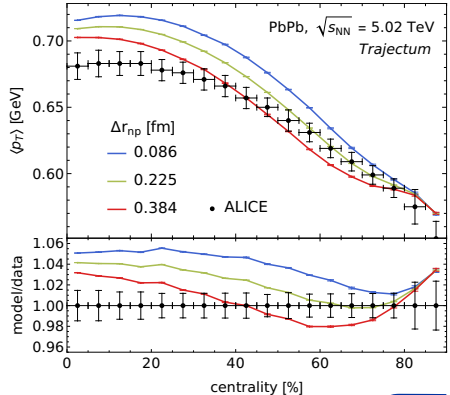
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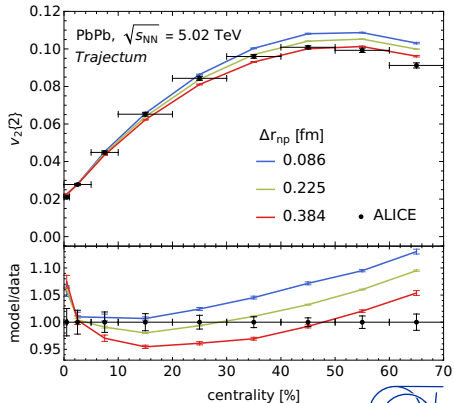
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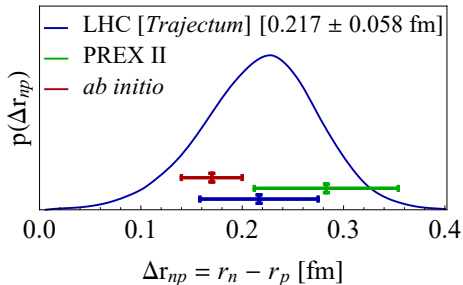
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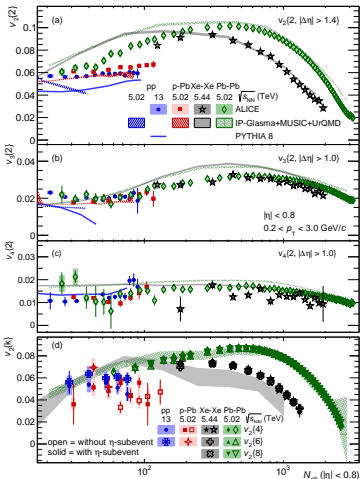


Bayesian analysis result using LHC data

- Resulting posterior for Δr_{np} is compatible with PREX II and *ab initio* nuclear theory.
- Slightly stronger constraint than PREX II ($\Delta r_{np} = 0.283 \pm 0.071$).
- Result is in principle improvable with better Bayesian analyses.
 - May be hard to do in practice.
 - The current analysis already took 2M CPUh.



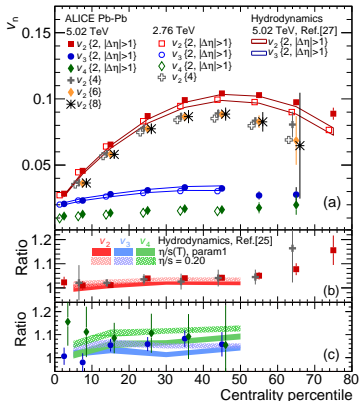
One fluid to rule them all?



- Anisotropic flow is present in a great range of system sizes:
 - PbPb,
 - High multiplicity pPb,
 - High multiplicity pp,
 - ...
- Is this a sign of hydrodynamics?
 - Hydrodynamical simulations seem to work reasonably well.
 - But can a system that small really behave hydrodynamically?
 - Initial state geometry is poorly understood.
- We need a precision test of hydrodynamics in small systems.



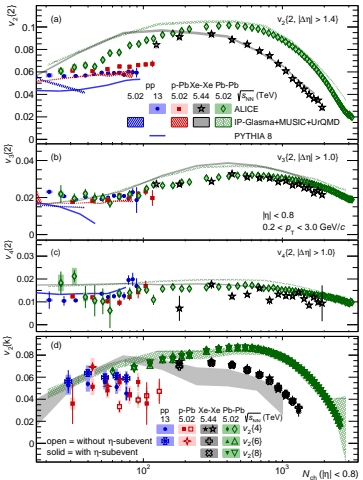
Recap: why do we believe PbPb is hydrodynamic?



- Not just the presence of $v_n\{k\}$.
- We understand where the $v_n\{k\}$ come from!
 - Hydrodynamics converts initial state anisotropic geometry into final state momentum anisotropy.
 - We understand very well what the initial geometry looks like!
- For pPb this is not the case.
 - There is $v_n\{k\}$ measured.
 - But we do not understand the initial geometry.
 - No clear interpretation of experimental results.



Posing a precise question

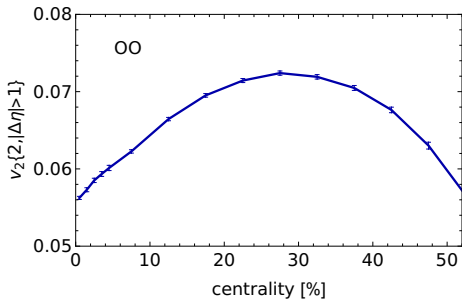


- Can we describe PbPb and a small system in a hydrodynamical model *with the same settings*?
 - Hydro model used should describe a wide range of PbPb observables.
- Can we find a quantity to predict which does not suffer from huge theoretical uncertainties? **Wishlist:**
 - Initial geometry under control.
 - Small sensitivity to proton substructure.
 - No longitudinal structure issues.
 - Quantifiable and small theory uncertainty.



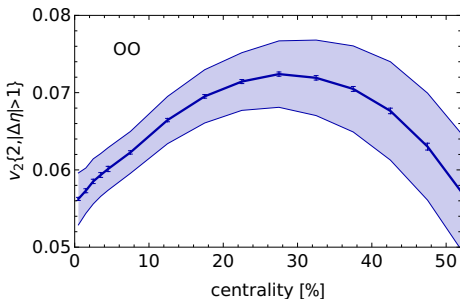
Can $^{16}\text{O}^{16}\text{O}$ collisions help?

- $^{16}\text{O}^{16}\text{O}$ collisions are planned at the LHC for 2025.
- Shape of the proton and longitudinal structure are not an issue, but...



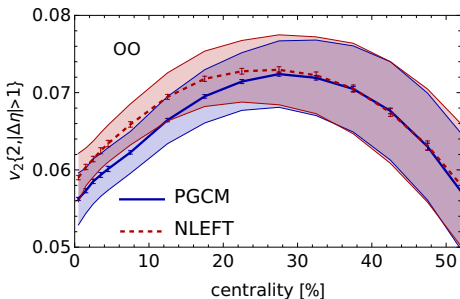
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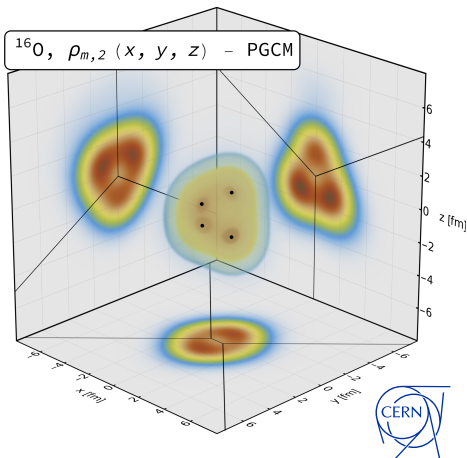
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- Magnitude of fluctuations in the initial state is poorly constrained.
- Different nuclear structure calculations give different answers!
- We have a handle on systematics, but errors are substantial.



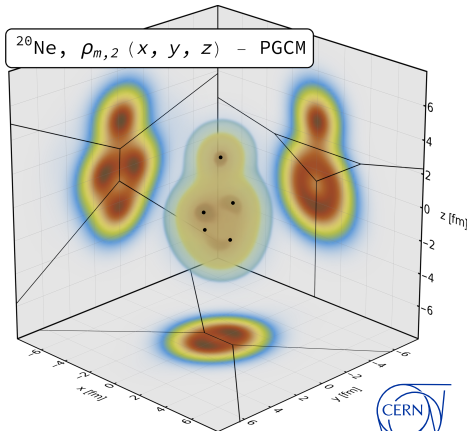
The nuclear bowling pin: ^{20}Ne

- We use both the PGCM and NLEFT frameworks for our nuclear structure input.
 - PGCM computes the average deformed densities.
 - NLEFT simulates an effective theory on a lattice.
- ^{16}O is shaped like an irregular tetrahedron.
- ^{20}Ne is close in size, but has the most extreme shape in the Segrè chart.
- Can we take a ratio between systems to cancel the uncertainties?



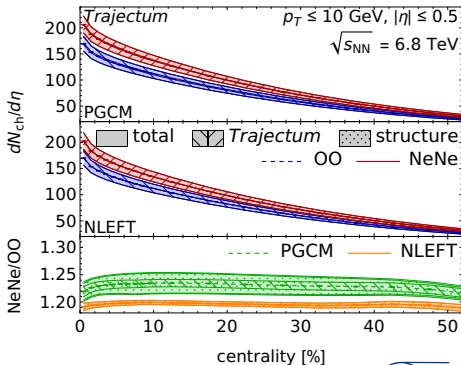
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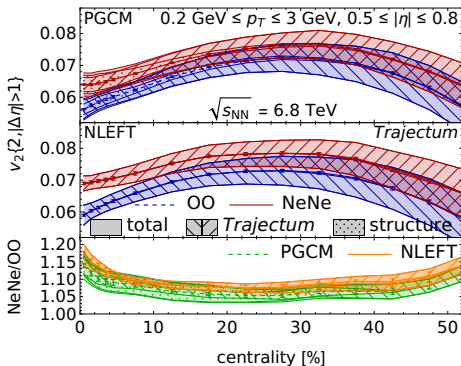
A careful look at uncertainties

- *Trajectum* systematic uncertainty contains contributions from:
 - Uncertainties in parameters.
 - Extrapolation to zero grid spacing.
- PGCM systematic uncertainty contains contributions from:
 - Sampling method: how to convert a density into a configuration.
 - Constraint application: order of operations in the PGCM computation.
- NLEFT systematic uncertainty contains contributions from:
 - Resolution of ambiguities from periodicity of the lattice.



Comparing ^{20}Ne to ^{16}O significantly reduces errors!

- NLEFT and PGCM are consistent within uncertainties.
- Ratio of $v_2\{2\}$ reaches percent level precision from 5% to 20% centrality!
- Difference of $\rho(v_2\{2\}^2, \langle p_T \rangle)$ has uncertainty reduced by up to a factor 6!
- Larger PGCM uncertainty is mostly due to ambiguity in how to generate configurations from densities.

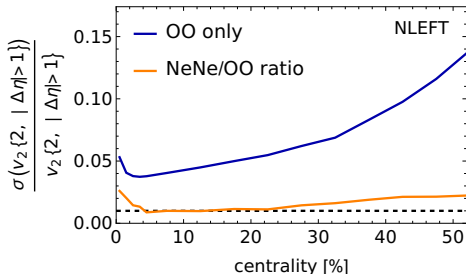


0–1%	$v_2\{2\}_{\text{NeNe}}/v_2\{2\}_{\text{OO}}$	$\rho_{2,\text{NeNe}} - \rho_{2,\text{OO}}$
NLEFT	1.170(8) _{stat.} (30) ^{Traj.} _{syst.} (0) ^{str.} _{syst.}	−0.121(14) _{stat.} (10) ^{Traj.} _{syst.} (0) ^{str.} _{syst.}
PGCM	1.139(6) _{stat.} (27) ^{Traj.} _{syst.} (28) ^{str.} _{syst.}	−0.124(10) _{stat.} (10) ^{Traj.} _{syst.} (29) ^{str.} _{syst.}



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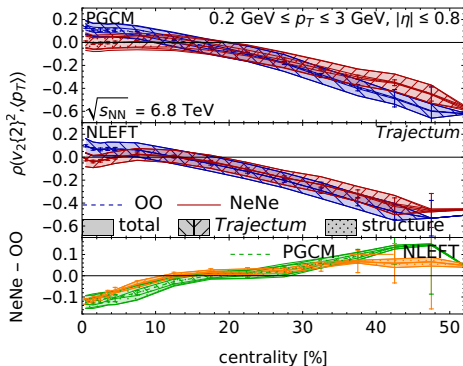


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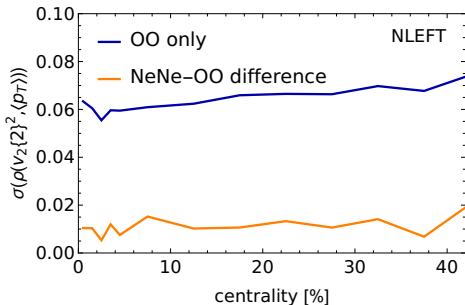


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Comparing ^{20}Ne to ^{16}O significantly reduces errors!

- NLEFT and PGCM are consistent within uncertainties.
- Ratio of $v_2\{2\}$ reaches percent level precision from 5% to 20% centrality!
- Difference of $\rho(v_2\{2\}^2, \langle p_T \rangle)$ has uncertainty reduced by up to a factor 6!
- Larger PGCM uncertainty is mostly due to ambiguity in how to generate configurations from densities.



0–1%	$v_2\{2\}_{\text{NeNe}}/v_2\{2\}_{\text{OO}}$	$\rho_{2,\text{NeNe}} - \rho_{2,\text{OO}}$
NLEFT	1.170(8) _{stat.} (30) ^{Traj.} (0) _{syst.} ^{str.}	−0.121(14) _{stat.} (10) ^{Traj.} (0) _{syst.} ^{str.}
PGCM	1.139(6) _{stat.} (27) ^{Traj.} (28) _{syst.} ^{str.}	−0.124(10) _{stat.} (10) ^{Traj.} (29) _{syst.} ^{str.}



Conclusions

- Let us take another look at our wishlist:

$v_n\{k\}$ in	$p\text{Pb}$	OO	NeNe/OO
Initial geometry under control	✗	✓	✓
Small sensitivity to proton substructure	✗	✓	✓
No longitudinal decorrelation issues	✗	✓	✓
Quantifiable theory uncertainty	✗	✓	✓
Small theory uncertainty	✗	$\geq 4\%$	$\geq 1\%$

- Theory has a much better handle on $^{16}\text{O}^{16}\text{O}$ compared to $p\text{Pb}$.
- Theory uncertainties can be substantially reduced by supplementing $^{16}\text{O}^{16}\text{O}$ collisions with $^{20}\text{Ne}^{20}\text{Ne}$ collisions.
 - $v_2\{2\}$ ratio can be predicted to 1% precision between 5% and 20% centrality.
 - Different nuclear structure calculations give consistent results.



TH Institute: Light Ions at the LHC

Light ion collisions at the LHC

Location: 4/3-006, CERN
Website: cern.ch/lightions

Date: Nov. 11-15, 2024



Topics covered in relation to small systems:

- Experimental highlights and projections
- Heavy flavour
- Hydrodynamics
- Initial conditions
- Jets
- Ultrapерipheral collisions
- Nuclear parton distribution functions
- Nuclear structure
- LHC accelerator opportunities

Organisers:

- Reyes Alemany Fernandez
- Giuliano Giacalone
- Qipeng Hu
- Govert Hugo Nijss
- Saverio Martini
- Wilke van der Schee
- Huichao Song
- Jing Wang
- Urs Wiedemann
- You Zhou



Backup

Backup

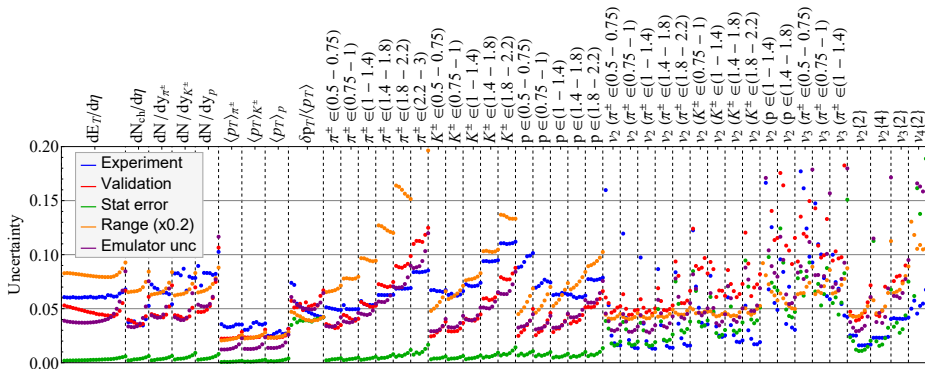


Bayesian analysis details

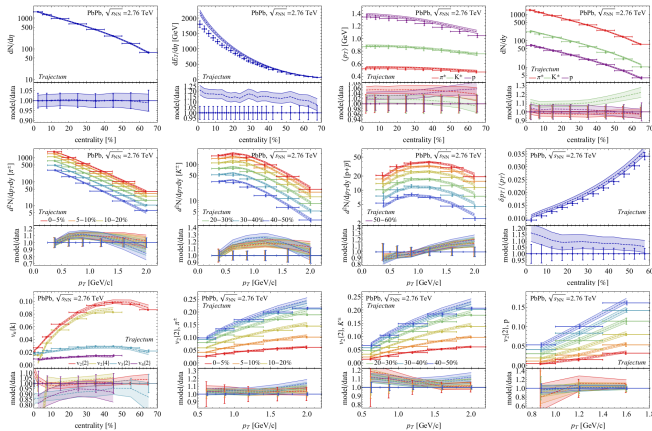
- 3000 design points.
- 18k events per design point.
- Every 15th design point has $10\times$ more statistics, enabling to emulate 'hard' observables such as $SC(n, m)$ and $\rho(v_2\{2\}^2, \langle p_T \rangle)$.



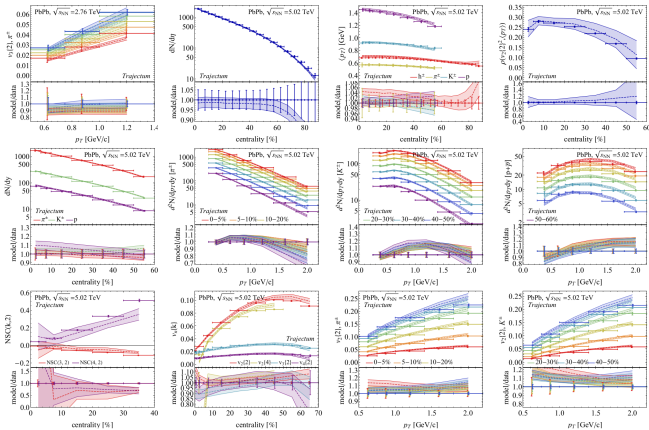
Error budget



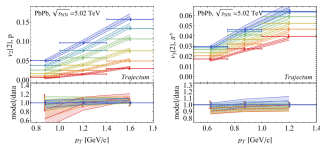
Posterior observables (1/3)



Posterior observables (2/3)



Posterior observables (3/3)



T_RENTo initial conditions

- Nucleons A and B become *wounded* with probability

$$P_{\text{wounded}} = 1 - \exp\left(-\sigma_{gg} \int d\mathbf{x} \rho_A(\mathbf{x}) \rho_B(\mathbf{x})\right), \quad \rho_A \propto \exp\left(\frac{-|\mathbf{x} - \mathbf{x}_A|^2}{2w^2}\right).$$

- Each wounded nucleon deposits energy into its nucleus's *thickness function* $\mathcal{T}_{A/B}$:

$$\mathcal{T}_{A/B} = \sum_{i \in \text{wounded } A/B} \gamma \exp(-|\mathbf{x} - \mathbf{x}_i|^2/2w^2),$$

with γ drawn from a gamma distribution with mean 1 and standard deviation σ_{fluct} .

- Actual formulas slightly modified because each nucleon has n_c constituents.



The T_RENTo phenomenological ansatz

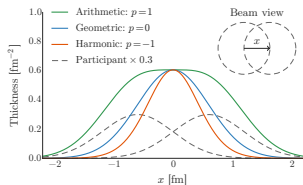
- The standard T_RENTo formula combines thickness functions of the two nuclei \mathcal{T}_A and \mathcal{T}_B into a *reduced thickness* \mathcal{T} , interpreted as an energy density:

$$\mathcal{T} \propto \left(\frac{\mathcal{T}_A^p + \mathcal{T}_B^p}{2} \right)^{1/p},$$

with p a parameter.

- Some useful limits:

p	-1	0	1
\mathcal{T}	$\frac{2}{\frac{1}{\mathcal{T}_A} + \frac{1}{\mathcal{T}_B}}$	$\sqrt{\mathcal{T}_A \mathcal{T}_B}$	$\frac{\mathcal{T}_A + \mathcal{T}_B}{2}$



Free streaming pre-hydrodynamic stage

- T_{REN}To creates matter at proper time $\tau = 0^+$.
- Propagate the matter using free streaming:

$$T^{\mu\nu}(x, y, \tau_{\text{hyd}}) = \frac{1}{\tau_{\text{hyd}}} \int d\phi \hat{p}^\mu \hat{p}^\nu \mathcal{T}(x - \tau_{\text{hyd}} \cos \phi, y - \tau_{\text{hyd}} \sin \phi),$$

with

$$\hat{p}^\mu = (1 \quad \cos \phi \quad \sin \phi),$$

giving us the stress tensor $T^{\mu\nu}$ at proper time $\tau = \tau_{\text{hyd}}$.

- Here τ_{hyd} is the time at which hydrodynamics is started.
- The factor $1/\tau_{\text{hyd}}$ is due to longitudinal expansion.



Basics of hydrodynamics

- Hydrodynamics is the ultimate effective theory. Knowledge of the underlying microscopic theory is completely summarized in transport coefficients.
- Only conservation laws survive: equation of motion is simply

$$\partial_\mu T^{\mu\nu} = 0.$$

- Not enough equations to close the system. Need additional assumption of *local thermal equilibrium*.
- We write $T^{\mu\nu}$ in terms of building blocks T , u^μ , $g^{\mu\nu}$ and ∂_μ .



Hydrodynamics in the 14-moment approximation

- Define ($g^{\mu\nu} = \text{diag}(1, -1, -1, -1)$):

$$\Delta^{\mu\nu} = g^{\mu\nu} - u^\mu u^\nu, \quad \nabla^\mu = \Delta^{\mu\nu} \partial_\nu, \quad D = u^\mu \nabla_\mu, \quad \sigma^{\mu\nu} = \nabla^{\langle\mu} u^{\nu\rangle},$$

with $\langle \rangle$ symmetrizing and removing the trace.

- We solve viscous hydrodynamics without currents, i.e.

$$\partial_\mu T^{\mu\nu} = 0, \quad T^{\mu\nu} = e u^\mu u^\nu - (P + \Pi) \Delta^{\mu\nu} + \pi^{\mu\nu},$$

- $\pi^{\mu\nu}$ and Π follow the 14-moment approximation:

$$\begin{aligned} -\tau_\pi \Delta_\alpha^\mu \Delta_\beta^\nu D \pi^{\alpha\beta} &= \pi^{\mu\nu} - 2\eta \sigma^{\mu\nu} + \delta_{\pi\pi} \pi^{\mu\nu} \nabla \cdot u \\ &\quad - \phi_7 \pi_\alpha^{\langle\mu} \pi^{\nu\rangle\alpha} + \tau_{\pi\pi} \pi_\alpha^{\langle\mu} \sigma^{\nu\rangle\alpha} - \lambda_{\pi\Pi} \Pi \sigma^{\mu\nu}, \\ -\tau_\Pi D \Pi &= \Pi + \zeta \nabla \cdot u + \delta_{\Pi\Pi} \nabla \cdot u \Pi - \lambda_{\Pi\pi} \pi^{\mu\nu} \sigma_{\mu\nu}. \end{aligned}$$



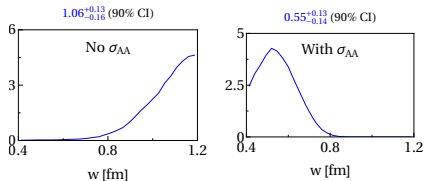
Particlization

- At the freeze-out temperature T_{sw} , we turn the fluid back into particles.
- Particles are sampled thermally, and boosted with the fluid velocity u^μ .
- We use the PTB prescription to match $\pi^{\mu\nu}$ and Π across the transition, so that $T^{\mu\nu}$ is smooth.
- After particlization, we use SMASH as a hadronic afterburner.



Fitting to the p Pb and PbPb cross sections

- In the TRÉNTó model, the nucleon size is described by the Gaussian radius w .
- Previous analyses favored $w \approx 1$ fm.
 - This leads to a 3σ discrepancy in σ_{PbPb} .
- Fitting to the p Pb and PbPb cross sections lowers w to 0.6 fm.
 - σ_{PbPb} discrepancy is reduced to 1σ .
 - Many other observables fit slightly worse.
- Smaller width is now compatible with our knowledge of the gluonic structure of the proton at low x .

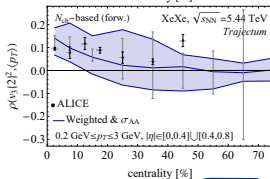
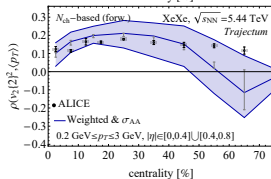
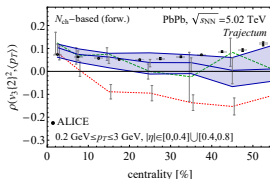
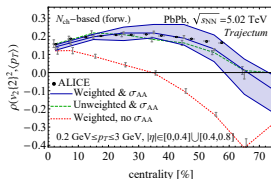


	σ_{PbPb} [b]	$\sigma_{p\text{Pb}}$ [b]
with σ_{AA}	8.02 ± 0.19	2.20 ± 0.06
without σ_{AA}	8.95 ± 0.36	2.48 ± 0.10
ALICE/CMS	7.67 ± 0.24	2.06 ± 0.08



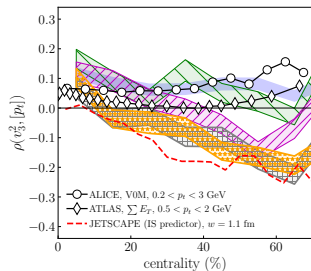
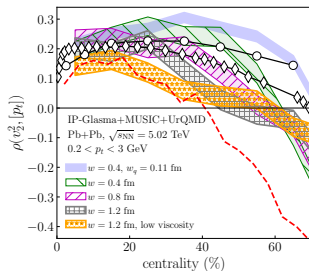
Implication for $\rho(v_n\{2\}^2, \langle p_T \rangle)$

- Pearson correlation coefficient $\rho(v_n\{2\}^2, \langle p_T \rangle)$ between $v_n\{2\}^2$ and $\langle p_T \rangle$ is sensitive to the nucleon size.
- Postdiction without fitting to σ_{PbPb} and σ_{pPb} is qualitatively wrong:
 - $\rho(v_2\{2\}^2, \langle p_T \rangle)$ goes negative already at 30% centrality.
 - $\rho(v_3\{2\}^2, \langle p_T \rangle)$ has the wrong sign.
- Fitting to σ_{PbPb} and σ_{pPb} results in a much improved agreement.



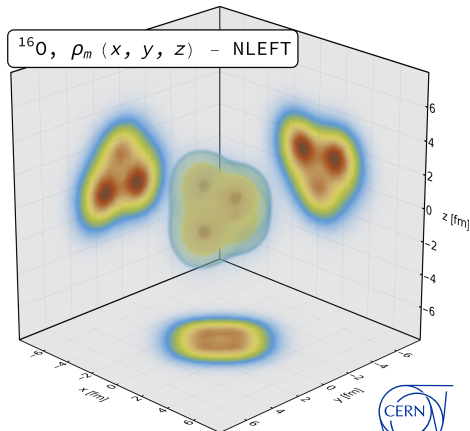
Nucleon width and $\rho(v_n\{2\}^2, \langle p_T \rangle)$

- Previous study shows that $\rho(v_n\{2\}^2, \langle p_T \rangle)$ depends strongly on the nucleon size w .



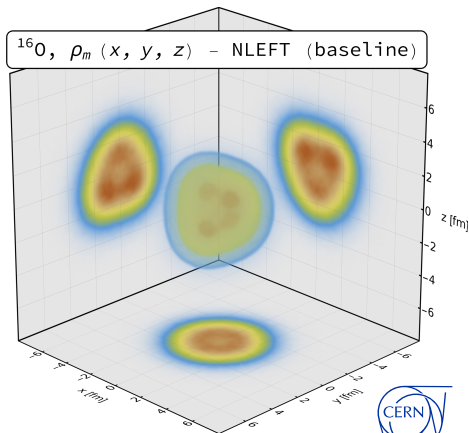
NLEFT densities

- We show the NLEFT densities for ^{16}O and ^{20}Ne .
- Densities are computed from configurations, requiring translation and rotation.
- This introduces biases, so we also show spherical configurations rotated in the same way to illustrate the size of this effect.



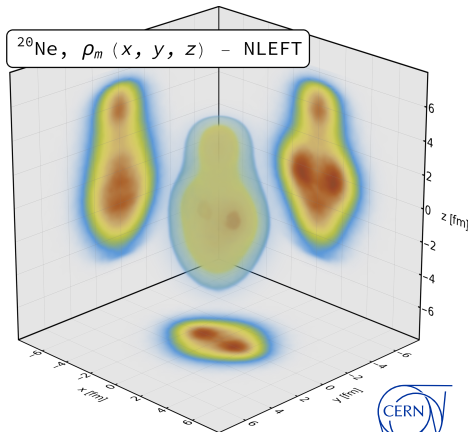
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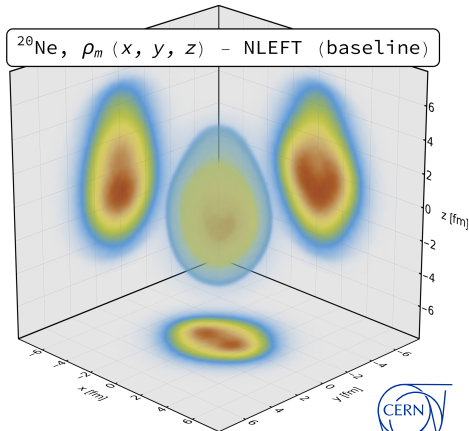
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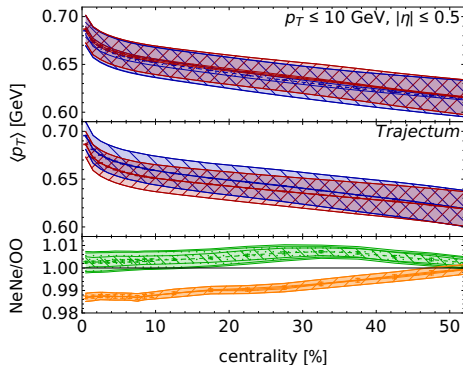
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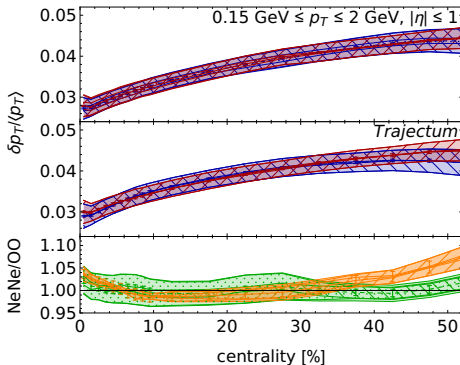
Other observables

- We show the NeNe/OO ratios for $\langle p_T \rangle$, $\delta p_T / \langle p_T \rangle$ and $v_3\{2\}$.
- Discrepancy in $\langle p_T \rangle$ between PGCM and NLEFT is due to the different nuclear charge radius.
- $\delta p_T / \langle p_T \rangle$ has interesting non-monotonic behavior for central collisions.



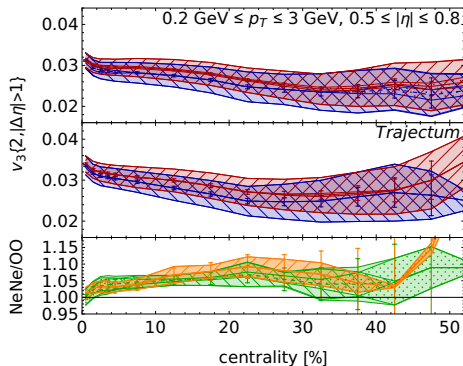
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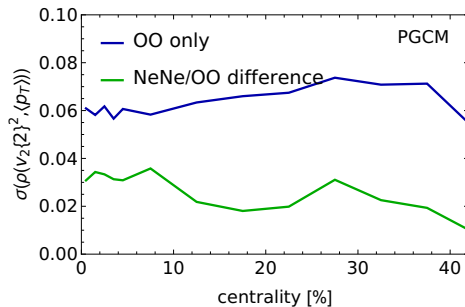
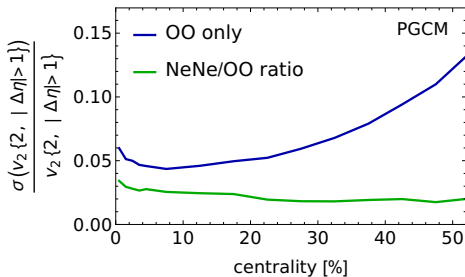


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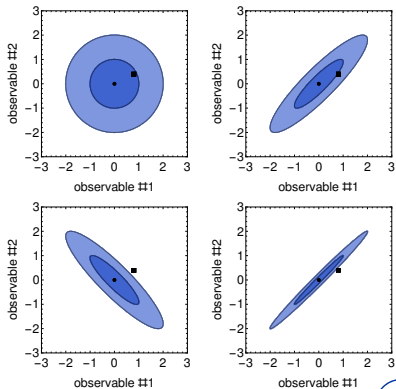


PGCM error ratios



Why weights?

- Higher p_T , higher centralities are harder to model theoretically.
- Experimental correlation matrix is not available.
 - Figure shows 1σ and 2σ regions for $\rho \in \{0, 0.9, -0.9, 0.99\}$, with standard deviations the same.
 - Same difference between theory and experiment can be within 1σ or outside of 2σ depending on ρ .
 - Correlated observable classes can be over/underimportant for the Bayesian analysis.



Definition of weights

- In the bayesian analysis, the probability of the data given the parameter point x is given by:

$$P(D|x) = \frac{1}{\sqrt{(2\pi)^m \det \Sigma}} \exp \left(-\frac{1}{2} (y - y_{\text{exp}})^T \Sigma^{-1} (y - y_{\text{exp}}) \right),$$

with y the vector of observables computed from x , y_{exp} the vector of the corresponding experimental data, and Σ the combined theory/experiment covariance matrix.

- We define weights by replacing

$$P(D|x) = \frac{1}{\sqrt{(2\pi)^m \det \Sigma}} \exp \left(-\frac{1}{2} (y - y_{\text{exp}})^T \omega \Sigma^{-1} \omega (y - y_{\text{exp}}) \right),$$

where ω is the diagonal matrix containing the weight for each observable.



Choice of weights

- We choose for weights ω :
 - 1/2 for every particle identified observable.
 - 1/2 for p_T -differential observables, and an additional $(2.5 - p_T[\text{GeV}])/1.5$ if $p_T > 1$ GeV.
 - $(100 - c[\%])/50$ if the centrality class c is beyond 50%.
- Weighting only worsens the average discrepancy slightly.
- Distribution of discrepancies makes more sense.

	$\langle (Y_{\text{theory}} - Y_{\text{experiment}}) / \sigma \rangle$				$\bar{\omega}$
	σ_{AA} & ω	ω	σ_{AA}	neither	
$dN_{\text{ch}}/d\eta$	0.55	0.60	1.23	1.22	1.00
$dN_{\pi^\pm, K^\pm, p^\pm}/dy$	0.76	0.70	0.60	0.57	0.48
$dE_T/d\eta$	1.59	1.51	0.82	0.77	0.48
$\langle p_T \rangle_{\text{ch}, \pi^\pm, K^\pm, p^\pm}$	0.66	0.60	0.88	0.72	0.46
$\delta p_T / \langle p_T \rangle$	0.56	0.62	0.51	0.58	0.49
$v_n\{k\}$	0.58	0.51	0.54	0.49	1.00
$d^2 N_{\pi^\pm} / dy dp_T$	1.19	1.07	0.86	0.92	0.20
$d^2 N_{K^\pm} / dy dp_T$	1.41	1.27	0.79	0.73	0.20
$d^2 N_{p^\pm} / dy dp_T$	1.35	1.21	0.73	0.67	0.25
$v_2^{\pi^\pm}(p_T)$	0.81	0.74	0.46	0.44	0.19
$v_2^{K^\pm}(p_T)$	0.92	0.89	0.55	0.55	0.19
$v_2^{p^\pm}(p_T)$	0.49	0.47	0.34	0.35	0.25
$v_3^{\pi^\pm}(p_T)$	0.65	0.57	0.69	0.57	0.24
average	0.89	0.83	0.69	0.66	
σ_{AA}	1.13	3.80	1.53	3.40	1.00



How much do weights change the posteriors?

