ILLINOIS

SETUP

In this work, we explore possible signatures of the phenomenon of α -clustering in ${}^{16}O{}^{16}O$ collisions at both RHIC and LHC energies [1-3].

We model the ${}^{16}O{}^{16}O{}$ initial state with Trento [4] using three different models of the ${}^{16}O{}$ nucleus:

- Smooth Woods-Saxon profile, without constituent quarks
- Smooth Woods-Saxon profile, with constituent quarks
- Nuclear lattice effective field theory profiles containing α -clustering [5]

Hydrodynamic modeling is performed using the iEBE-VISHNU package [6]. Transport coefficients are parameterized as in [7,8]. Fig. 1 shows initial energy densities alongside maximum Knudsen and inverse Reynolds numbers for shear and bulk viscous stresses, using [9]:

 $\operatorname{Kn}_{\pi} = \tau_{\pi} \sqrt{\sigma_{\mu\nu} \sigma^{\mu\nu}}, \, \operatorname{Kn}_{\Pi} = \tau_{\Pi} \theta, \, \operatorname{Re}_{\pi}^{-1} = \sqrt{\pi_{\mu\nu} \pi^{\mu\nu}} / P, \, \operatorname{Re}_{\Pi}^{-1} = |\Pi| / P$

In general, ${}^{16}O{}^{16}O$ collisions yield large Knudsen and inverse Reynolds numbers, implying that the hydrodynamic description in ${}^{16}O{}^{16}O$ is pressed to its limits.



Figure 1: Left: Initial energy densities for three different initial-state models at RHIC (top) and LHC (bottom) energies. The same random seed is used in each panel. **Right**: Centrality class and time dependence of maximum Knudsen or inverse Reynolds number distributions in the transverse plane. The hydrodynamic evolution is clearly pushed to its limits at early times.

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FINDING *a*-clustering in Ultra-relativistic ¹⁶O¹⁶O collisions

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 α -clustering in ${}^{16}O{}^{16}O$ collisions quantitatively affect the centrality dependence of the $v_n\{k\}$ at the same order of magnitude as the inclusion of constituent quarks within individual nucleons. The hydrodynamic modeling of ${}^{16}O{}^{16}O$ collisions exhibits large gradients and viscous stresses, as parameterized by the Knudsen and inverse Reynolds numbers. Obtaining a qualitatively clean signature of α -clustering will likely require more sensitive techniques, such as the use of a flow-based principal component analyses [10].

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RESULTS



Figure 2: Left: Average transverse momentum $\langle p_T \rangle$, flow coefficients $v_n \{N\}$, and the ratio $v_2 \{4\} / v_2 \{2\}$, for all three initial-state models at RHIC energies. A total of 10^4 events were generated and all quantities are plotted as functions of centrality. **Right**: Same as the lefthand side, but at LHC energies.

We find that the effects of α -clustering are rather weak in $\langle p_T \rangle$, but detectable in flow observables, particularly in $v_2 \{k\}$. The inclusion of subnucleonic degrees of freedom in the initial state leads to an enhancement of v_2 at centralities about $\sim 15\%$, while the effects α -clustering are of a similar order of magnitude and are confined to centralities below ~ 20%. The effects on the ratio $v_2 \{4\} / v_2 \{2\}$ behave similarly, but the three initial state models considered here appear to be consistent with one another within statistical errors.

CONCLUSIONS

