

Effect of clustered nuclear geometry to azimuthal anisotropy and flow fluctuations in O+O collisions at the LHC

N. Mallick, S. Pasad, R. Sahoo, **G.G. Barnaföldi**

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Ref.: *arXiv:2407.15065 (Submitted to PLB)*

V4HEP#3, Prague, Czech Republic, 3rd October 2024

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REN

Wigner

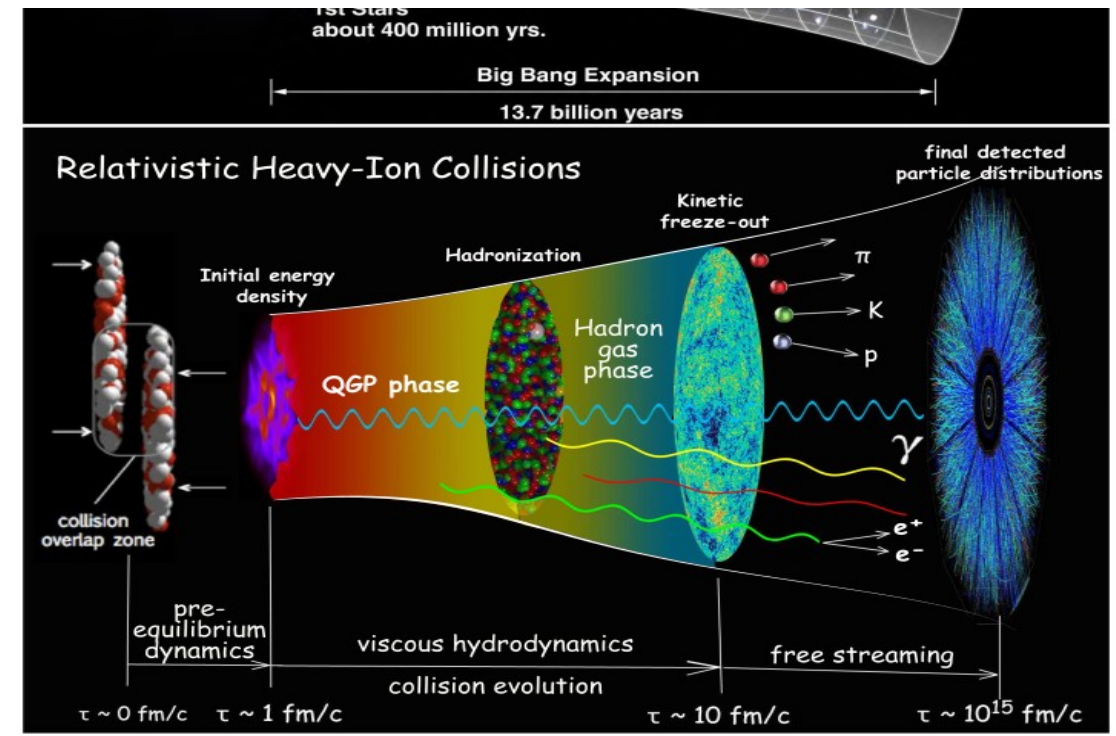
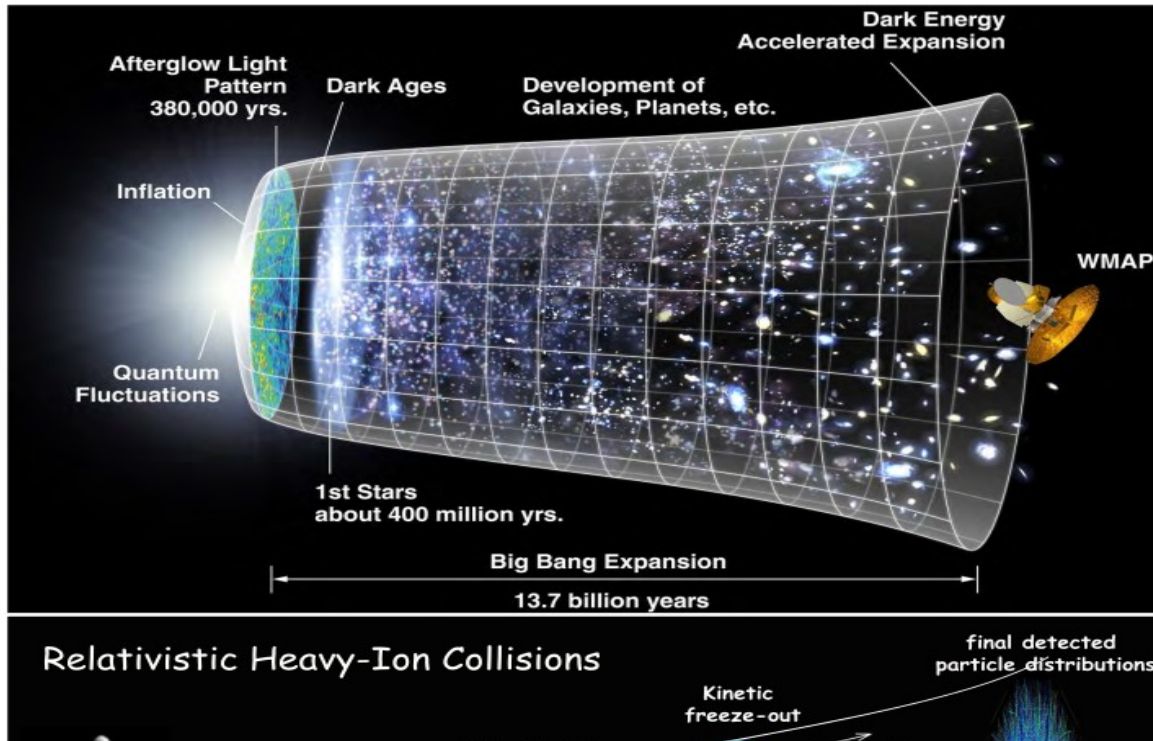
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Motivation & definitions

Primordial matter in heavy-ion collisions

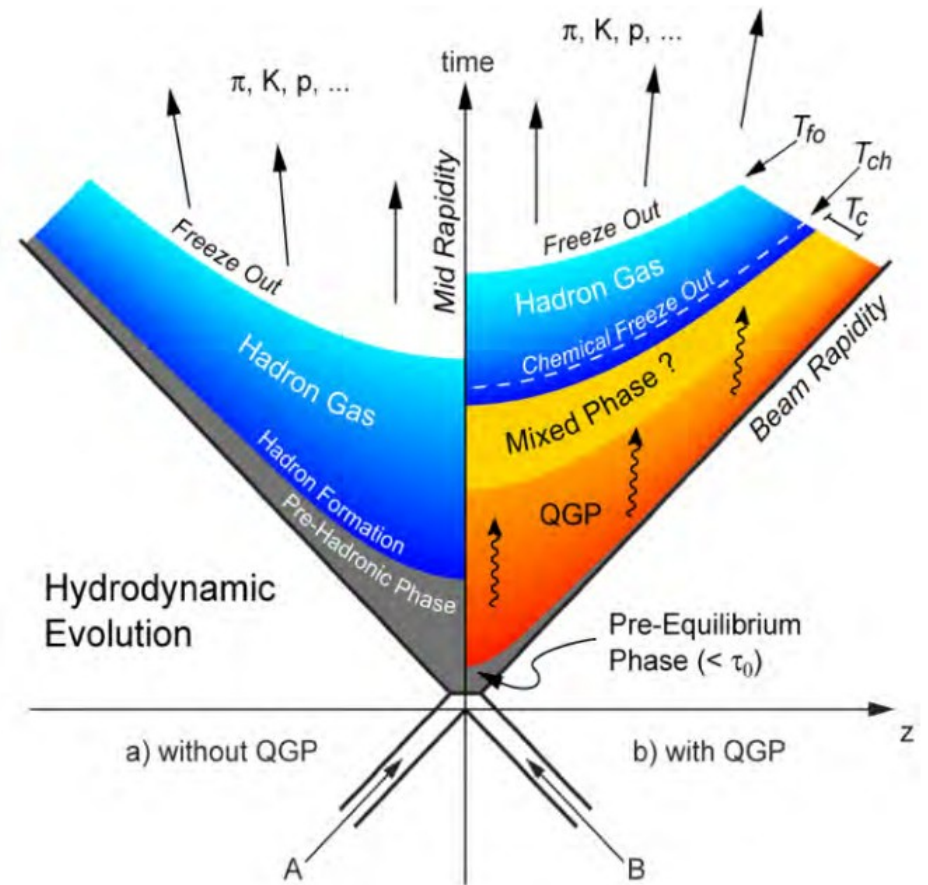
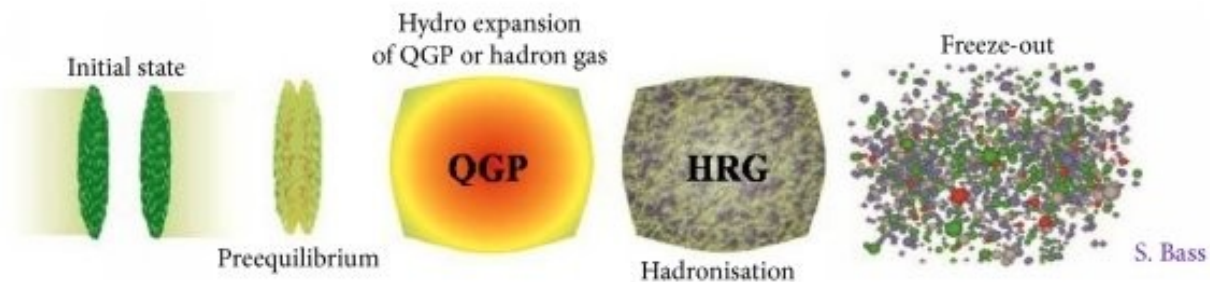
- **Quark-Gluon Plasma (QGP) research**



Primordial matter in heavy-ion collisions

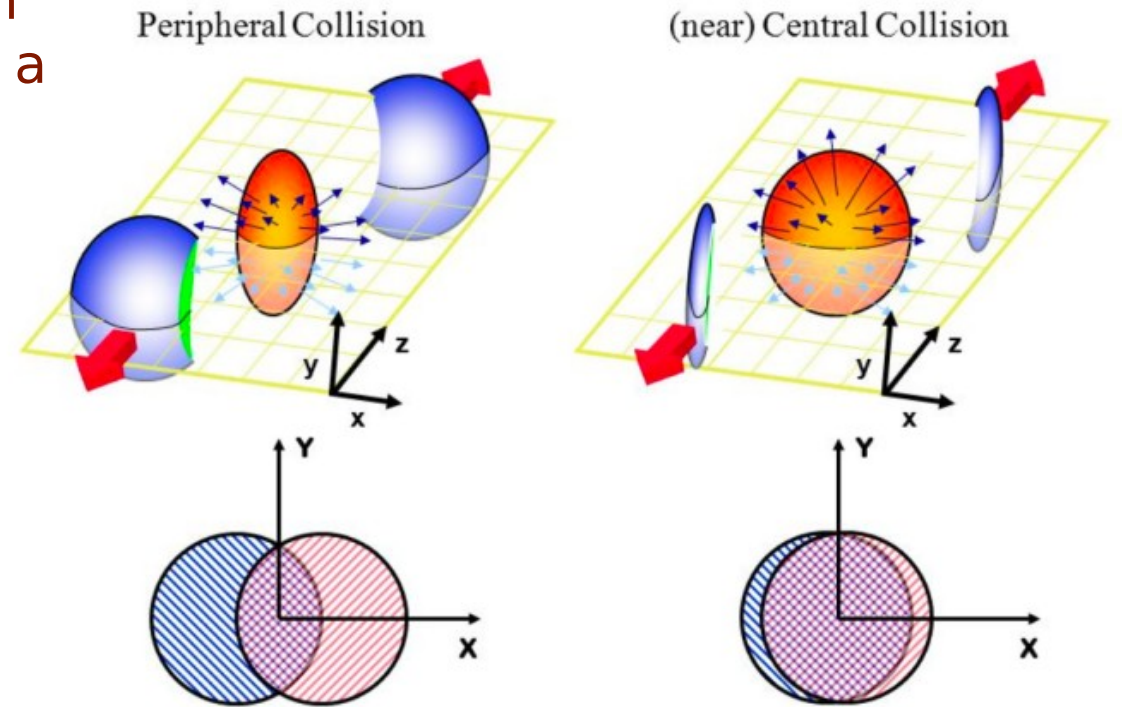
- **QGP in experimental vs theory points**

- By colliding heavy-ions we can form small drop of the hot & dense primordial matter
- No direct observations, just **signatures**: jet-quenching, correlations, collective effects, **(anisotropic) flow...**
- Need a complex description, including QCD phenomenology, hydrodynamics, (non-equilibrium) thermodynamics



Flow (v_n) in heavy-ion collisions

- **Experimental point:**
 - Flow describes the azimuthal momentum space anisotropy of particle emission for a non-central heavy-ion collision.

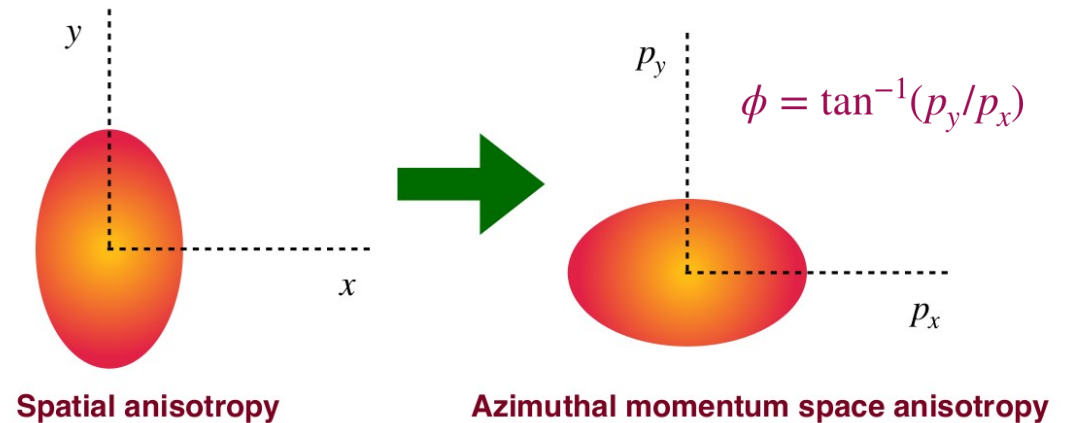


Flow (v_n) in heavy-ion collisions

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- The n^{th} harmonic coefficient of the Fourier expansion of azimuthal momentum distribution:

$$E \frac{d^3N}{dp^3} = \frac{d^2N}{p_T dp_T dy} \frac{1}{2\pi} \left(1 + 2 \sum_{n=1}^{\infty} v_n \cos[n(\phi - \psi_n)] \right)$$



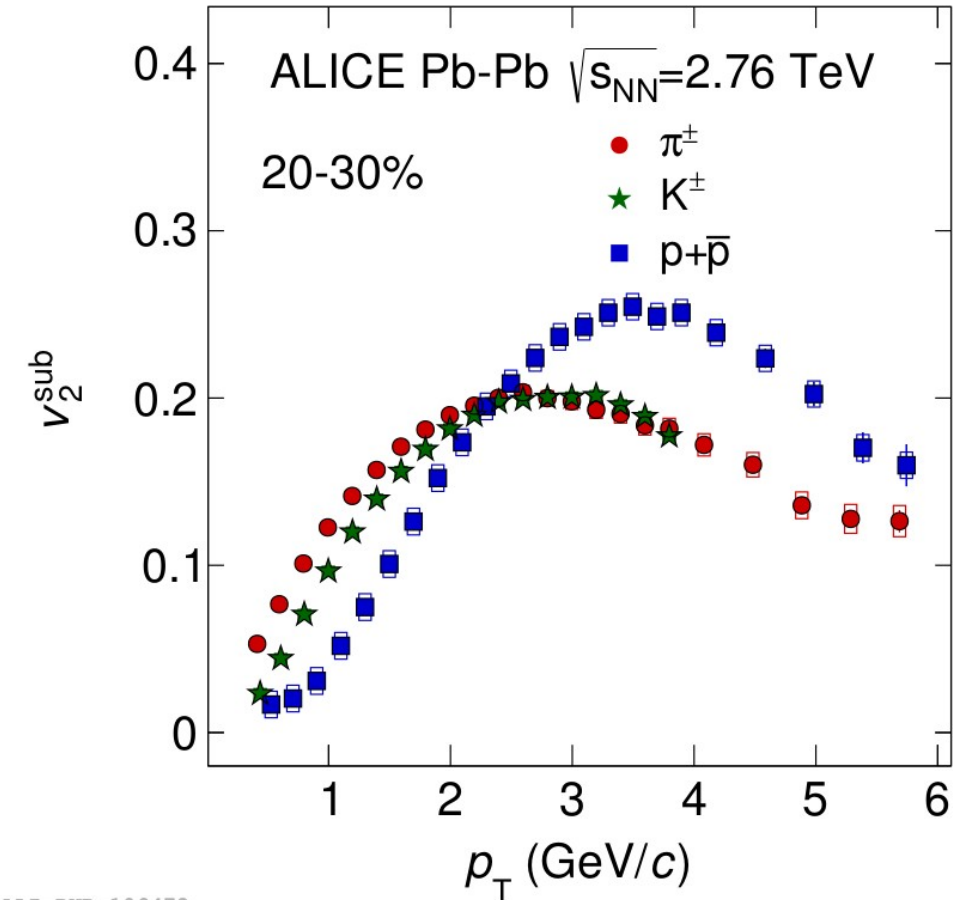
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- The $v_2(p_T, y) = \langle \cos(2(\phi - \psi_2)) \rangle$ directly reflects the initial spatial anisotropy of the nuclear overlap region in the transverse plane.



ALI-PUB-109472

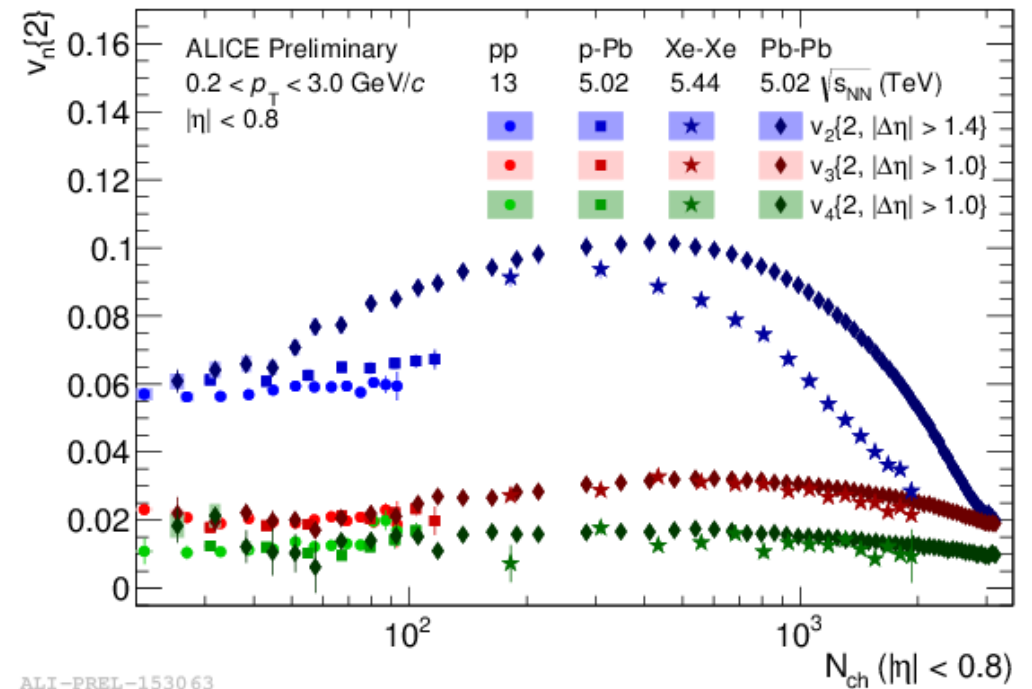
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- The $v_2(p_T, y) = \langle \cos(2(\phi - \psi_2)) \rangle$ directly reflects the initial spatial anisotropy of the nuclear overlap region in the transverse plane.
- Higher flow components can be measured



Future Nuclear Collisions at LHC

- **LHC Schedule with new nuclear collisions**

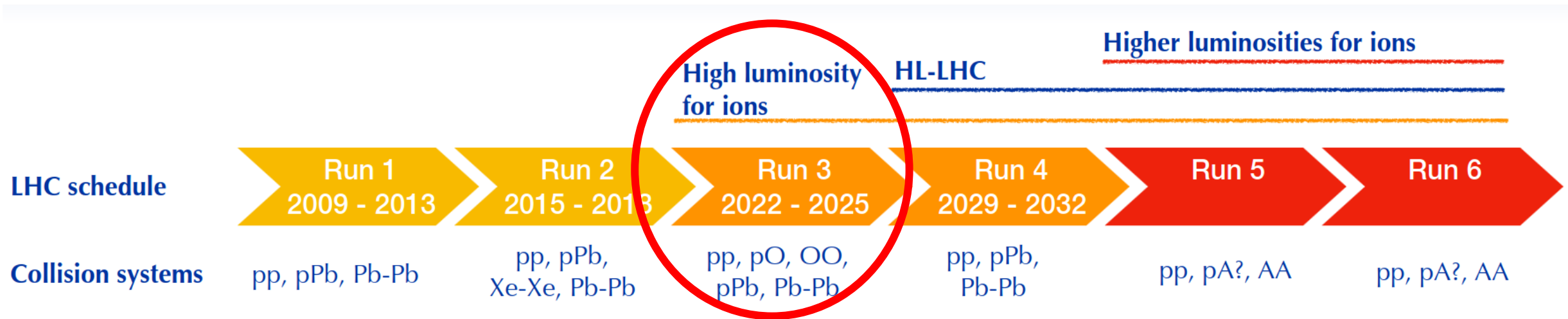
- Run 2: XeXe
- Run 3: pO & OO



Future Nuclear Collisions at LHC

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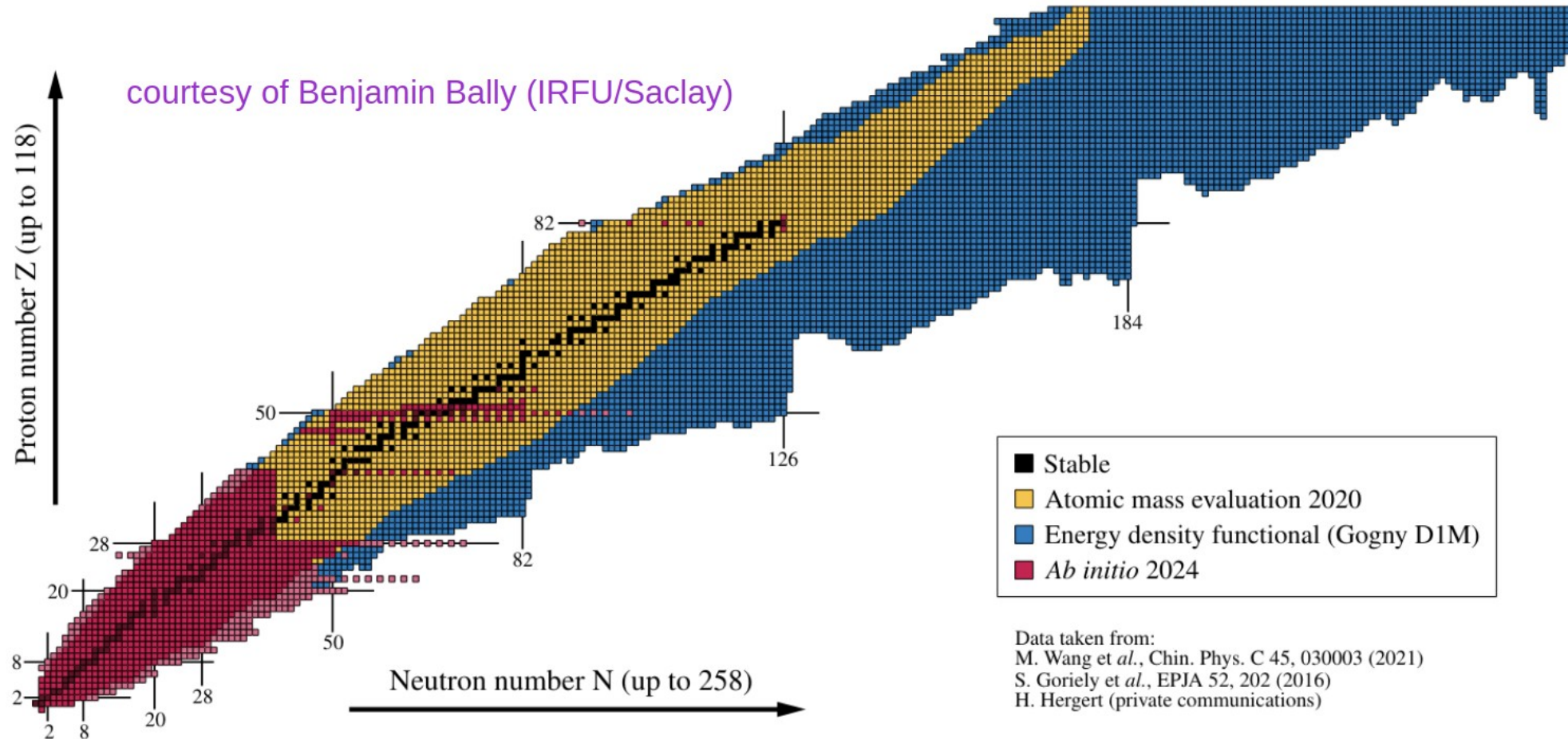
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Nuclei & nuclear structure

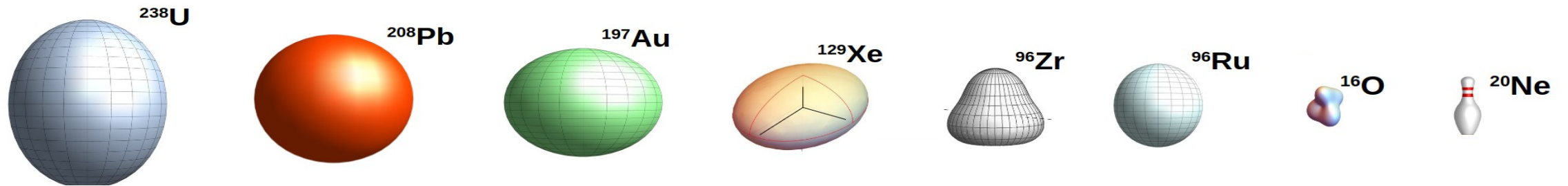
Nuclei for Future Nuclear Collisions

- **High-mass and deformed nuclei are in the focus:**



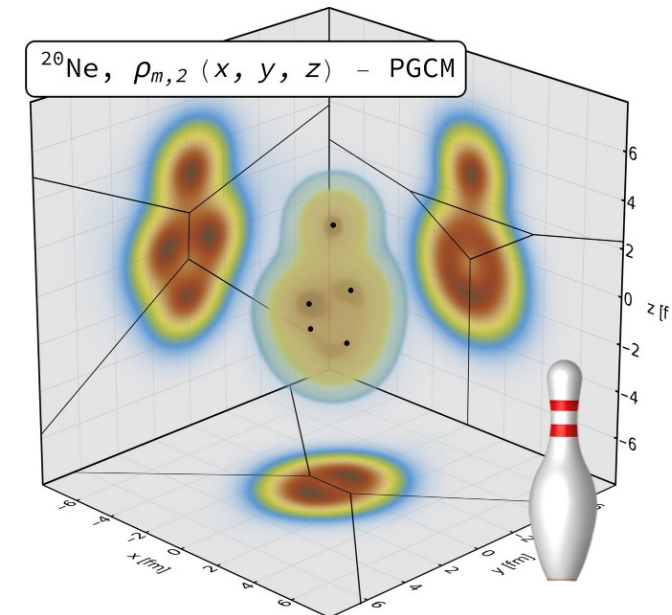
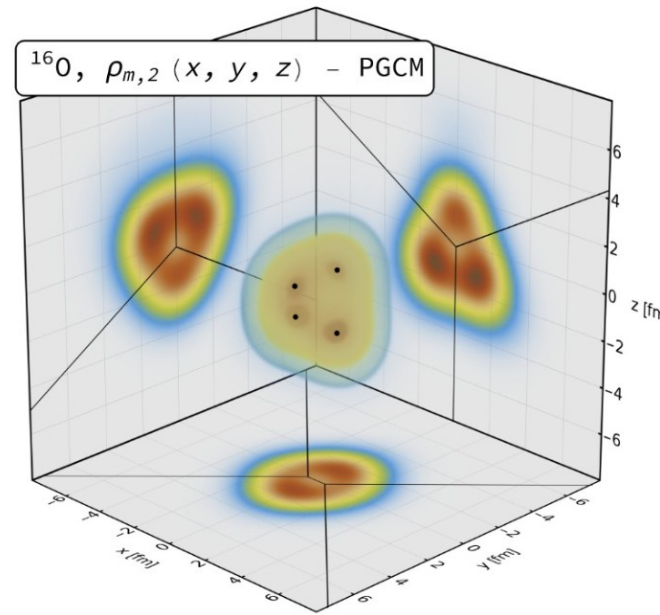
Nuclei for Future Nuclear Collisions

- **Experimental possibilities & interest**
 - Large deformed nuclei: uranium, gold, xenon
 - Smaller zirconium, rubidium, oxygen, neon



Nuclei for Future Nuclear Collisions

- **Oxygen and Neon are unique**
 - Oxygen is a double magic nucleus, since both shells are closed shell. In cluster model Tetrahedron shape.
 - Neon, has bowling pin shape, even more complicated geometry



Ancillary files (details):

- NLEFT_dmin_0.5fm_negativeweights_Ne.h5
- NLEFT_dmin_0.5fm_negativeweights_O.h5
- NLEFT_dmin_0.5fm_positiveweights_Ne.h5
- NLEFT_dmin_0.5fm_positiveweights_O.h5
- PGCM_clustered_dmin0_Ne.h5
- PGCM_clustered_dmin0_O.h5
- PGCM_uniform_dmin0_Ne.h5
- PGCM_uniform_dmin0_O.h5

[arXiv:2402.05995](https://arxiv.org/abs/2402.05995)

The shape of the oxygen

Modeling the oxygen

- **Woods-Saxon (WS)**

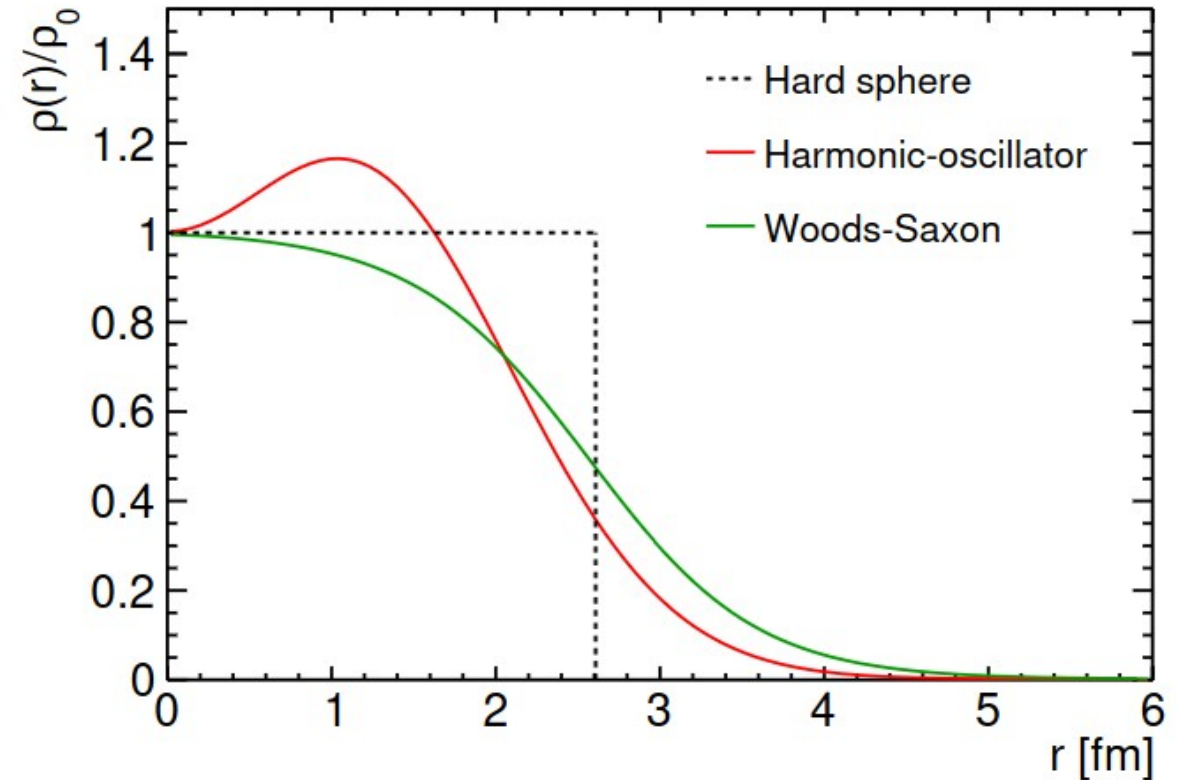
$$\rho(r) = \rho_0 \left[1 + \alpha \left(\frac{r}{a} \right)^2 \right] \exp\left(\frac{-r^2}{a^2} \right)$$

- **Harmonic oscillator (HO)**

$$\rho(r) = \frac{\rho_0 (1 + w (\frac{r}{r_0})^2)}{1 + \exp(\frac{r-r_0}{a})}$$

- **Normalization:**

$$\int \rho(r) d^3r = 4\pi \int \rho(r) r^2 dr = Ze$$

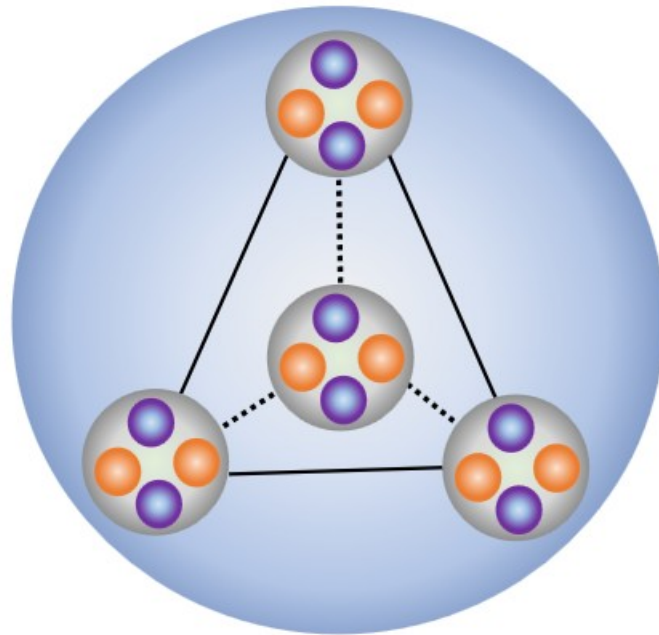


The shape of the oxygen

Nuclear structure description

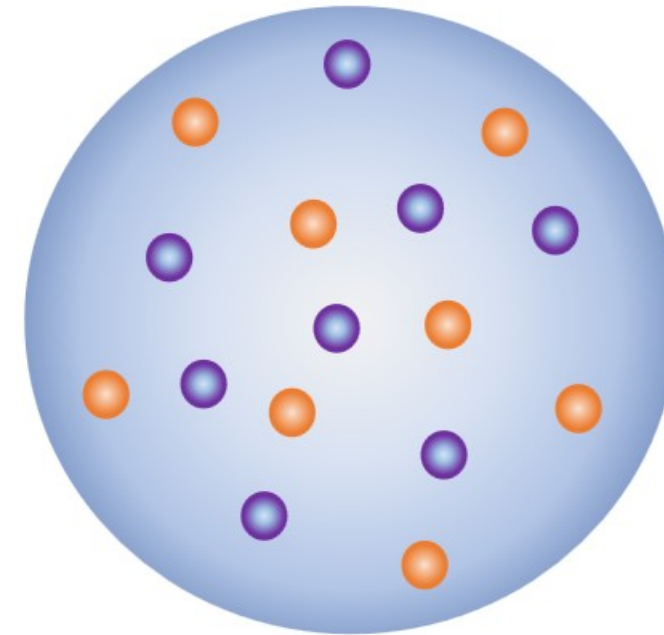
– Cluster model vs.

Non-cluster model (Woods-Saxon)



● Proton ● Neutron ● α -particle

α - clustered Oxygen (^{16}O) nucleus



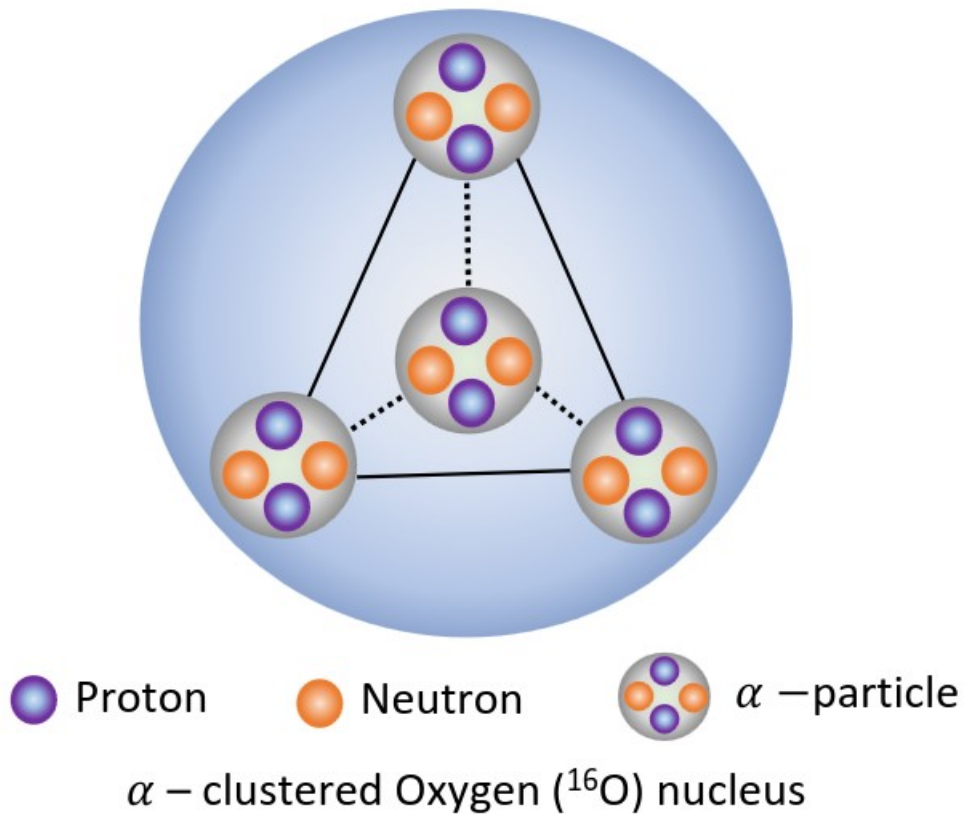
● Proton ● Neutron

Non-clustered Oxygen (^{16}O) nucleus

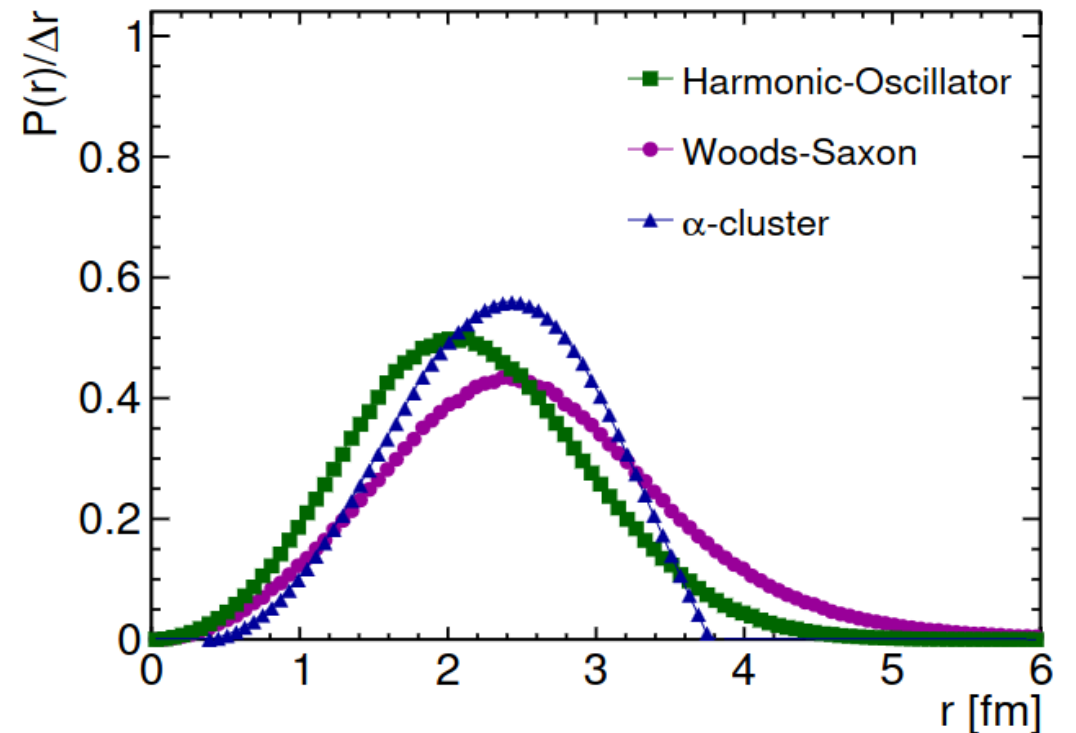
The shape of the oxygen

Nuclear structure description

- Cluster model vs WS & HO



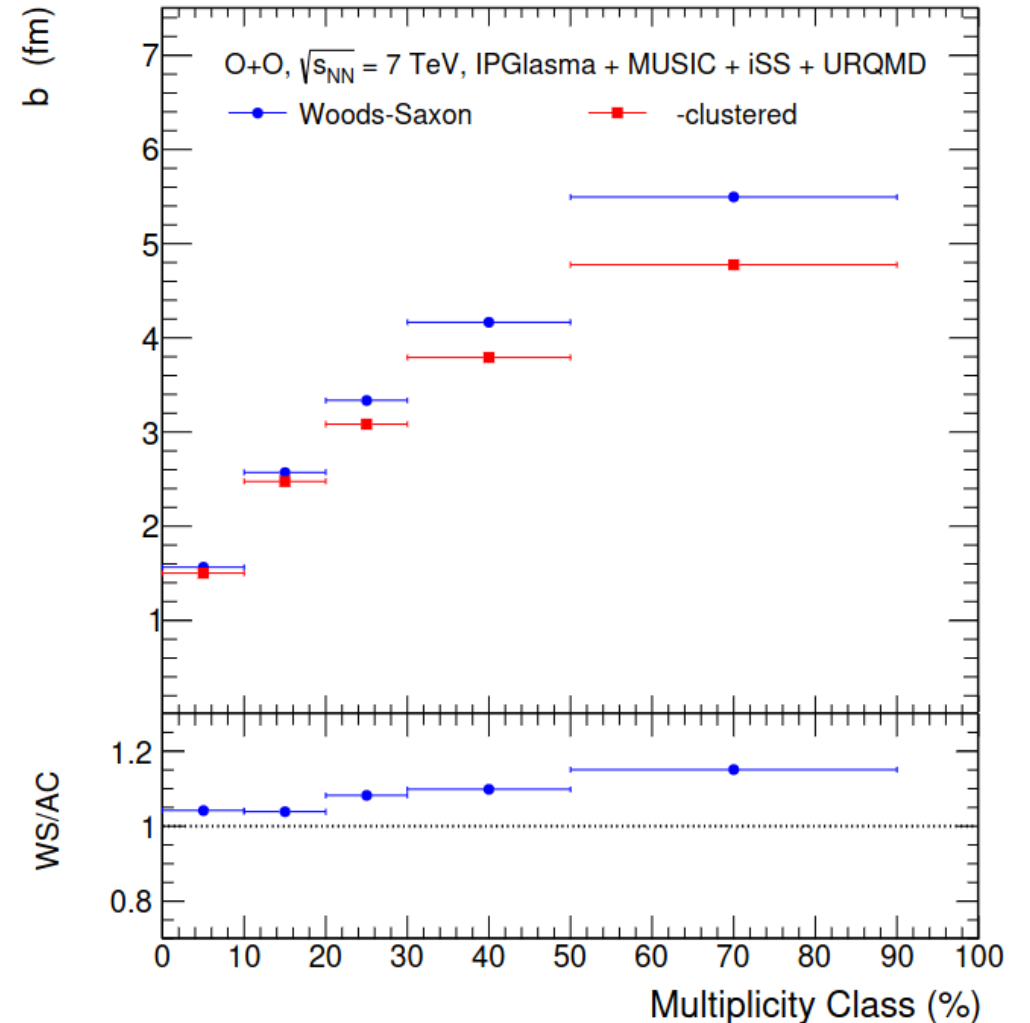
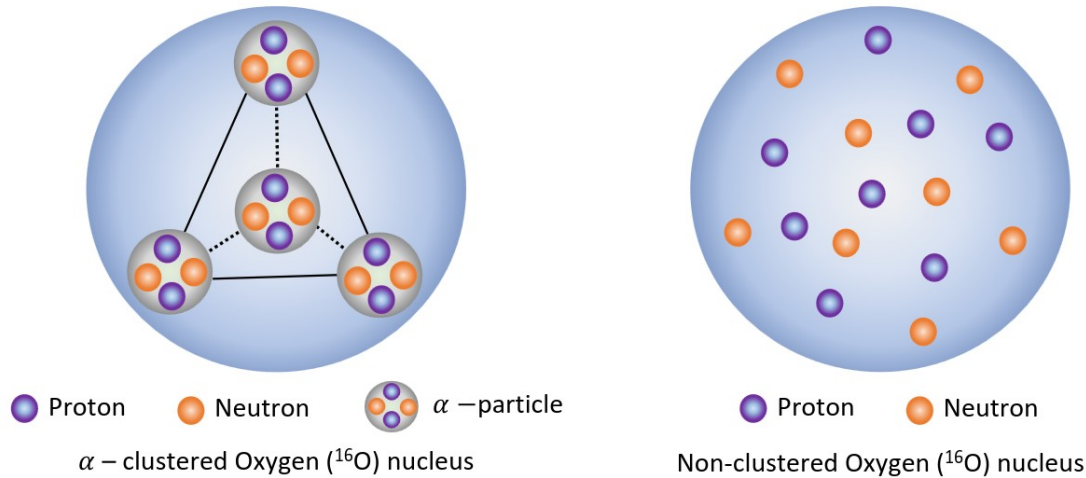
Probability of the radial position in O



The shape of the OO collision

Nuclear structure description

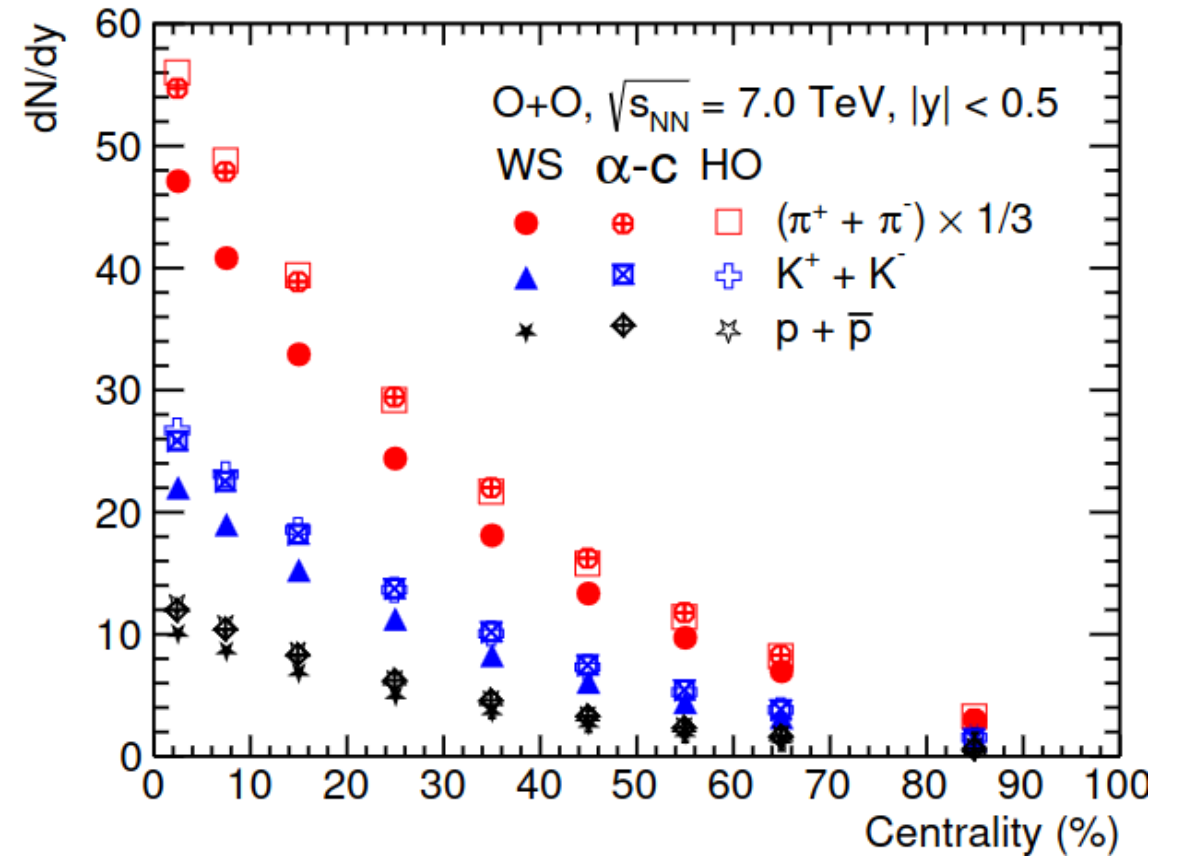
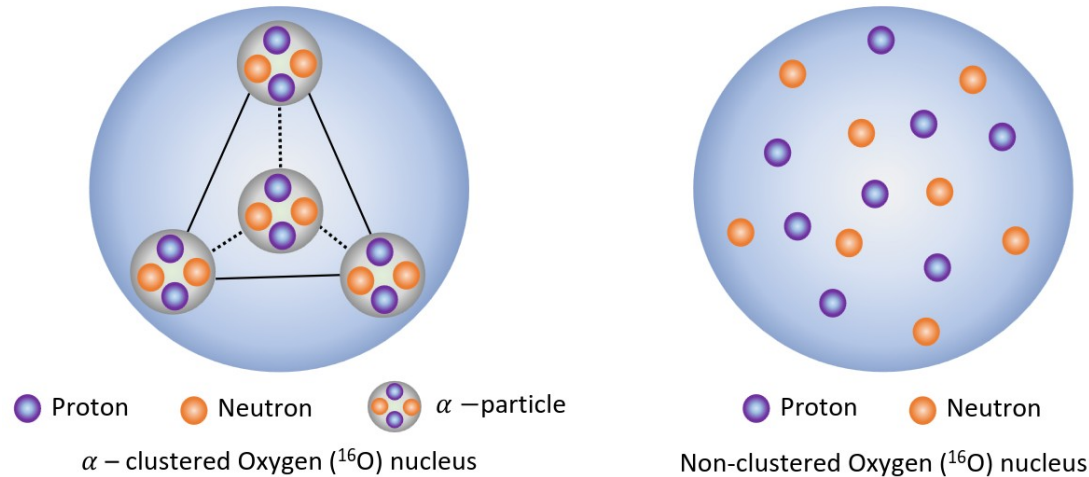
- Cluster model vs WS



The shape of the OO collision

Nuclear structure description

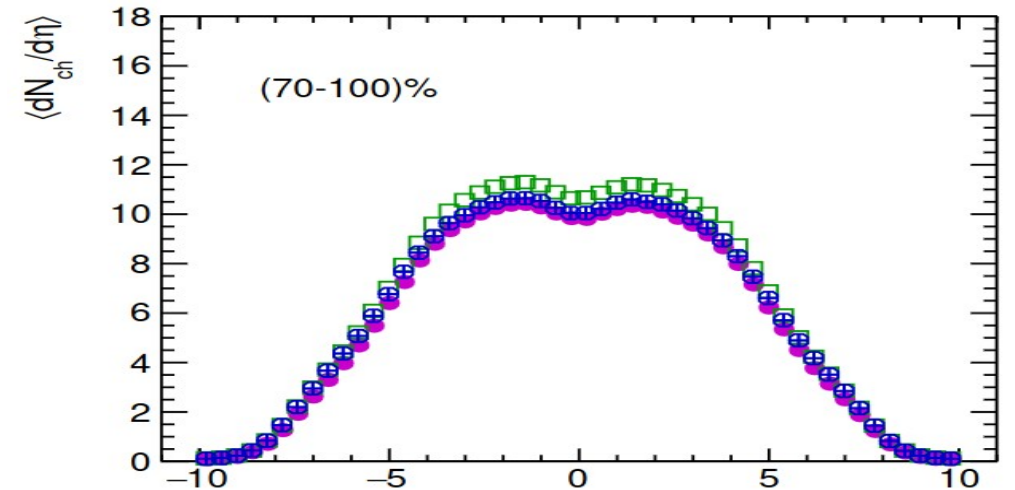
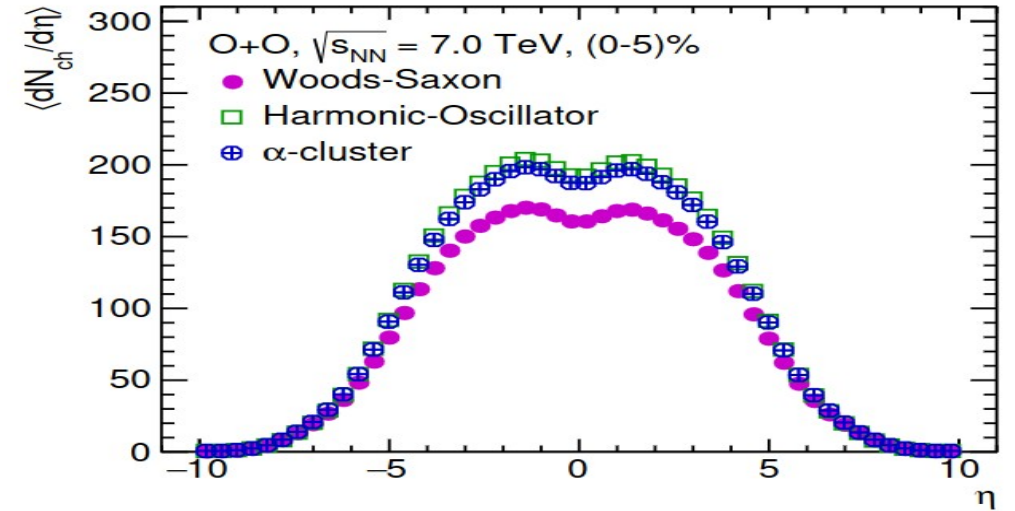
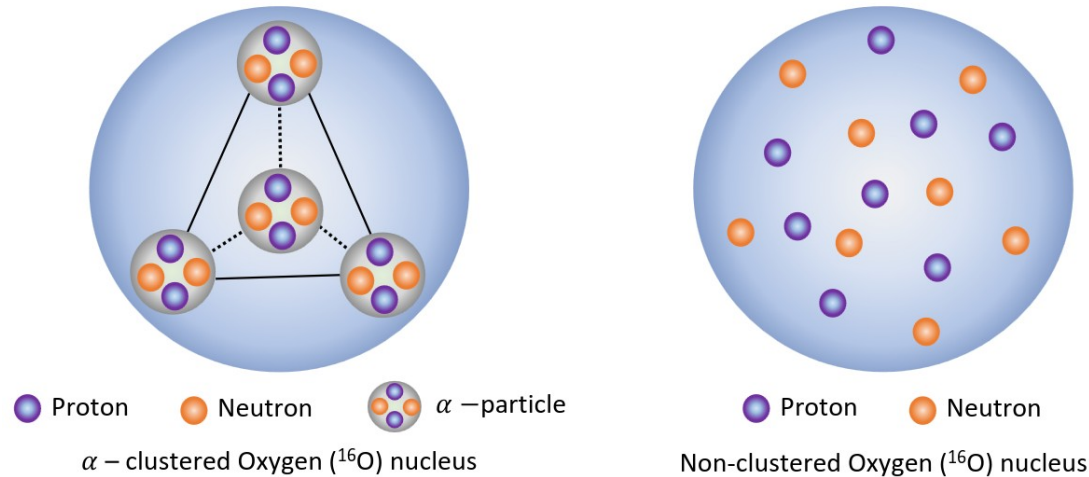
- Cluster model vs WS



The shape of the 00 collision

Nuclear structure description

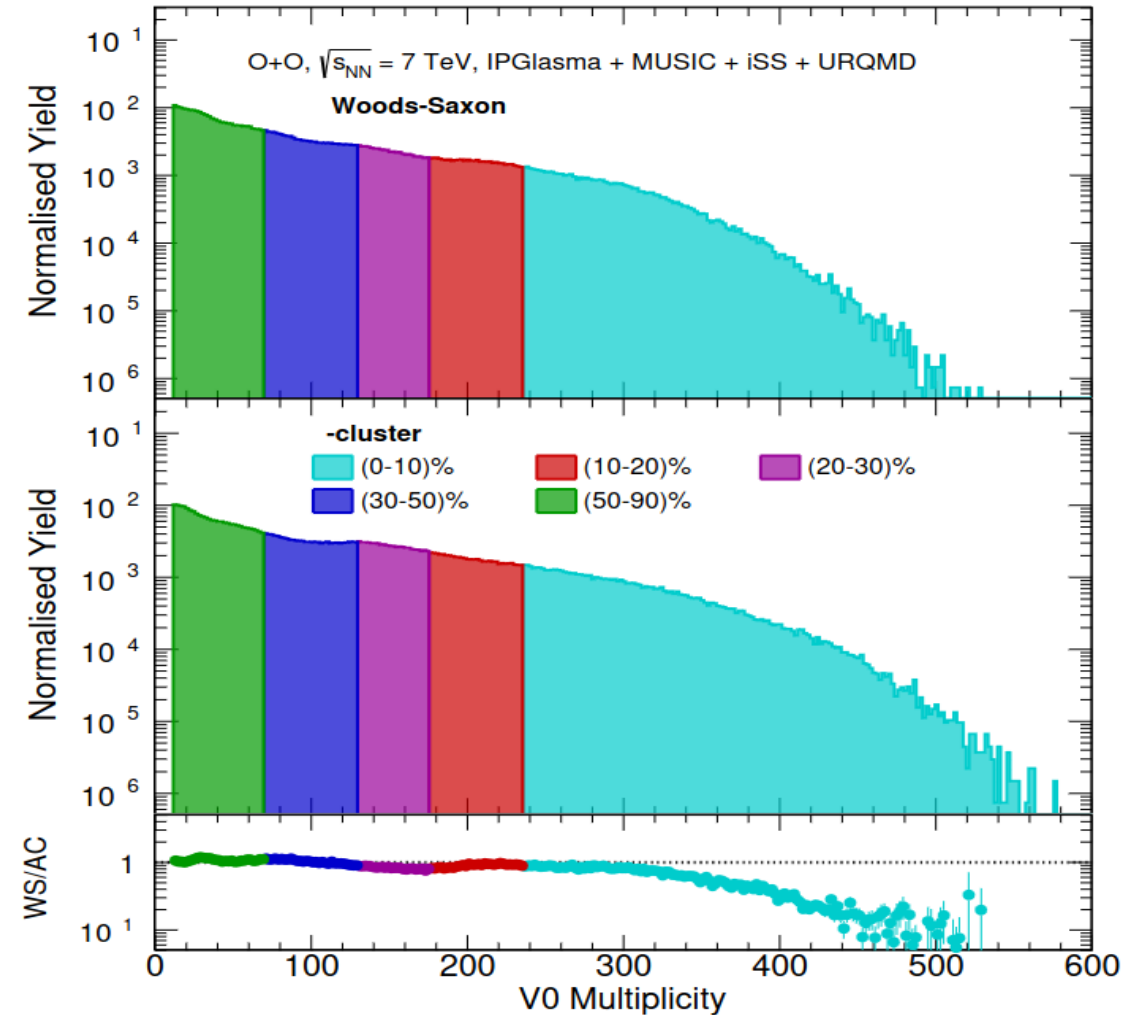
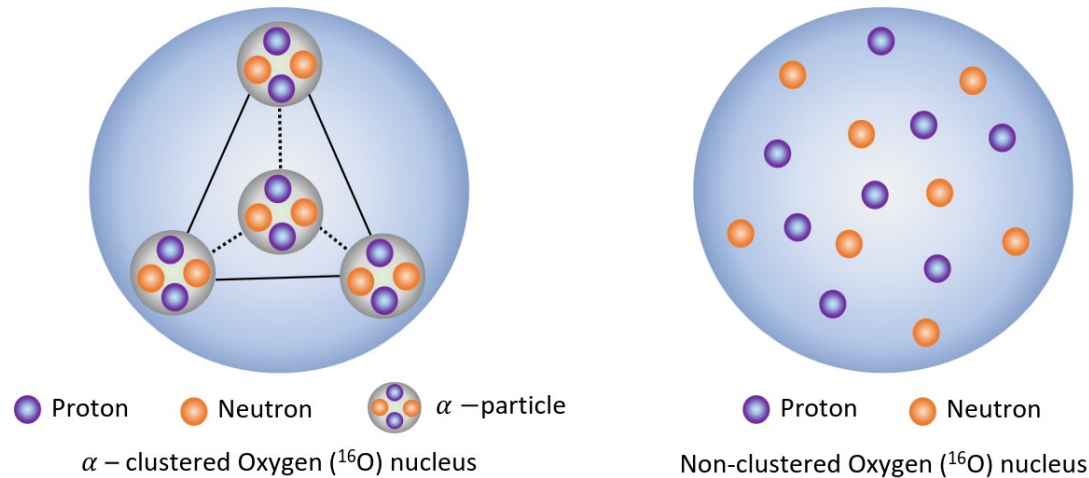
– Cluster model vs WS



The shape of the OO collision

Nuclear structure description

- Cluster model vs WS



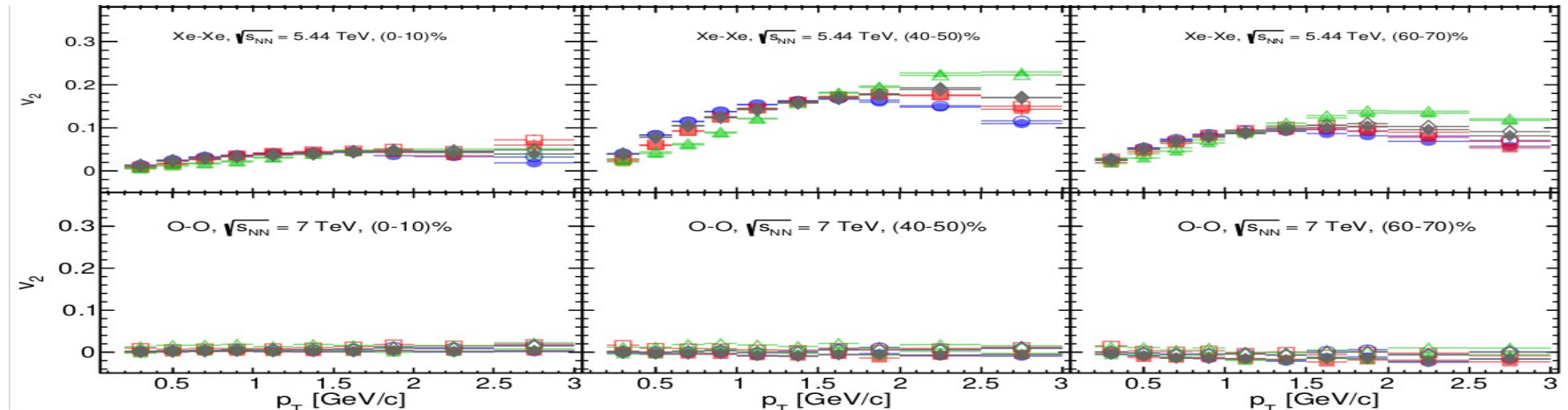
Calculating the flow in small systems

Calculating the flow

Event plane and average method

$$v_n = \langle \cos[n(\phi - \psi_n)] \rangle$$

- Need to determine the event plane, which fails for small nuclei:



The Model

- **A full hydro & Boltzmann transport with viscosity:**

- IPGlasma
- MUSIC
- iSS
- URQMD

$$\langle 2 \rangle = \frac{|Q_n|^2 - M}{M(M-1)},$$

$$\langle 4 \rangle = \frac{|Q_n|^4 + |Q_{2n}|^2 - 2 \cdot \text{Re}[Q_{2n} Q_n^* Q_n^*]}{M(M-1)(M-2)(M-3)} - 2 \frac{2(M-2) \cdot |Q_n|^2 - M(M-3)}{M(M-1)(M-2)},$$

$$c_n\{2\} = \langle\langle 2 \rangle\rangle,$$

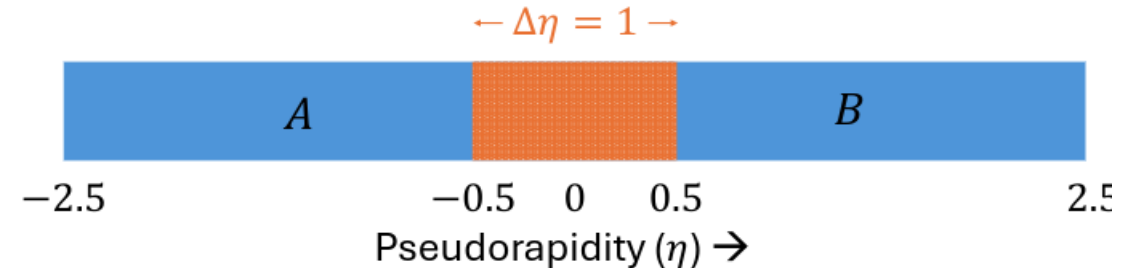
$$c_n\{4\} = \langle\langle 4 \rangle\rangle - 2 \cdot \langle\langle 2 \rangle\rangle^2.$$

$$v_n\{2\} = \sqrt{c_n\{2\}},$$

$$v_n\{4\} = \sqrt[4]{-c_n\{4\}}.$$

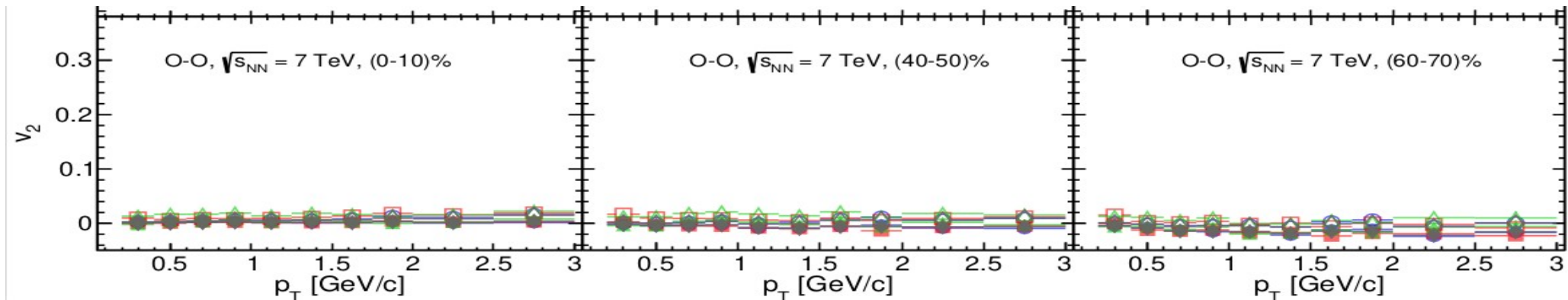
- **Kinematical settings are:**

- Energy (c.m.): 7 TeV O+O
- Pseudorapidity: $|\eta| < 2.5$
- Transverse momentum: $0.2 < p_T < 5.0$ GeV/c
- Pseudorapidity gap: $|\Delta\eta| > 1.0$

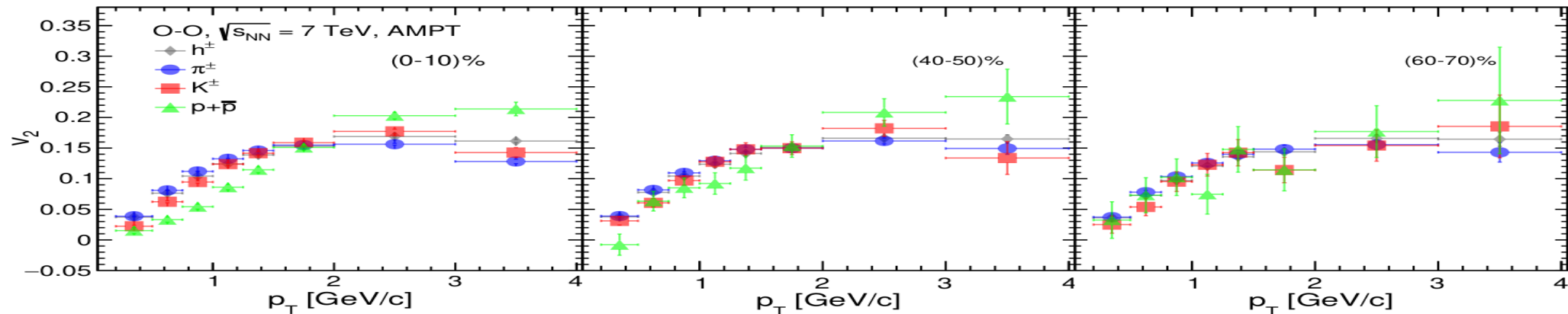


Calculating the flow

Event plane and average method



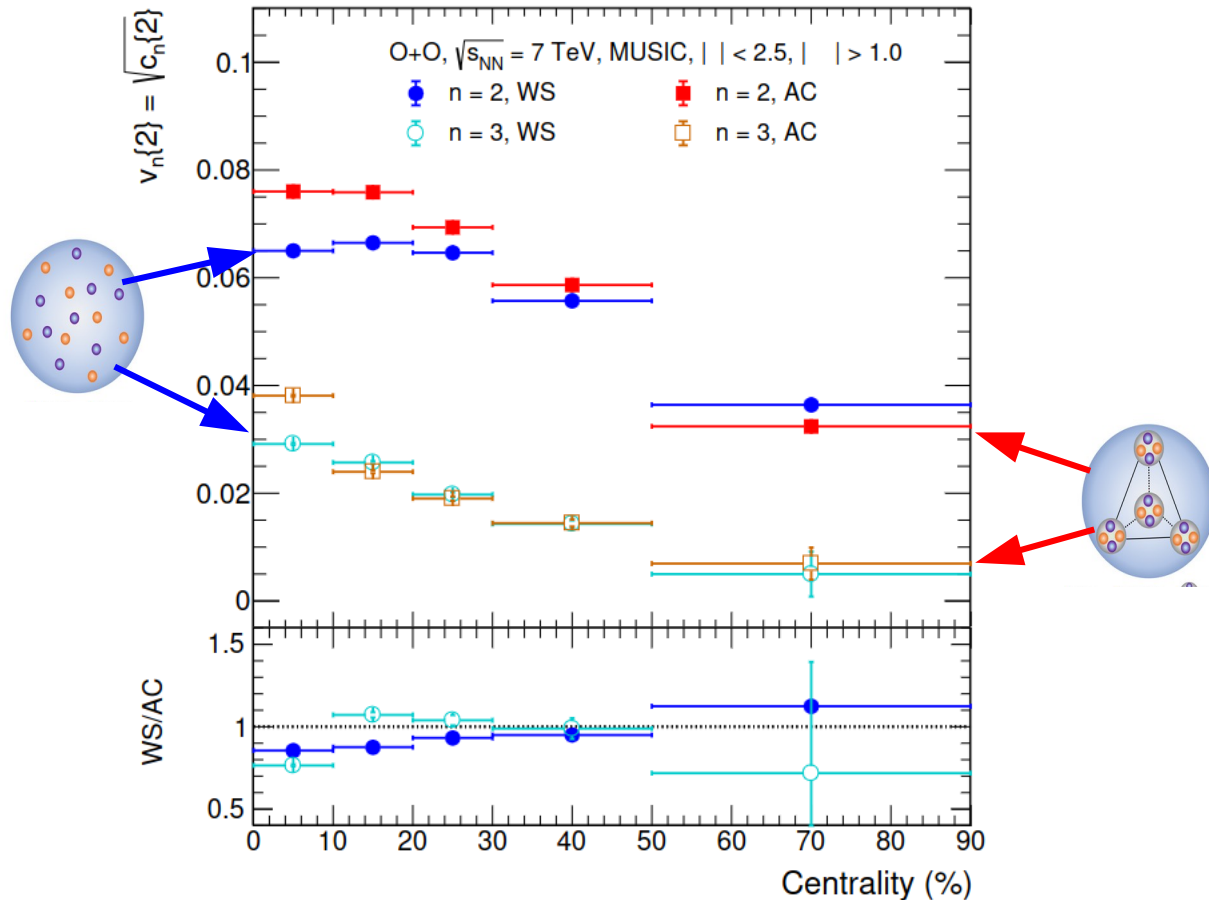
Multiparticle Q-cumulant method



Flow in oxygen-oxygen (OO)

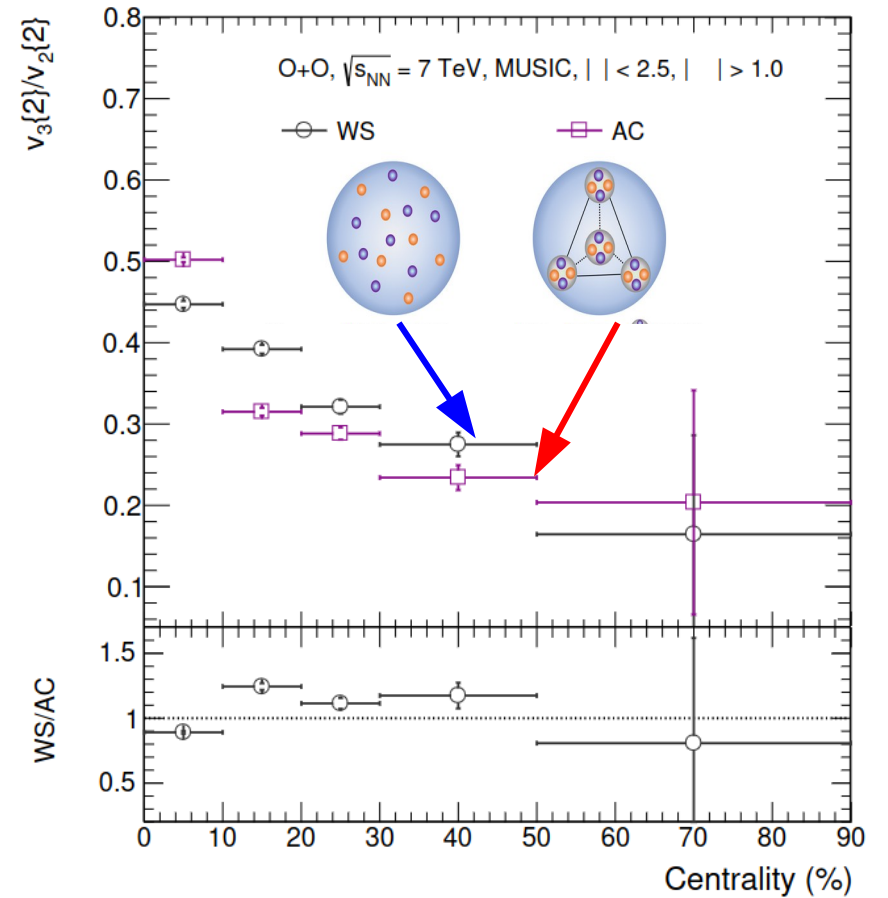
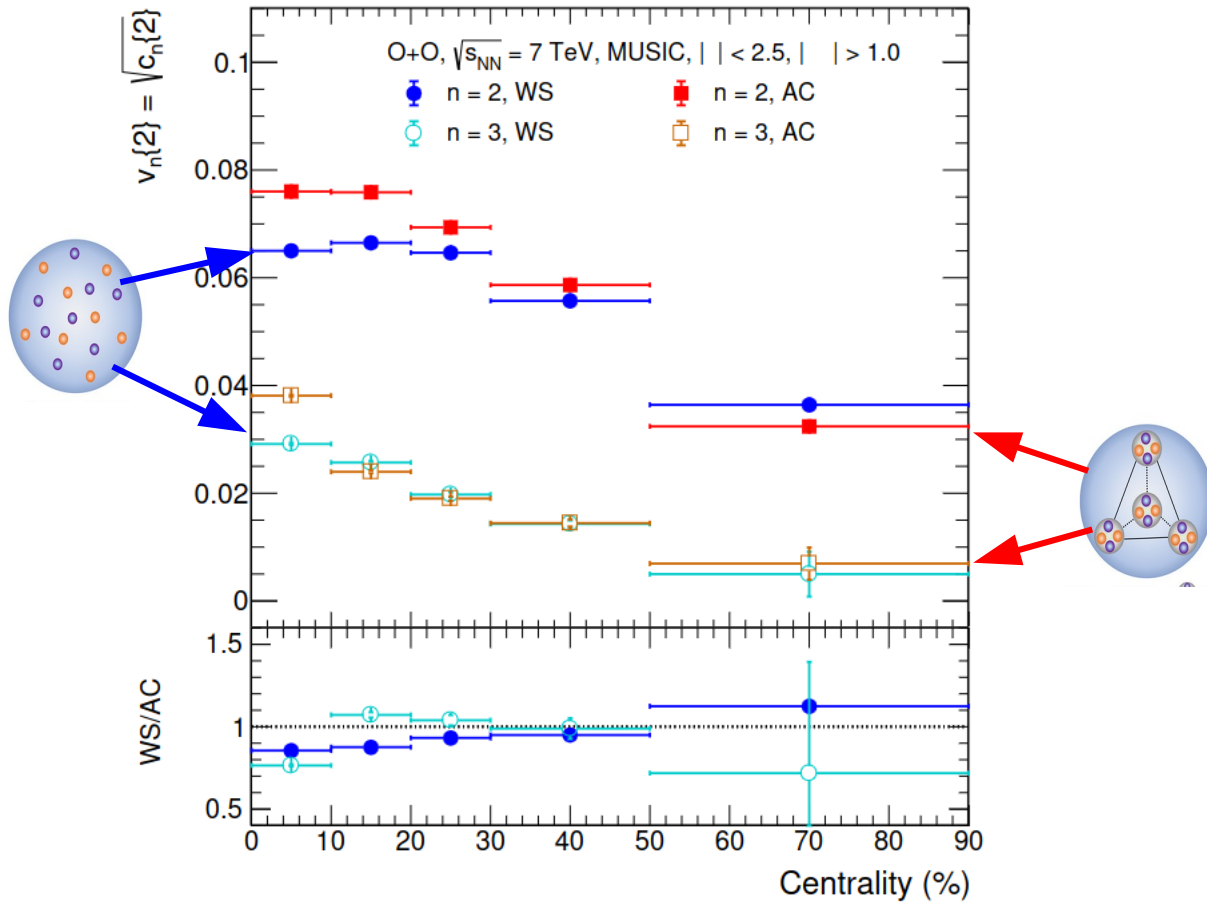
Flow components in O+O @ 7TeV

2-cummulants based calculation of v_2 & v_3



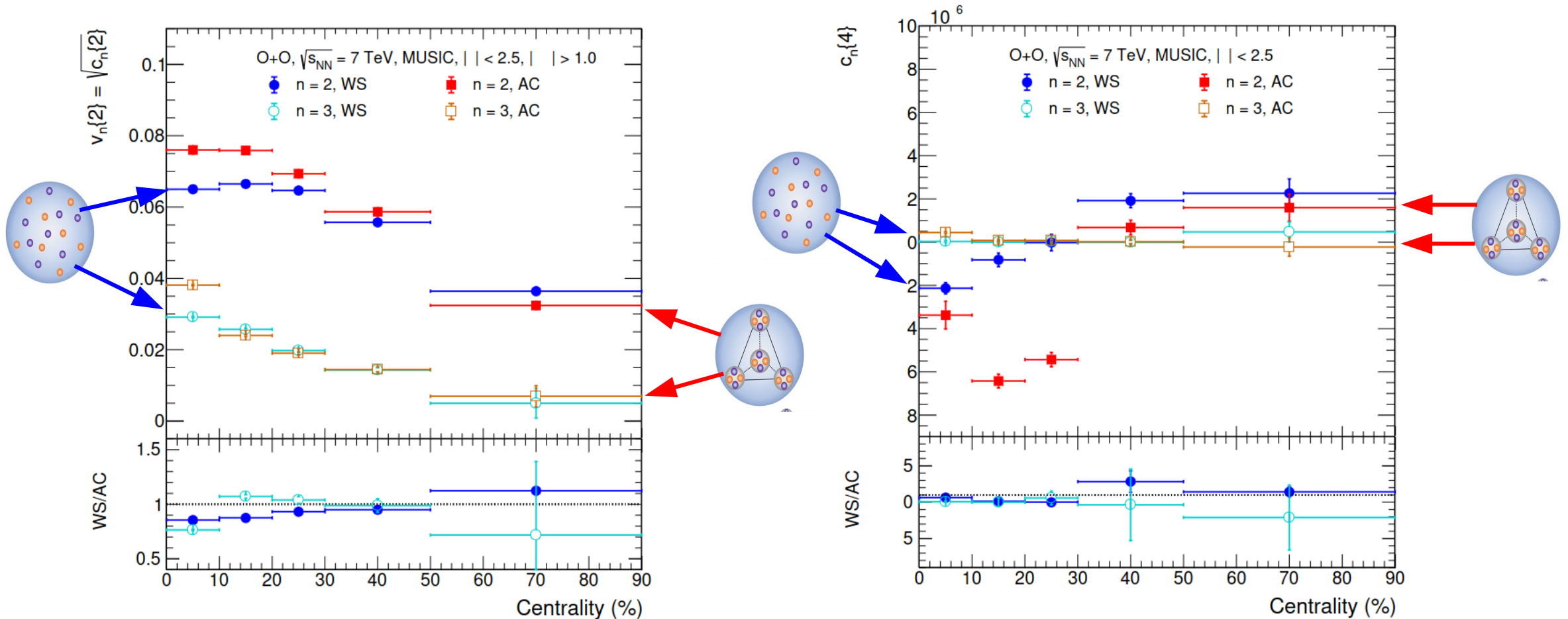
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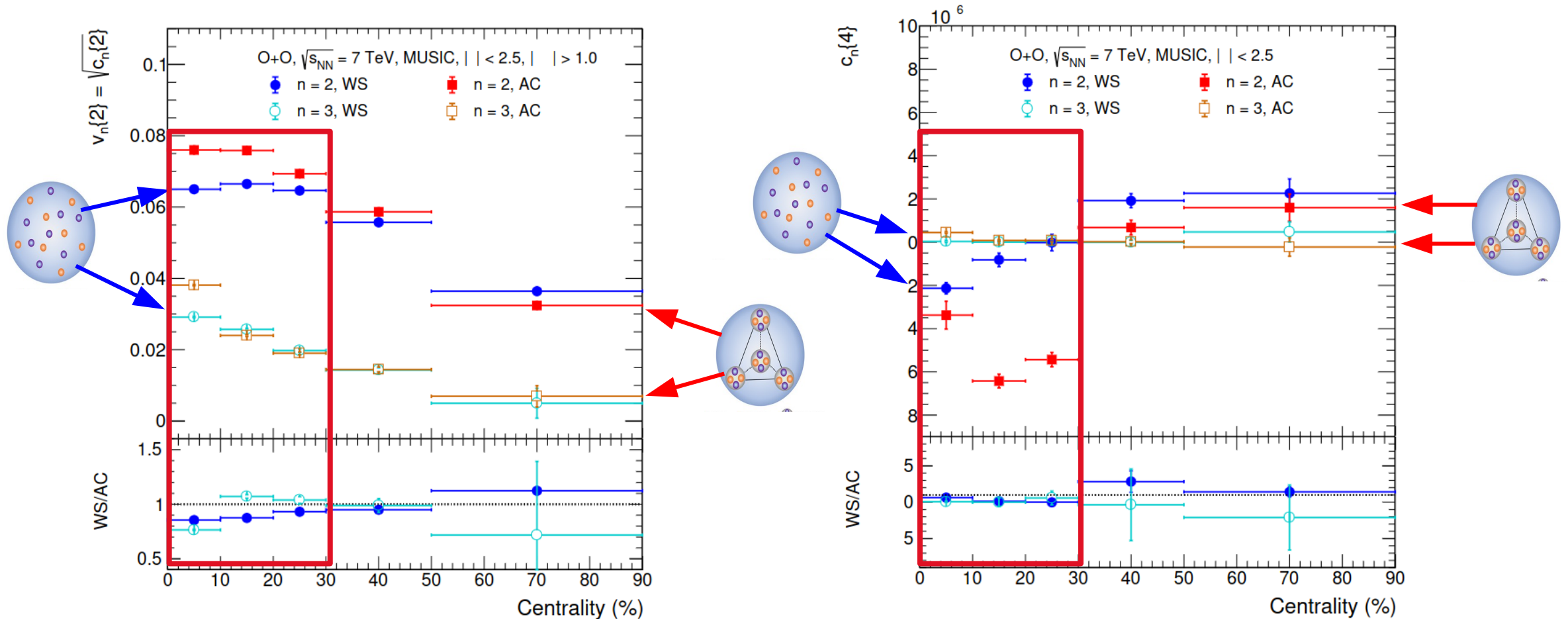
Flow components in O+O @ 7TeV

2- & 4-cummulants based v_n & c_n calculations



Flow components in O+O @ 7TeV

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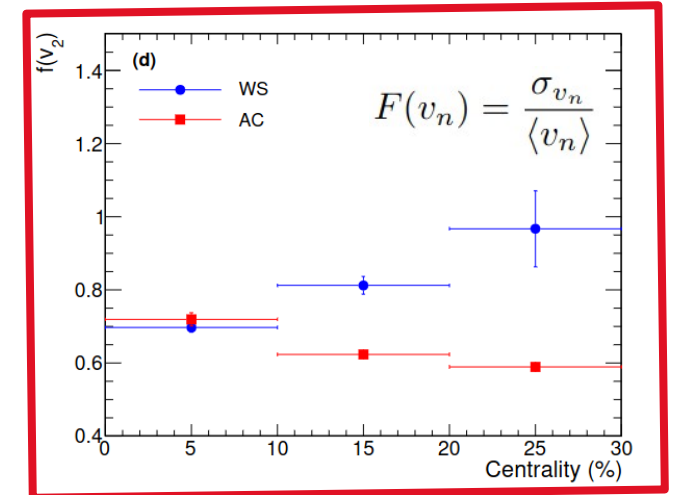
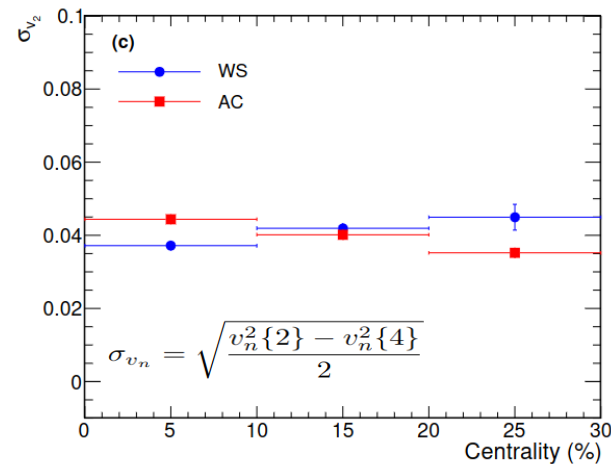
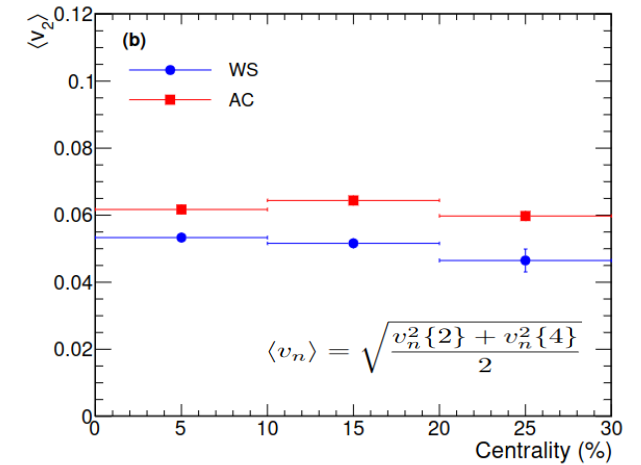
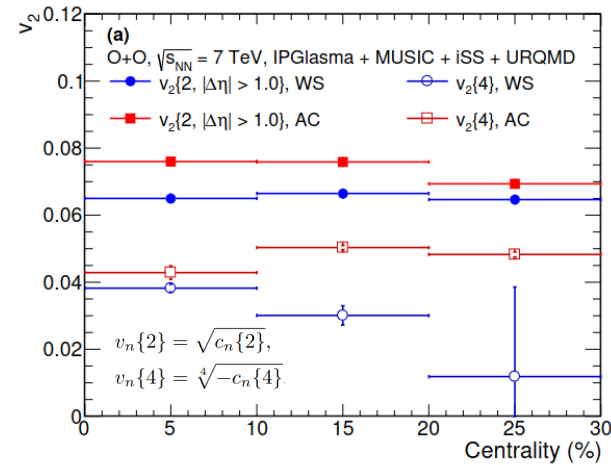


Flow components in O+O @ 7TeV

2- & 4-cummulants based calculations

- Flow and fluctuation measures changed significantly in the most central 0-30% regime
- Alpha-cluster has larger values, than Wood-Saxon profile
- Higher cummulants has higher effect at larger centrality
- Clearly visible on the relative measure:

$$F(v_n) = \frac{\sigma_{v_n}}{\langle v_n \rangle}$$



Conclusions

- **In a IPGlasma+MUSIC+iSS+URQMD = “realistic model”**
 - It is possible to calculate the flow for small system like OO
 - event plane method fails
 - 2- & 4-cummulants can be calculated for v_2
 - v_3 can not be calculated for 4-cummulant
 - Need for a kinematical cut to reduce non-flow
- **Nuclear structure has consequences on the flow**
 - Nuclear structure matters in the calculations
 - Alpha Cluster method is stronger than Woods-Saxon
 - Relevant difference is in centra O+O collisions
 - **Comparable with the size of the alpha cluster**

Thank You!

Can we prove the model' validity
in heavy-ion collisions?

Calculating the flow

Event plane and average method

$$v_n = \langle \cos[n(\phi - \psi_n)] \rangle$$

Multiparticle Q-cumulant method

- Flow vector $Q_n = \sum_{j=1}^M e^{in\phi_j}$

- The 2- and 4-particle cumulants are:

$$\langle 2 \rangle = \frac{|Q_n|^2 - M}{M(M-1)},$$

$$\langle 4 \rangle = \frac{|Q_n|^4 + |Q_{2n}|^2 - 2 \cdot \text{Re}[Q_{2n} Q_n^* Q_n^*]}{M(M-1)(M-2)(M-3)} - 2 \frac{2(M-2) \cdot |Q_n|^2 - M(M-3)}{M(M-1)(M-2)},$$

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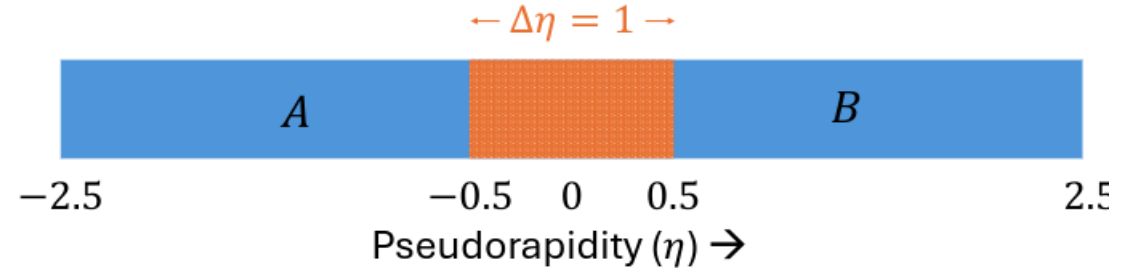
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$$v_n\{2\} = \sqrt{c_n\{2\}},$$
$$v_n\{4\} = \sqrt[4]{-c_n\{4\}}.$$

Calculating the flow

Suppressing the non-flow contribution:

- Kinematical cut: 2 sub-events, A&B are introduced, with a rapidity gap:



$$\langle 2 \rangle_{\Delta\eta} = \frac{Q_n^A \cdot Q_n^{B*}}{M_A \cdot M_B} \quad \longrightarrow \quad v_n\{2, |\Delta\eta|\}(p_T) = \frac{d_n\{2, |\Delta\eta|\}}{\sqrt{c_n\{2, |\Delta\eta|\}}}$$

- Differential flow cummulants:

$$\begin{aligned} d_n\{2\} &= \langle\langle 2' \rangle\rangle, \\ d_n\{4\} &= \langle\langle 4' \rangle\rangle - 2\langle\langle 2' \rangle\rangle\langle\langle 2 \rangle\rangle \end{aligned} \quad \longrightarrow \quad d_n\{2, |\Delta\eta|\} = \langle\langle 2' \rangle\rangle_{\Delta\eta}$$

- Mean and the fluctuations of the flow & ratio:

$$\langle v_n \rangle = \sqrt{\frac{v_n^2\{2\} + v_n^2\{4\}}{2}}$$

$$\sigma_{v_n} = \sqrt{\frac{v_n^2\{2\} - v_n^2\{4\}}{2}}$$

$$F(v_n) = \frac{\sigma_{v_n}}{\langle v_n \rangle}$$

Flow components in O+O @ 7TeV

2-cummulants based $v_n(p_T)$ calculations

Centrality

