

METHODS FOR SYSTEMATIC STUDY OF NUCLEAR STRUCTURE IN HIGH-ENERGY COLLISIONS

OR: CHANGING NUCLEI BY SHIFTING NUCLEONS

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Work in progress with Jean-Yves Ollitrault and Mauricio Hippert

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Intersection of nuclear structure and high-energy nuclear collisions
January 27, 2023

INTRODUCTION

- Systematic study of nuclear properties requires changing nuclear parameters and studying how observables change
- Small changes in parameters \implies small change in observables
- \implies Huge statistics required?
- No! It's possible to determine **change** in observables (or relative observable ratios) much more precisely than absolute value

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PROCEDURE USED UNTIL NOW

- 1 Choose set of nuclear parameters
- 2 Sample distribution to generate discrete nuclear configurations
- 3 Collide nuclei and compute observables
- 4 Choose new set of nuclear parameters
- 5 Generate new set of nuclear configurations from new distribution
- 6 Perform collisions and compute observables
- 7 Take ratios of observables, with independent statistical uncertainty for numerator and denominator

BETTER PROCEDURE

- 1 Generate discrete nuclear configurations **once**.
- 2 For each desired parameter set, modify configurations to obey new distribution by making small **shifts to nucleon positions**
 - Statistical uncertainty in observable ratios can be drastically reduced
 - Can be used to systematically study short-range **correlations** in addition to 1-body distribution

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ASIDE: STEP + GAUSS DISTRIBUTION

- For these numerical results, we use an alternative to a Woods-Saxon
- Not necessary, but has nice properties and makes some things easier
- Nucleon position is sum of two random vectors sampled from:
 - 1 3D step $P_s(\mathbf{x}) \sim \Theta(R_s - r)$
 - 2 3D Gaussian $P_g(\mathbf{x}) \sim e^{-\frac{r^2}{2w^2}}$
- Rough rule of thumb:

$$R_s(R, a) \simeq R \left[1 + 1.5 \left(\frac{a}{R} \right)^{1.8} \right]$$

$$w(R, a) \simeq 1.83 a$$

$$\rho_c(\mathbf{x}) = \int P_s(\mathbf{z}) P_g(\mathbf{x} - \mathbf{z}) d^3z$$

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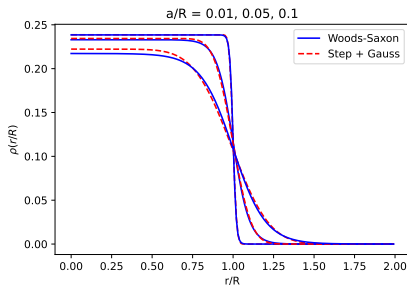
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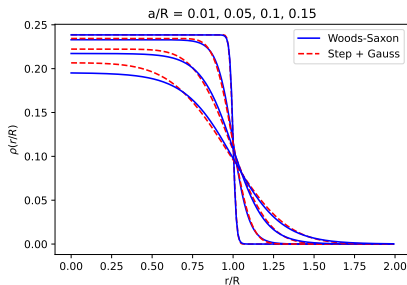
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STEP+GAUSS DISTRIBUTION ADVANTAGES

BENEFITS OF STEP+GAUSS

- Can directly modify Woods-Saxon parameters R, a without using the to-be-described methods
- No need for acceptance/rejection
- Trivial relation between point nucleon density and charge density
- Nice analytic properties — smooth at origin

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CHANGING NUCLEAR SHAPE

- 1-body nucleon distribution parameterized as

$$\rho(r) \propto \frac{1}{1 + e^{\frac{r-R}{a}}}$$
$$\tilde{\rho}(r, \theta, \phi) \propto \frac{1}{1 + e^{\frac{r-R-R\sum_{\ell,m}\beta_{\ell,m}Y_{\ell,m}}{a}}} = \rho(r - R\sum_{\ell,m}\beta_{\ell,m}Y_{\ell,m})$$

- Define continuous parameter t that takes you from spherical ($t = 0$) to desired deformed distribution ($t = 1$)

$$\rho(\vec{x}, t) \equiv \rho(r - t\sum_{\ell,m}R\beta_{\ell,m}Y_{\ell,m})$$

- Idea: change nuclear properties by shifting the position of nucleons

$$\implies \frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \vec{v}) = 0$$

- Start with uncorrelated nucleons satisfying $\rho(r)$, end with uncorrelated nucleons satisfying $\rho(r - R\sum_{\ell,m}\beta_{\ell,m})$

ANGULAR DEFORMATION

$$\rho(\vec{x}, t) \equiv \rho(r - t \sum_{\ell, m} R\beta_{\ell, m} Y_{\ell, m})$$

$$0 = \frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \vec{v})$$

- One solution (at $t = 0$):

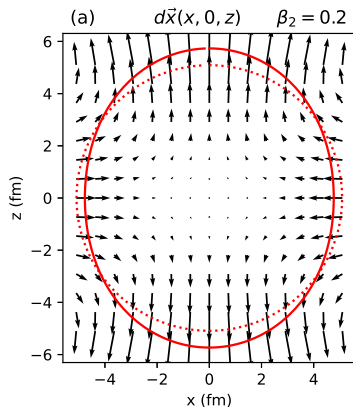
$$\vec{v} = \nabla \Phi(\vec{x})$$

$$\Phi = \sum R\beta_{\ell, m} f_{\ell, m}(r) Y_{\ell, m}$$

$$0 = f''_{\ell, m} + f'_{\ell, m} \left(\frac{2}{r} + \frac{\rho'}{\rho} \right) - \frac{\ell(\ell + 1)}{r^2} f_{\ell, m} - \frac{\rho'}{\rho}$$

$$0 = f_{\ell, m}(r \rightarrow 0)$$

$$1 = f'_{\ell, m}(r \rightarrow \infty)$$



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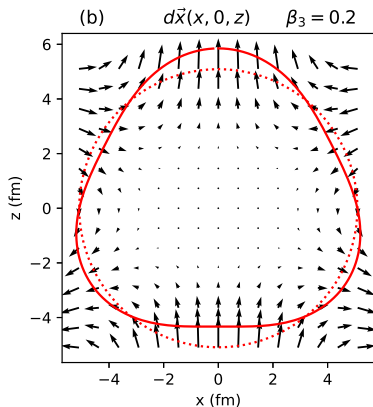
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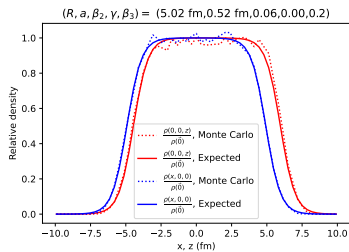
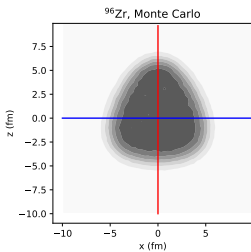
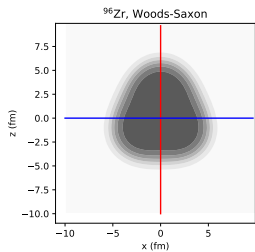
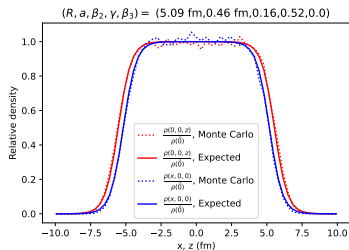
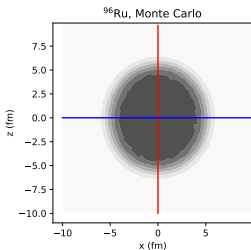
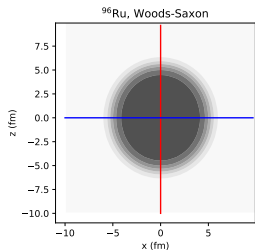
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NUMERICAL RESULTS (100K NUCLEI)



OUTLINE

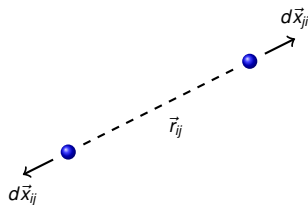
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SHORT-RANGE CORRELATIONS

- Short-range interactions cause particles to be correlated

$$\rho_2(\vec{x}_1, \vec{x}_2) = \rho(\vec{x}_1)\rho(\vec{x}_2) [1 + C(\vec{r}_{12})]$$

- Idea: induce correlation C from uncorrelated set by shifting particles



$$d\vec{x}_i = \sum_{j \neq i} d\vec{x}_{ij} = \sum_{j \neq i} \frac{1}{2} (\tilde{r}_{ij} - r_{ij}) \hat{r}_{ij}$$

FINDING $dr = \tilde{r} - r$

- Conserve pairs:

$$\int_0^r d^3 r' = \int_0^{\tilde{r}} d^3 r' (1 + C(\vec{r}'))$$

- Invert relation to solve for \tilde{r}
- For simplicity, we implemented a step function correlation function with variable length $C_{\text{length}} \geq 0$ and strength $C_{\text{strength}} \geq -1$

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$$(r^3 - \tilde{r}^3) = 3 \int_0^{\tilde{r}} dr' r'^2 C(r')$$

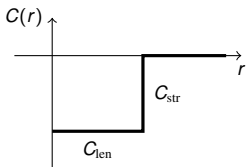
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VALID CORRELATION FUNCTIONS

- Note that the number of pairs is fixed:

$$\begin{aligned} \rho(\vec{x}_1)\rho(\vec{x}_2) [1 + C(\vec{r}_{12})] &= \rho_2(\vec{x}_1, \vec{x}_2) \\ \implies \int d^3x_1 d^3x_2 \rho(\mathbf{x}_1)\rho(\mathbf{x}_2)C(\vec{r}_{12}) &= 0 \end{aligned}$$

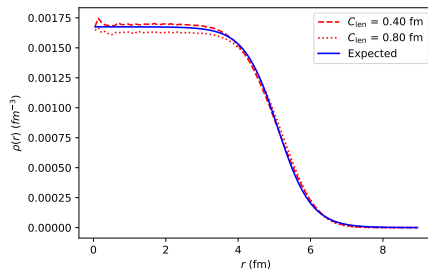
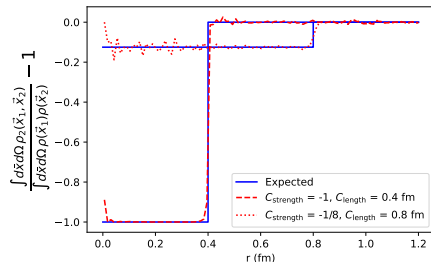
- Respecting sum rule important for maintaining fixed 1-body distribution
- If nominal short-range correlation doesn't satisfy, we add constant

$$\begin{aligned} C(r) &= C_{\text{short}}(r) + C_{\infty} \\ C_{\infty} &\simeq -C_{\text{vol}} \int d^3x \rho(\mathbf{x})^2 \end{aligned}$$

ADVANTAGES

Besides statistical speedup:

- Can study correlation of arbitrary shape – not just exclusion distance
- No problems with triaxial nuclei
- Better control over 2-body and 1-body distributions

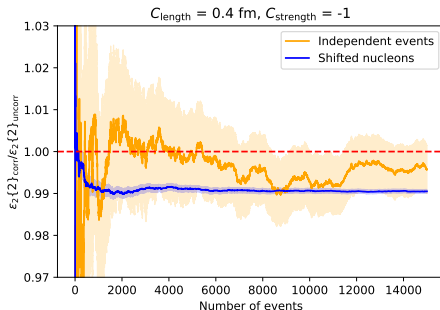


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PRELIMINARY BENCHMARKS

HOW MUCH BENEFIT CAN YOU GET?

- Simple benchmark: participant Glauber model at $b = 0$. Ratios of eccentricities $\varepsilon_n\{2\}$, (baseline)/(baseline + change in 1 parameter)
- Compare our method to naive method with independent nuclei
- Question: If I get a statistical uncertainty with N events in the naive case, how many events N/F do I need in order to get at least as small statistical uncertainty, when using this method? F = improvement factor.



PRELIMINARY BENCHMARKS

Par.	Param. Change	$\varepsilon_2\{2\}$ Change	Improv. Factor	Avg. Shift
C_{len}^3	(0.2 fm) ³	0.13%	2900	0.002 fm
C_{len}^3	×2	0.27%	1100	0.005 fm
C_{len}^3	×4	0.53%	350	0.009 fm
C_{len}^3	(0.4 fm) ³	1.1%	180	0.017 fm
C_{len}^3	×2	2.0%	98	0.032 fm
C_{len}^3	×4	3.8%	54	0.059 fm
C_{len}^3	(0.8 fm) ³	7.3%	25	0.11 fm
C_{len}^3	×2	14%	13	0.19 fm

TAKEAWAYS

- Significant improvement possible
- Main limitation: nucleon shift can change participant ↔ spectator
- Larger differences in nuclei ⇒ reduced improvement factor
- Exact numbers will depend on centrality, model, etc.

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Par.	Param. Change	$\varepsilon_n\{2\}$ Change	Improv. Factor	Avg. Shift
β_2	0.005	0.02%	170	0.008 fm
β_2	0.01	0.10%	100	0.02 fm
β_2	0.02	0.39%	42	0.03 fm
β_2	0.05	2.3%	12	0.08 fm
β_2	0.1	8.8%	4.7	0.17 fm
β_2	0.2	31%	2.1	0.33 fm
β_3	0.01	0.05%	79	0.01 fm
β_3	0.05	1.6%	13	0.06 fm
β_3	0.1	6.3%	5.0	0.12 fm
β_3	0.2	23%	2.2	0.25 fm

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SUMMARY

- Can significantly reduce statistical demands by correlating statistical fluctuations — change nuclear properties by shifting nucleons
- Allows for efficient systematic study of nuclear structure
- Allows for arbitrary Woods-Saxon parameters ($R, a, \{\beta_{\ell,m}\}$) and short-range correlation function $C(\vec{r})$
- Statistical improvements depend on context — better improvement for smaller changes in nuclei — but always an improvement over standard method
- Article and Python code to generate nuclei to appear soon
- Warning: must synchronize other fluctuations in collision model — impact parameter, orientation of nuclei, etc.

EXTRA SLIDES