

# Collective phenomena with machine learning

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CoE in Quark Matter  
YoctoLHC

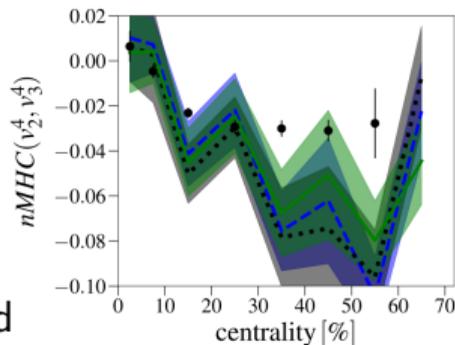
Particle Physics Day 24.11.2022



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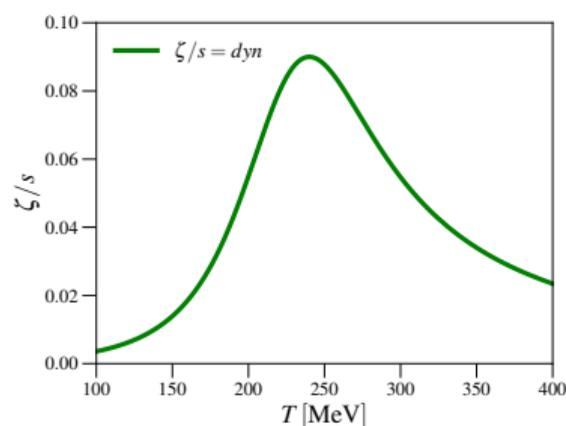
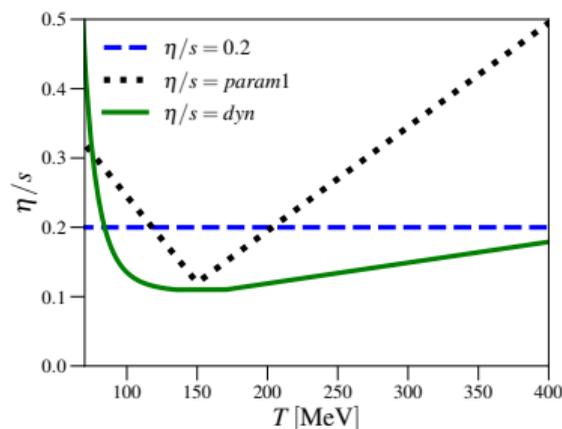


- Heavy ion collisions provide a way to probe matter properties of QGP
- Shear and bulk viscosities of QGP can be constrained from the measured data by the means of Bayesian analysis
- Measured multi-particle correlations require millions of simulated collision events to obtain enough statistics for reliable comparison with the data
- One event  $\sim 0.5$  CPU hours  $\implies \sim 10^6$  CPU hours per viscosity parametrization
- Problem: to perform Bayesian analysis one needs observables for  $\sim 10^2$  parametrizations, i.e. total of  $\sim 10^8$  CPU hours!
- Solution: Use machine learning to speed up the process



# The theory framework

- Initial state from pQCD+saturation **EKRT-framework**
- 2nd-order viscous fluid dynamics with **shear and bulk viscosities**
  - $\eta/s = 0.2$  and  $\eta/s = param1$  from earlier EbyE EKRT works with  $T_{dec} = 100$  MeV
- Here, we convert the fluid into particle spectrum by calculating Cooper-Frye integral at the decoupling surface determined by **dynamical freeze-out** conditions
  - Purely hydrodynamic description  $\implies$  Continuous parametrization of transport coefficients across all phases of strongly interacting matter



# Dynamical freeze-out

- Fluid dynamics applicable when expansion rate ( $\theta$ )  $\lesssim$  scattering rate ( $1/\tau_\pi$ ) and mean free path ( $\tau_\pi$ )  $\lesssim$  size of the system ( $R$ )  
 $\implies$  Dynamical decoupling conditions:

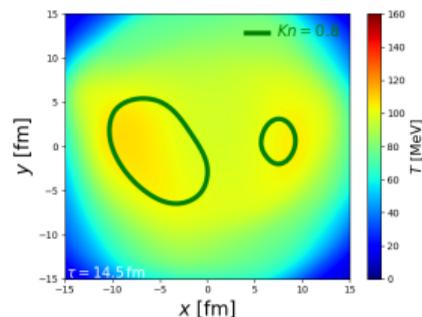
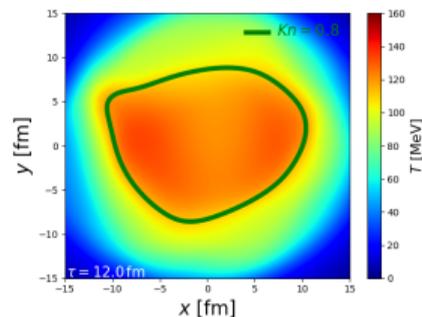
## Knudsen number

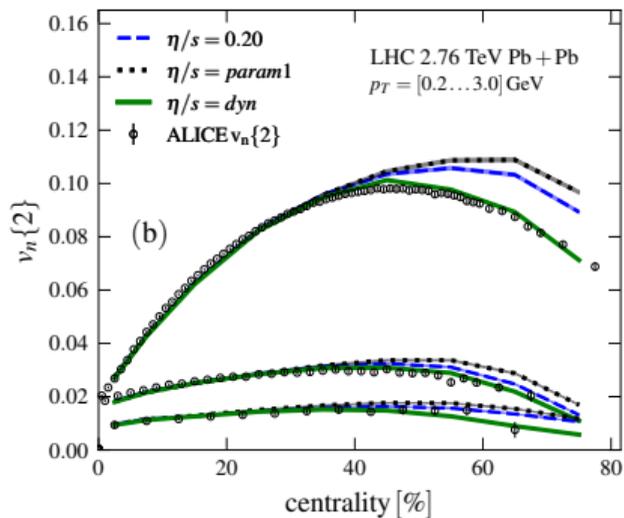
$$Kn \equiv \frac{\text{exp. rate}}{\text{scat. rate}} = \tau_\pi \theta = C_{Kn}$$

## Global size of the system

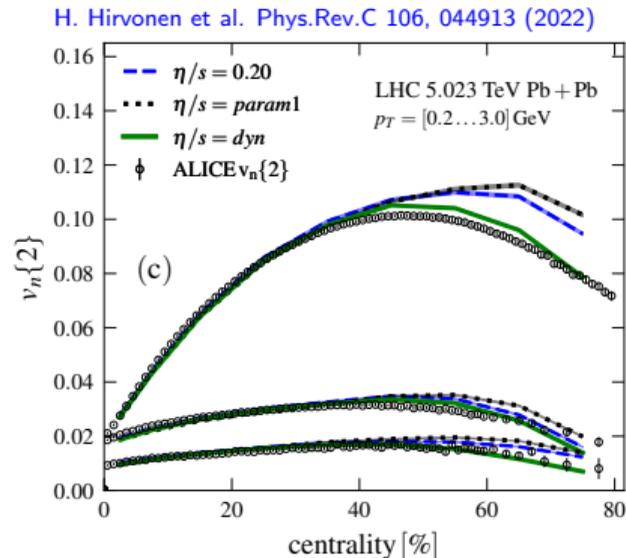
$$\frac{\gamma \tau_\pi}{R} = C_R, \quad R = \sqrt{A/\pi}$$

- $C_{Kn}$  and  $C_R$  are free parameters, fitted from data
- $A$  is the area in which  $Kn < C_{Kn}$  and  $T < 150$  MeV
- Allow multiple separate areas with different  $R$





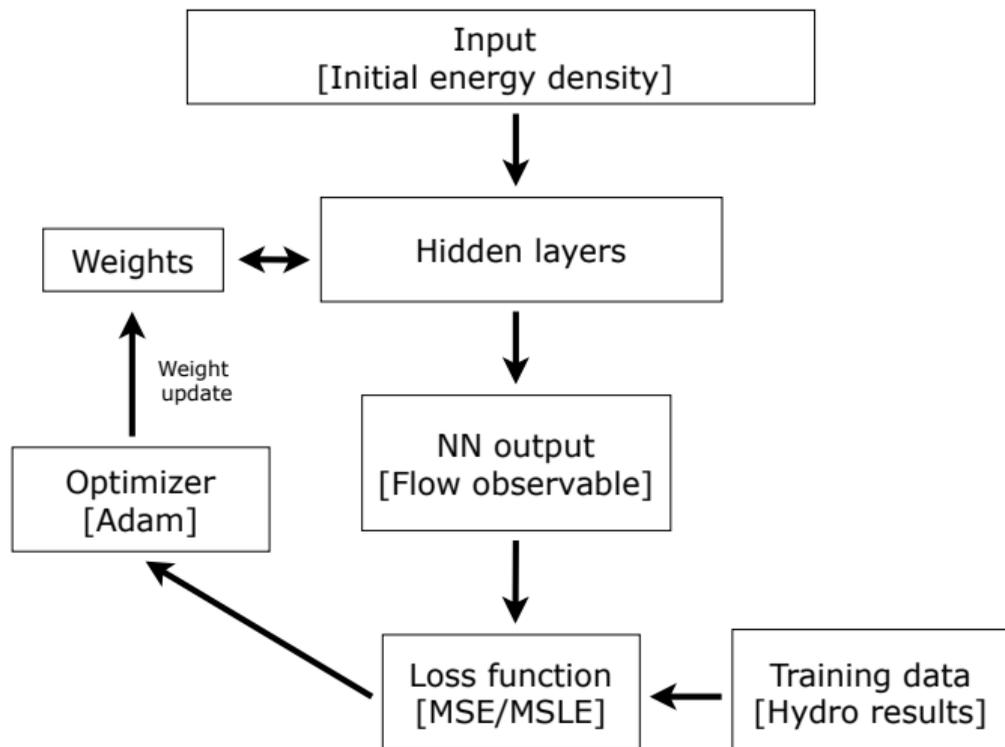
$$\frac{dN}{dyd\phi} = \frac{1}{2\pi} \frac{dN}{dy} \left( 1 + \sum_{n=1}^{\infty} v_n \cos[n(\phi - \Psi_n(p_T))] \right)$$



$$v_n\{2\} = \sqrt{\langle v_n^2 \rangle_{ev}}$$

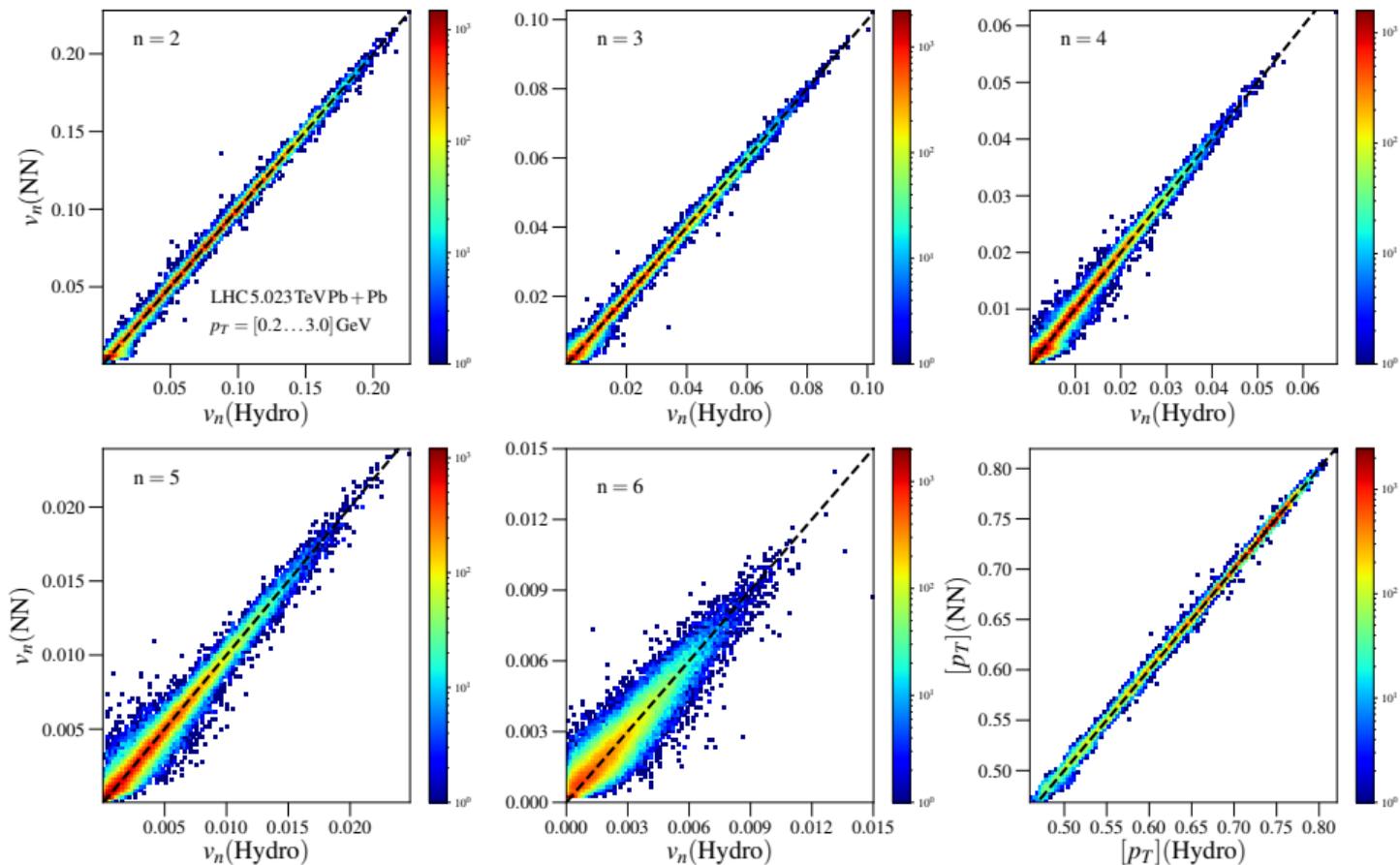
- Dynamical freeze-out decreases amount of flow in peripheral collisions and improves agreement with the measurements

- Layers structure is implemented as a modified version of DenseNet [G. Huang et al. arXiv:1608.06993](#)
  - Very deep network structure with 128 convolutional layers
  - In total of  $\approx 5.4\text{M}$  trainable parameters
- Training data is for one fixed viscosity parametrization

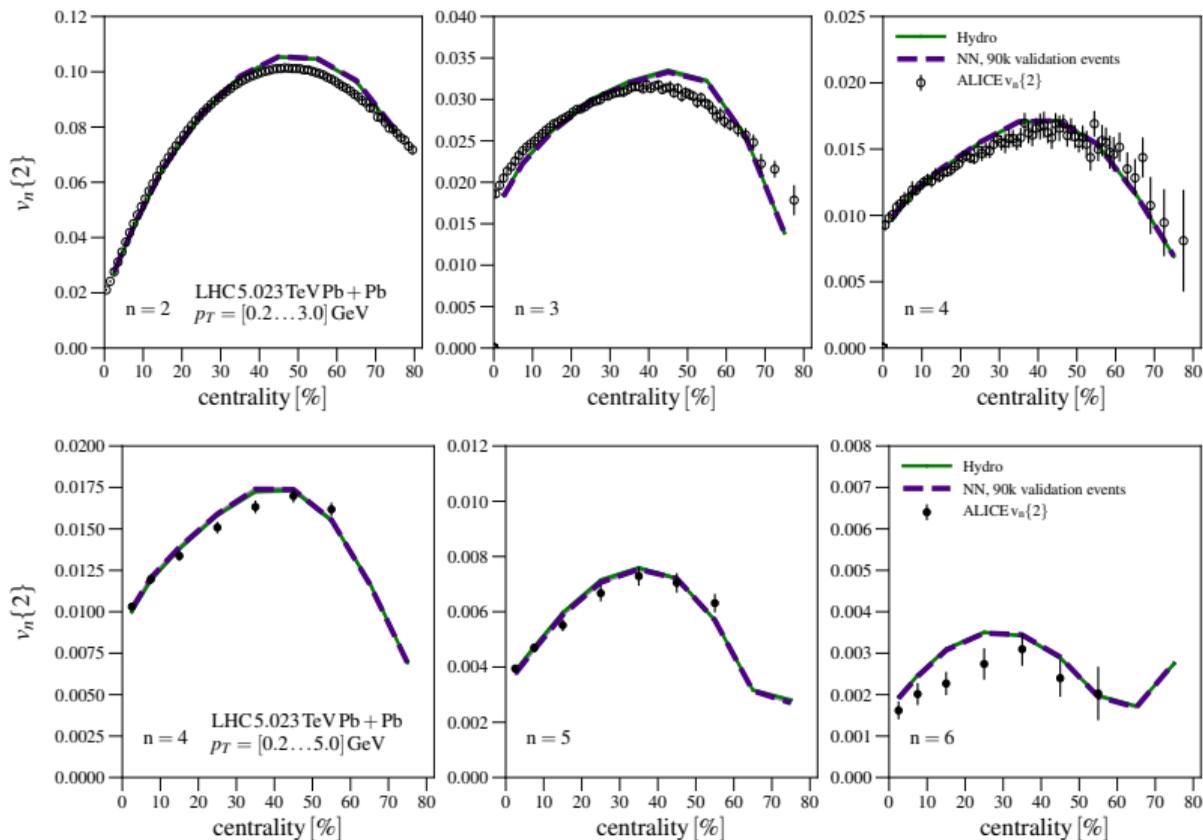


- Separate network trained for each  $p_T$ -integrated observable:  
 $v_2, v_3, v_4, v_5, v_6, [p_T], dN_{ch}/d\eta$
- Single network trained with multiple different  $p_T$  ranges for an observable
- In total of 20000 training events: 5000 from each collision system
  - 200 GeV Au+Au
  - 2.76 TeV Pb+Pb
  - 5.023 TeV Pb+Pb
  - 5.44 TeV Xe+Xe (deformed nuclei)
- Training data augmented using random flips, rotations and translations
- Training one network takes  $\approx 1$  GPU hour with NVIDIA V100 32GB GPU
- After training, NN can generate  $\sim 10^6$  events/hour with GPU
  - Factor of  $10^5$  faster than doing full simulations using CPU!

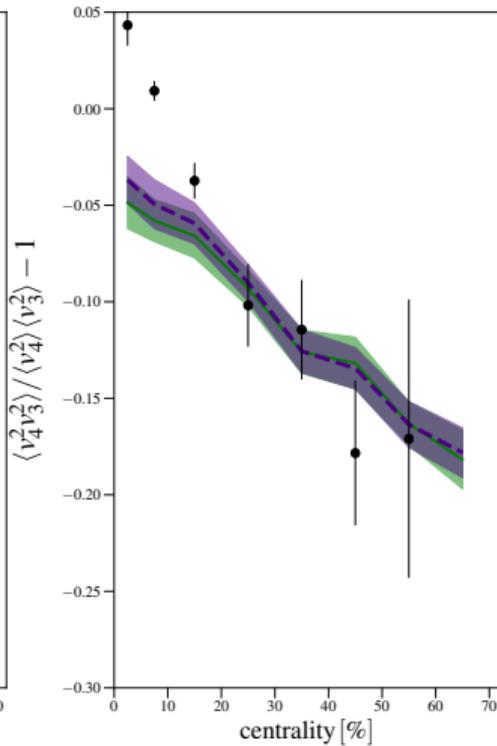
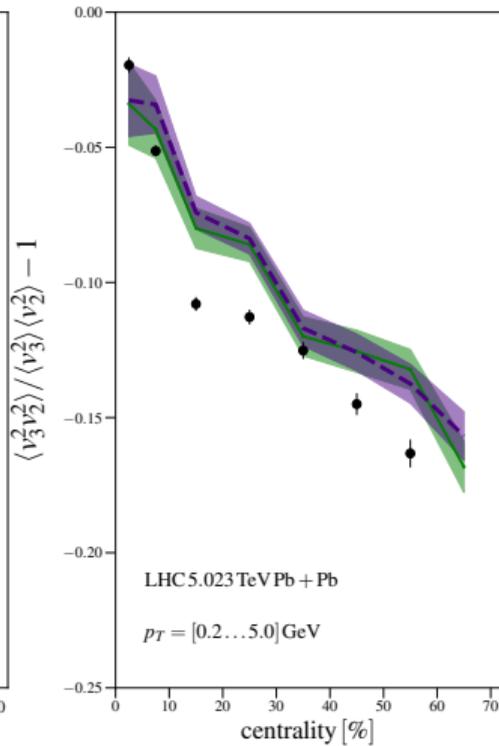
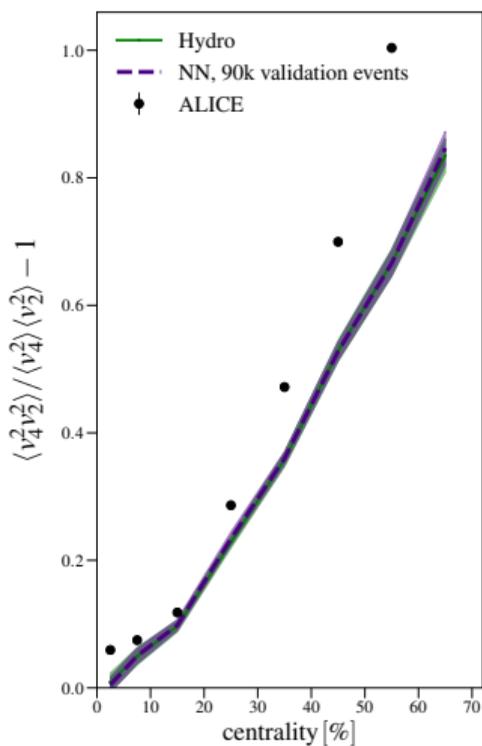
# Validation tests: Errors with 90k validation events



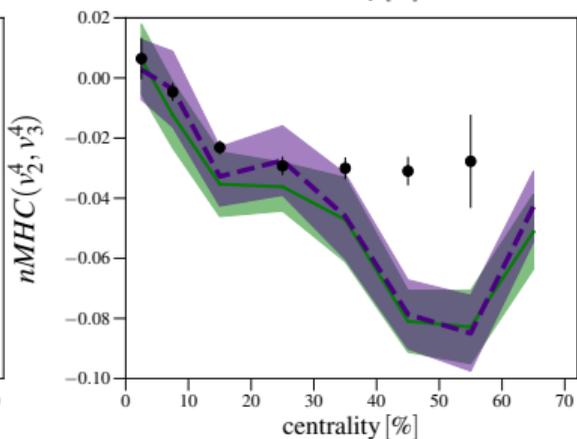
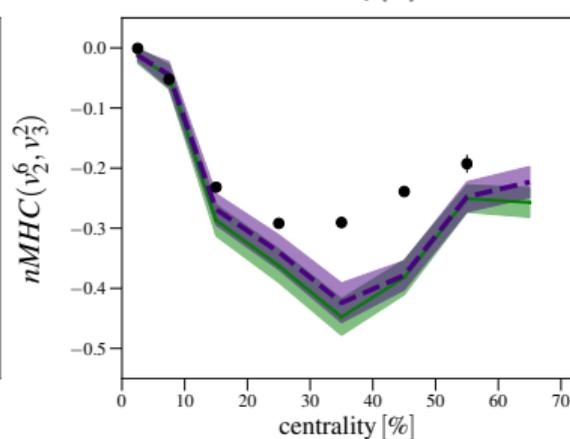
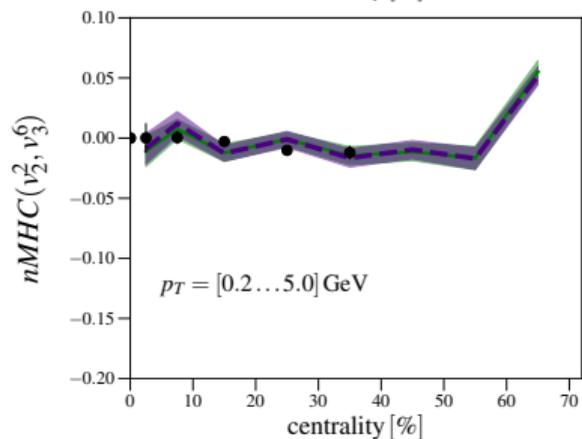
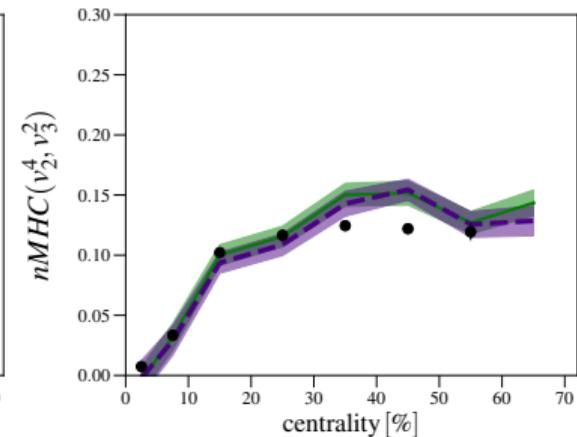
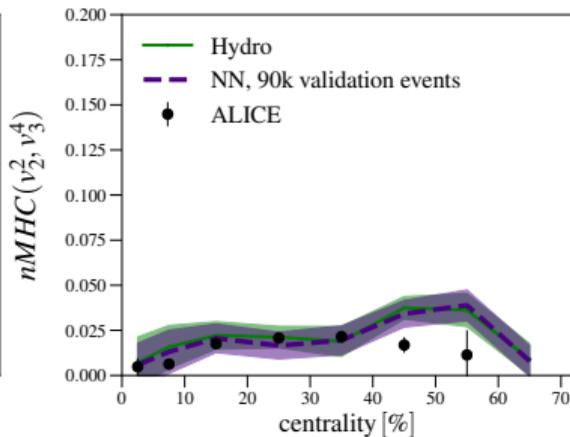
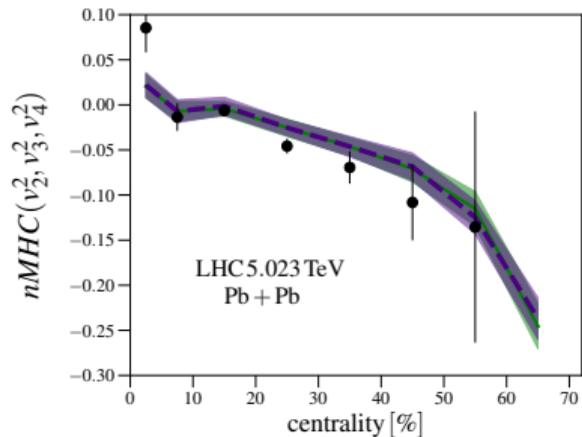
# Validation tests: Flow coefficients



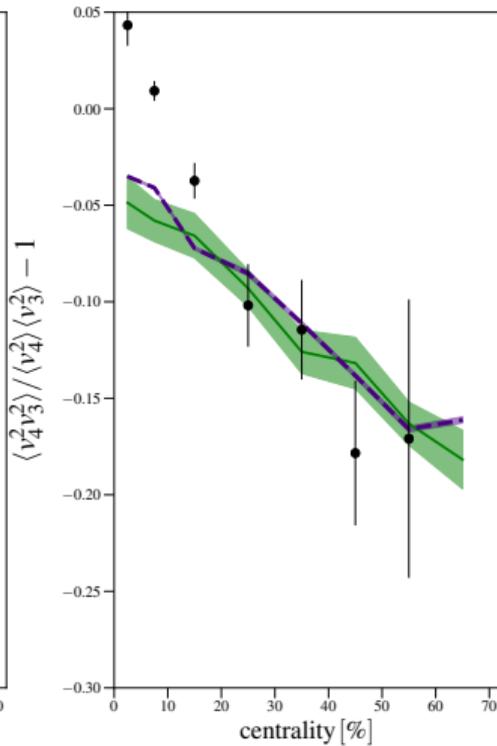
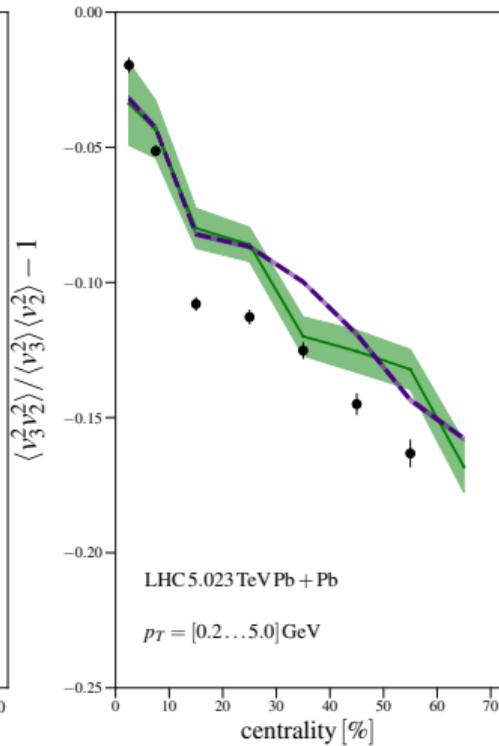
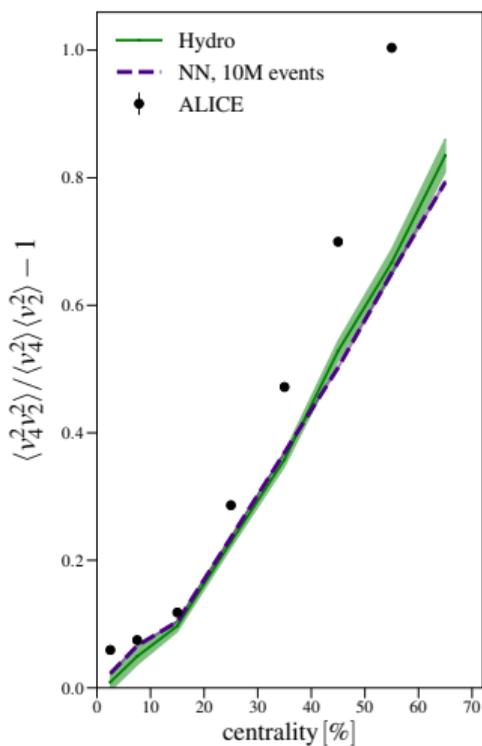
# Validation tests: Four-particle flow correlations



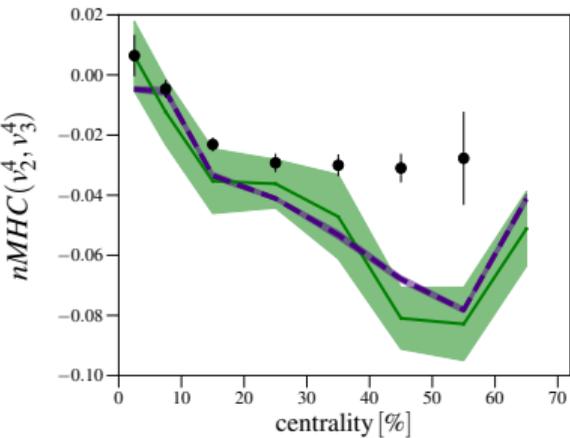
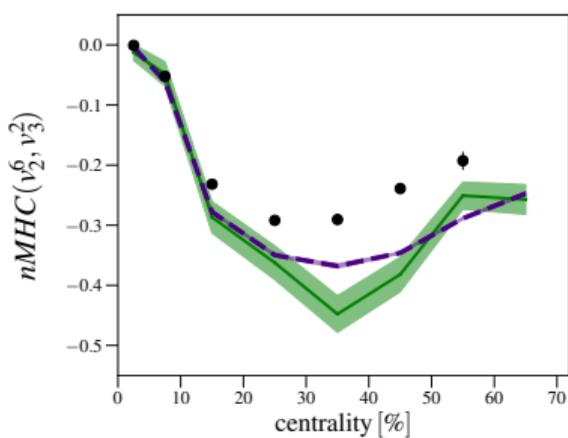
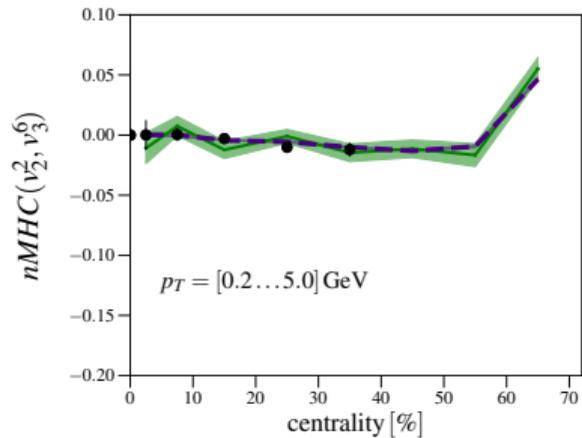
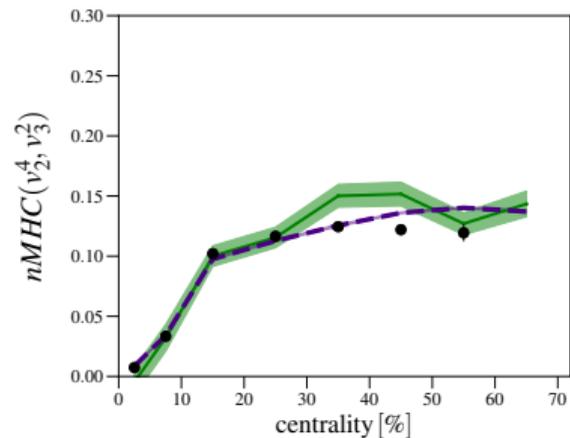
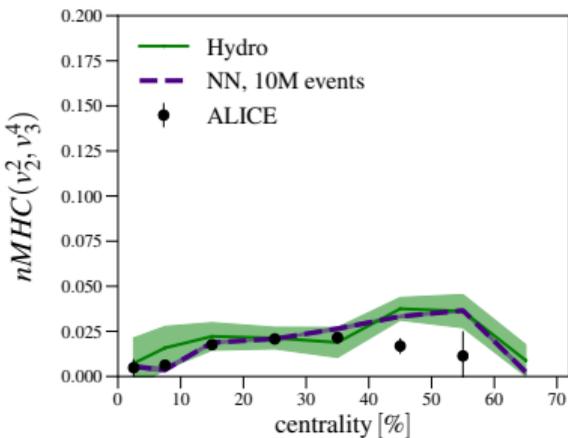
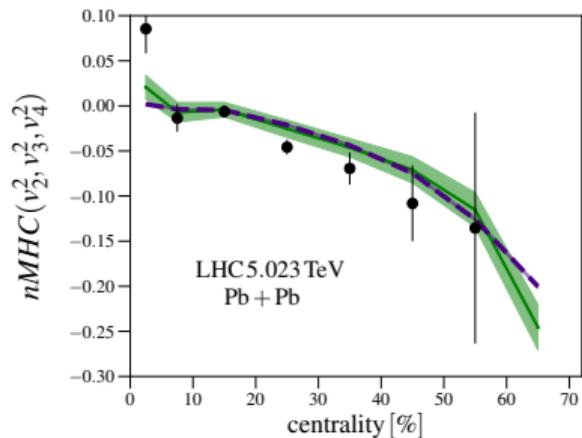
# Validation tests: Six- and eight-particle flow correlations



# NN predictions: Four-particle flow correlations



# NN predictions: Six- and eight-particle flow correlations



- Dynamical decoupling  $\implies$  Clear improvement in centrality dependence of flow coefficients
- Using neural network to predict flow observables from initial energy density reduces computation time by many orders of magnitude
  - Speedup achieved while maintaining good accuracy, even for multi-particle correlations
  - Currently works only for fixed viscosity parametrization
- In future: extend the neural network for arbitrary viscosity parametrizations and use it to perform Bayesian analysis