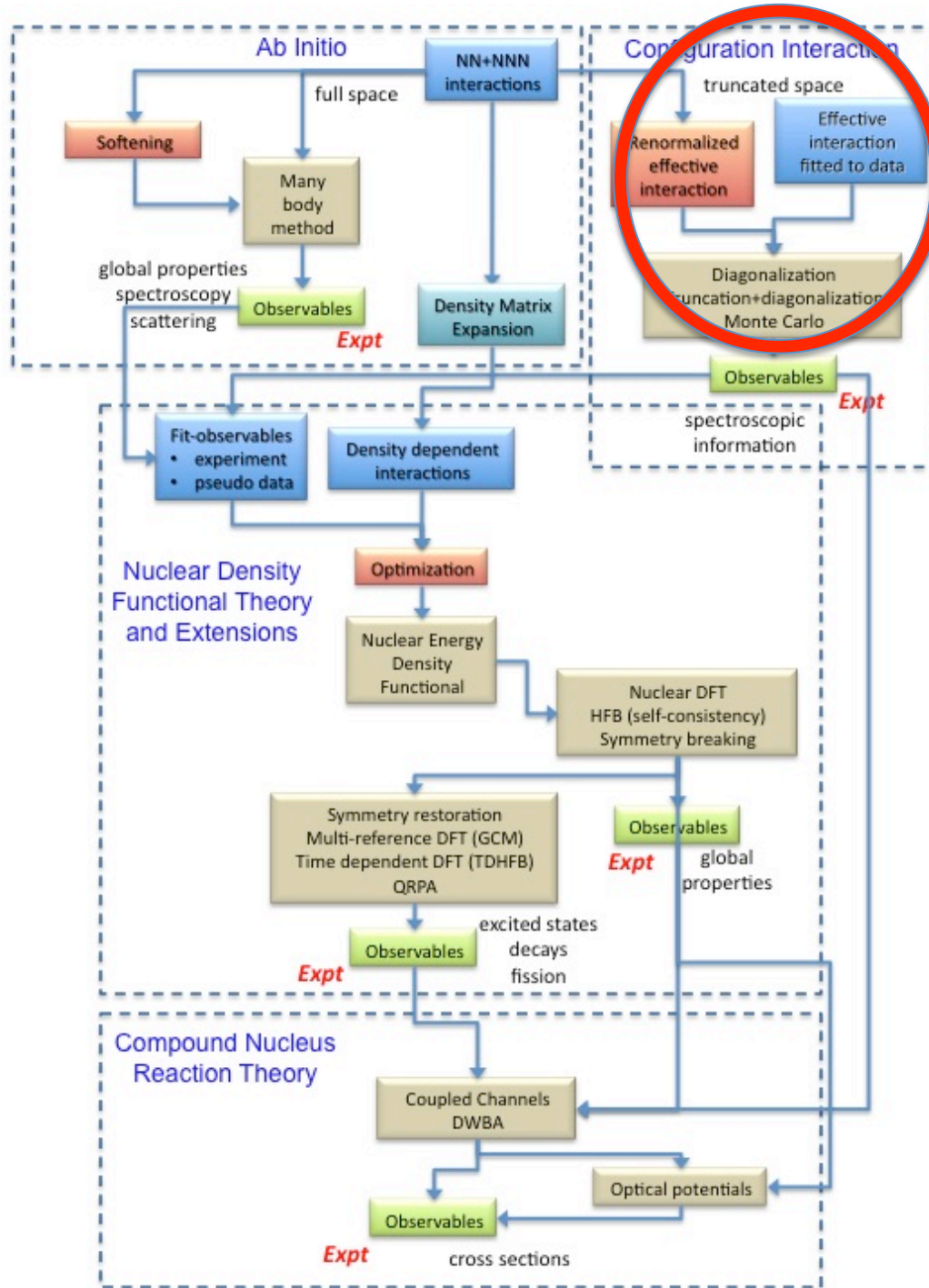


*UNEDF Annual Meeting 2010*

# Making Effective Interactions More Effective

Applications of the `BIGSTICK CI` code  
to 2-species (up/down) fermions at unitarity:

- (1) General effective interaction
- (2) Center of mass without exact factorization



## What goes into a shell-model CI calculation

Configuration-interaction (CI) calculations in a shell-model basis:

Solve  $\hat{H}|\Psi\rangle = E|\Psi\rangle$  in a Slater determinant basis:  $|\Psi\rangle = \sum_{\alpha} c_{\alpha}|\alpha\rangle$

where each Slater determinant is built from single-particle states with good angular momentum  $j, m$  (but arbitrary radial wavefunction).

The Hamiltonian is input in second quantization:

$$\hat{H} = \sum \varepsilon_a \hat{n}_a + \frac{1}{4} \sum V_{abcd} \hat{a}_a^+ \hat{a}_b^+ \hat{a}_d \hat{a}_c$$

The single-particle energies and two-body matrix elements are computed externally to the CI code and read in as **a file of numbers**.

**The BIG QUESTION:**

***What are these numbers? How do we get them?***

## Introduction: Conquering Empirical Interactions

**Naive** use of *ab initio* interactions fail to describe data.

1. “Hard core” makes calculations troublesome.
2. Tractable model space is too small.
3. Need 3 body forces, but only use 2-body forces.

## Introduction: Conquering Empirical Interactions

One creates a renormalized *effective interaction* which implicitly account for the sums to high-momentum states, e.g., Brueckner G-matrices.

**Modern approaches use unitary transformations**

A renormalized effective interaction is numerically more tractable, but still doesn't give the right spectrum.

Therefore one often tweaks a renormalized realistic interaction in order to make it agree better with data.

cf Brussaard and Glaudemans, Ch.7

more recent: Brown and Richter, PRC 74 034315 (2006) (“USDA”, “USDB”) and others...

## Introduction: Conquering Empirical Interactions

Therefore one often tweaks a renormalized realistic interaction in order to make it agree better with data.

Given a Hamiltonian  $\mathbf{H}$ , compute some set of levels (over many nuclei)  $\{|\alpha\rangle\}$  with energies  $E_\alpha$ ; let  $E_\alpha^0$  be the experimental (target) energies.

Want to minimize  $\chi^2 = \sum_{\alpha} (E_\alpha^0 - E_\alpha)^2$

Let  $\hat{H} \rightarrow \hat{H} + \sum_i \delta c_i \hat{H}_i$

and  $E_\alpha \rightarrow E_\alpha + \sum_i \delta c_i \frac{\partial E_\alpha}{\partial c_i}$

Hellmann-Feynman theorem:

$$\frac{\partial E_\alpha}{\partial c_i} = \langle \alpha | \hat{H}_i | \alpha \rangle$$

## Introduction: Conquering Empirical Interactions

$$\frac{\partial \chi^2}{\partial \delta c_i} = 0 \quad \longrightarrow \quad \sum_j \left( \sum_\alpha \frac{\partial E_\alpha}{\partial c_i} \frac{\partial E_\alpha}{\partial c_j} \right) \delta c_j = \sum_\alpha \frac{\partial E_\alpha}{\partial c_i} (E_\alpha^0 - E_\alpha)$$

This has the form  $\mathbf{B}^T \mathbf{B} \vec{c} = \mathbf{B}^T \delta \vec{E}$

Formally the solution is  $\vec{c} = \left( \mathbf{B}^T \mathbf{B} \right)^{-1} \mathbf{B}^T \delta \vec{E}$  but

$B_{\alpha i} = \frac{\partial E_\alpha}{\partial c_i}$  may be singular or nearly so

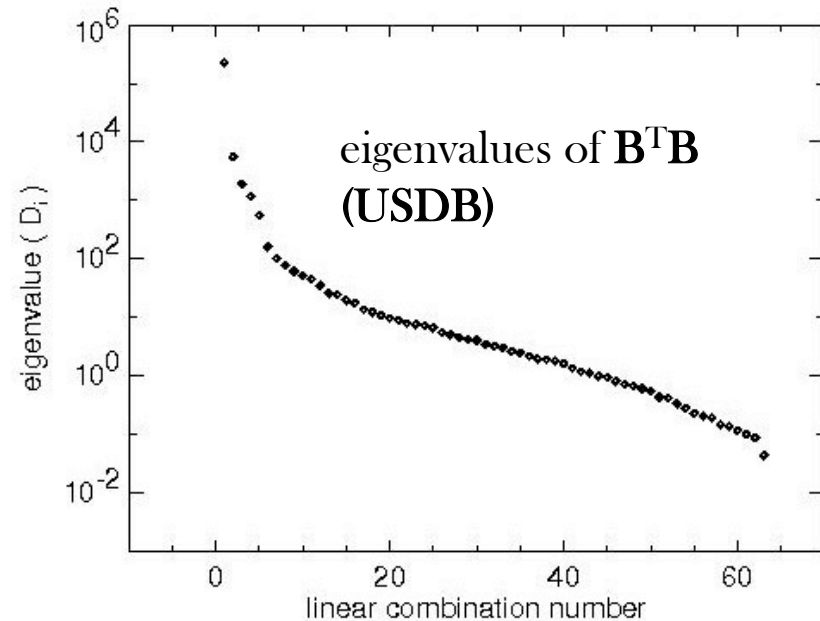
Thus one does a singular value decomposition—find the eigenvalues of  $\mathbf{B}^T \mathbf{B}$  and truncate.

## Part 2: SVD analysis of random and non-random interactions



Let's review: Given an interaction and a set of states  $\{|\alpha\rangle\}$ , one can use the Hellmann-Feynman theory to compute the sensitivity of the spectrum to perturbations in the Hamiltonian

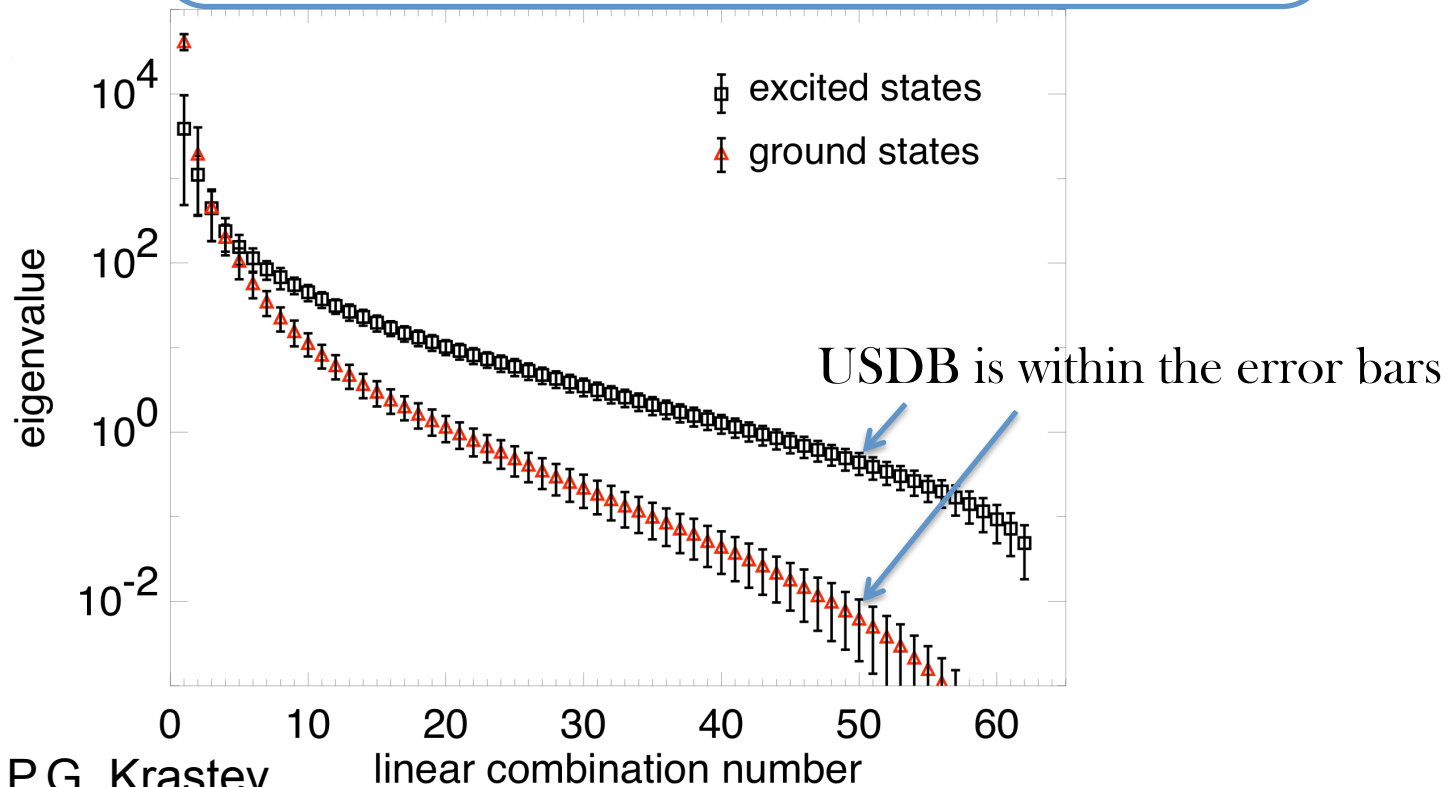
$$B_{\alpha i} = \frac{\partial E_{\alpha}}{\partial c_i} = \langle \alpha | \hat{H}_i | \alpha \rangle$$



## Part 2: SVD analysis of random and non-random interactions



Is there something *special* about the nuclear interaction? What about other interactions? Suppose we take a *random* interaction?



Johnson and P.G. Krastev,  
PRC **81**, 054303 (2010).

## Interlude:

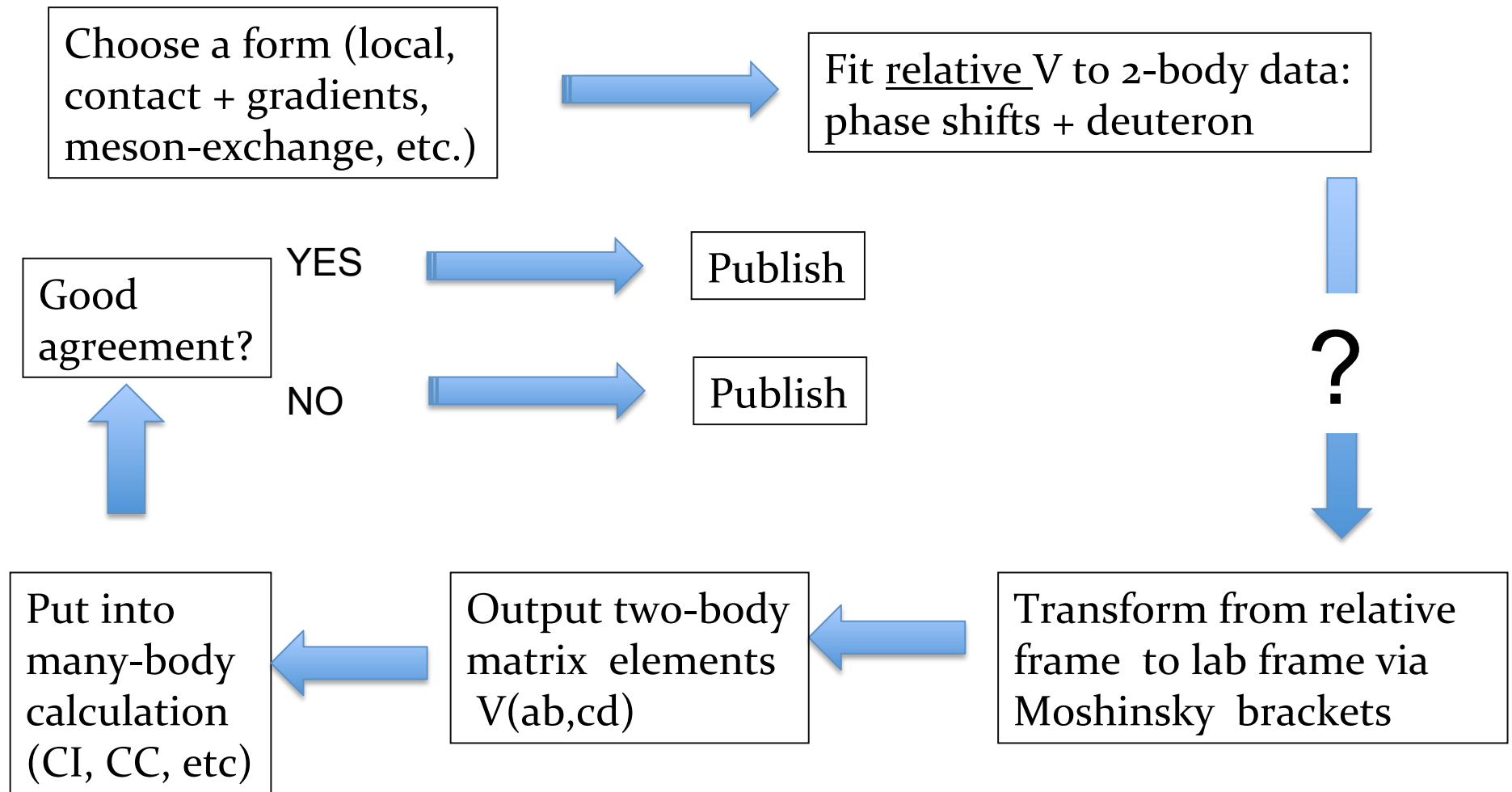
What about “realistic” effective nuclear interactions?

Q: What does it mean to be “realistic”?

A: Match experimental data!

Interlude: What about "realistic" effective nuclear interactions?

Life cycle of a realistic interaction:



Interlude: What about "realistic" effective nuclear interactions?

Life cycle of a realistic interaction:

Fit relative  $V$  to 2-body data:  
phase shifts + deuteron

Here is where one needs to  
"renormalize" the short-range/  
high momentum part of the  
interaction



**Today this renormalization is accomplished via unitary transformations that preserve two-body data (phase shifts, bound states)**

Transform from relative frame to lab frame via Moshinsky brackets

Interlude: What about "realistic" effective nuclear interactions?

Some common unitary transformations are Okubo-Lee-Suzuki,  $V_{\text{low-k}}$ , and the similarity renormalization group (SRG).

They all have the same goal: soften the short-range/high- $p$  behavior while preserving two-body (on-shell) data. **In other words, they modify the off-shell behavior, which can only be seen in many-body ( $A = 3$  and higher) systems.**

There have been some other attempts to choose different off-shell behavior, e.g., UCOM, INOY and JISP16 interactions.

Interlude: What about "realistic" effective nuclear interactions?

They all have the same goal: soften the short-range/high- $p$  behavior while preserving two-body (on-shell) data. **In other words, they modify the off-shell behavior, which can only be seen in many-body ( $A = 3$  and higher) systems.**

$$\hat{H}_{eff} = \hat{U}^{-1} \hat{H} \hat{U} = e^{-\hat{A}} \hat{H} e^{\hat{A}}$$



Can we choose the **best** generator  $A$  of the unitary transformation...  
the same way we fitted semi-empirical interactions?

Part 3: Cracking the off-shell degrees of freedom in in "realistic" interactions

A Modest Proposal:

$$\hat{H}_{eff} = \hat{U}^{-1} \hat{H} \hat{U} = e^{-\hat{A}} \hat{H} e^{\hat{A}}$$

We can expand the antisymmetric operator  $\mathbf{A}$  in a series of "base" operators:  $\hat{A} = \sum_i c_i \hat{A}_i$

Then we can find perturbations of the unitary transformation  $\hat{H}_{eff} \approx \hat{H} + \sum_i c_i [\hat{H}, \hat{A}_i]$



Then we compute

$$B_{\alpha i} = \frac{\partial E_{\alpha}}{\partial c_i} = \langle \alpha | [\hat{H}, \hat{A}_i] | \alpha \rangle$$

and do SVD as before...

Part 3: Cracking the off-shell degrees of freedom in "realistic" interactions

$$\hat{H}_{eff} = \hat{U}^{-1} \hat{H} \hat{U} = e^{-\hat{A}} \hat{H} e^{\hat{A}}$$

This is *just like* the SVD fits to semi-empirical interactions such as USDB, GXPF1, etc, except

USDB etc: work in lab frame, perturb Hamiltonian

New: we perturb the generators of the unitary transformation in the relative frame

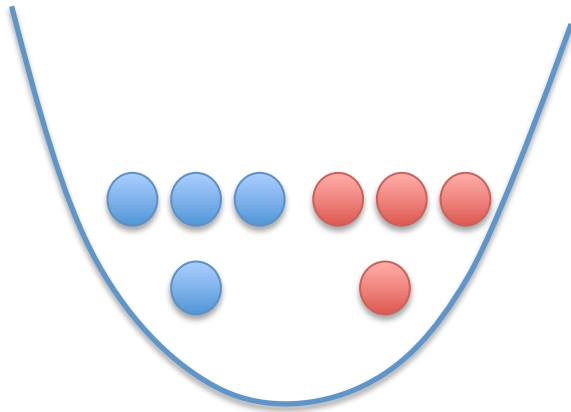


$$B_{\alpha i} = \frac{\partial E_{\alpha}}{\partial c_i} = \langle \alpha | [\hat{H}, \hat{A}_i] | \alpha \rangle$$

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Sample application: cold atomic gases at unitarity in a harmonic trap

$$\hat{H} = \sum_i -\frac{\hbar^2}{2m} \nabla_i^2 + \frac{1}{2} m \Omega^2 r_i^2 - V_0 \sum_{i < j} \delta(\vec{r}_i - \vec{r}_j)$$



$V_0$  tuned for infinite scattering length  
(cutoff-dependent)

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Sample application: cold atomic gases at unitarity in a harmonic trap

Only  $s$ -wave channel in relative coordinates 
$$\hat{H} = \sum_i -\frac{\hbar^2}{2m} \nabla_i^2 + \frac{1}{2} m \Omega^2 r_i^2 - V_0 \sum_{i < j} \delta(\vec{r}_i - \vec{r}_j)$$

Use ABF regularization

Alhassid, Bertsch, Fang, PRL100,

230401(2008)

in relative frame with  
harmonic oscillator basis  
up to  $N\hbar\Omega$

$$\langle n'l | \hat{H}_{rel} | nl \rangle$$

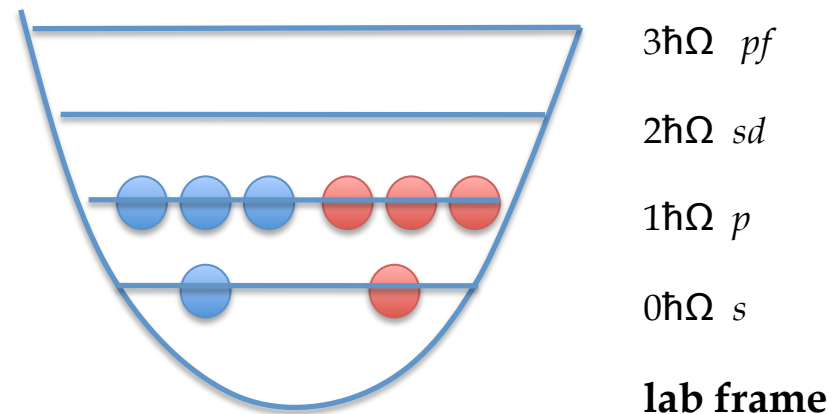
If cutoff at  $10\hbar\Omega$  ( $q=5$ ) then a  
6x6 symmetric matrix

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Sample application: cold atomic gases at unitarity in a harmonic trap

Then need to transform from the relative frame to the lab frame (also in harmonic oscillator basis) using Talmi-Brody-Moshinsky brackets

So there are *two* parameters for the system: the cutoff in the relative frame and the number of h.o. shells in the lab frame.

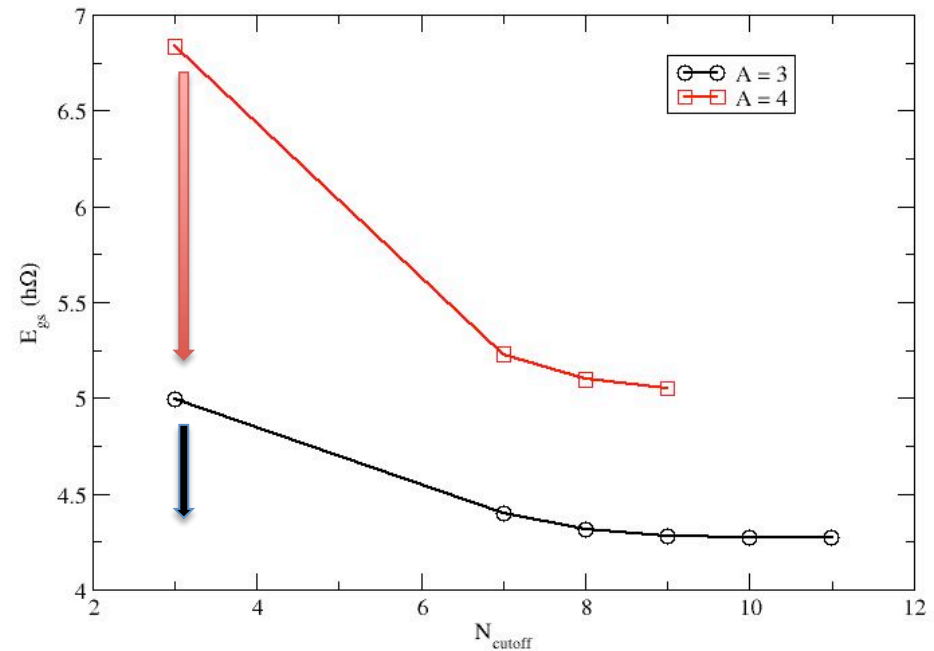


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Sample application: cold atomic gases at unitarity in a harmonic trap

Use ABF regularization  
with cutoff of  $10\hbar\Omega$   
(in relative  $s$ -channel).

Slow convergence  
in CI calculations.



Part 3: Cracking the off-shell degrees of freedom in "realistic" interactions

Sample application: cold atomic gases at unitarity in a harmonic trap

$$\hat{H}_{eff} = \hat{U}^{-1} \hat{H} \hat{U} = e^{-\hat{A}} \hat{H} e^{\hat{A}}$$

In the relative frame, with a  $8\hbar\Omega$ , then  $\mathbf{H}$ ,  $\mathbf{A}$ , and  $\mathbf{U}$  are all  $5 \times 5$  matrices. (Then go from relative to lab via Moshinsky).

There are thus 10 generators of  $\mathbf{A}$ .

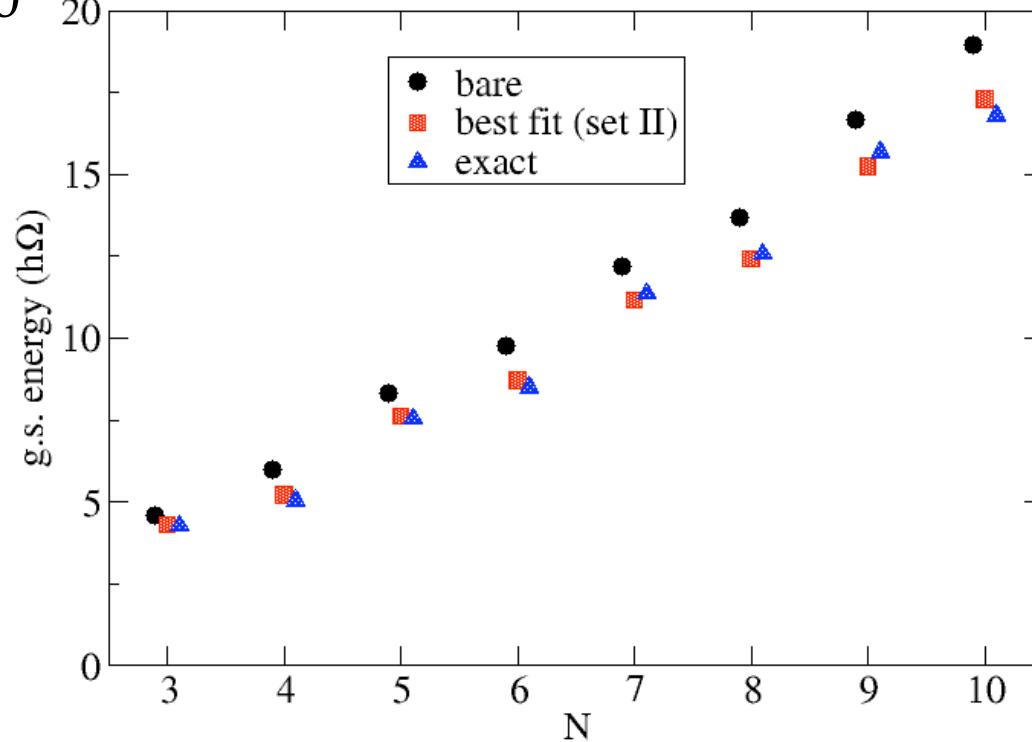
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Part 3: Cracking the off-shell degrees of freedom in "realistic" interactions

Sample application: cold atomic gases at unitarity in a harmonic trap

Using all generators, fit to g.s.  
energies for  $N = 3-10$

starting rms = 2.32  
final rms = 0.25



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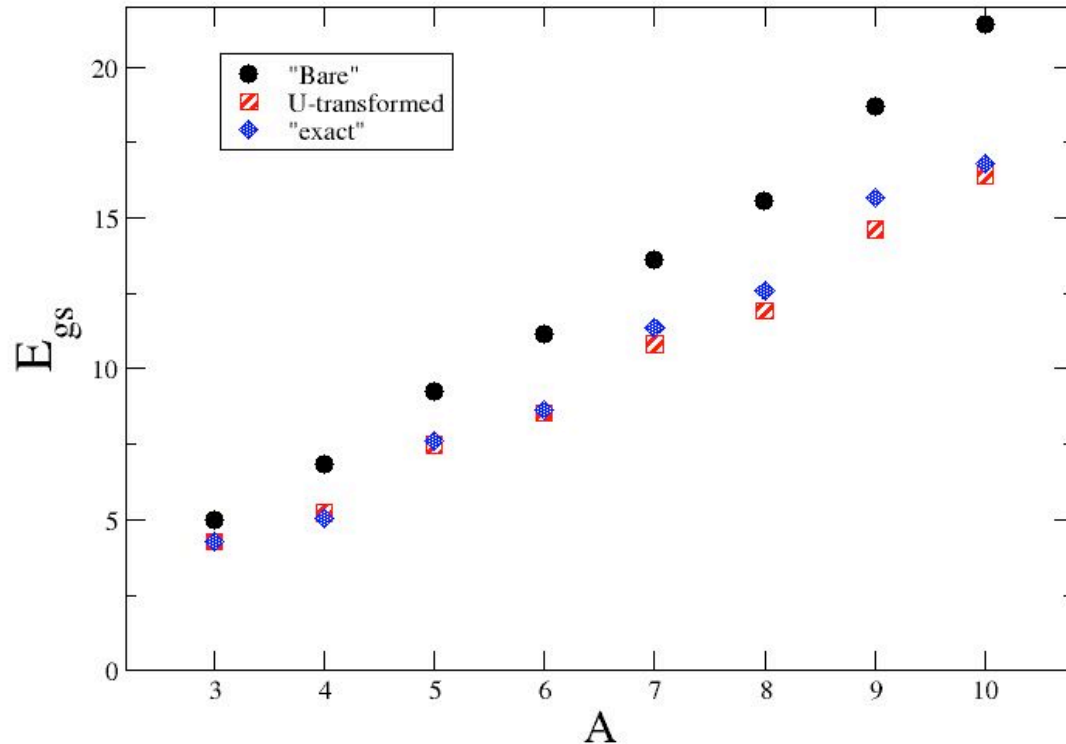
Part 3: Cracking the off-shell degrees of freedom in "realistic" interactions

Sample application: cold atomic gases at unitarity in a harmonic trap

Using only 1 generator ( $d/dr$ ) (very much like UCOM)

Fit to  $A = 3, 1^-, 0^+$   
 $A = 4, 0^+, 1^+, 2^+$

starting rms = 2.32  
final rms = 0.58



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I have developed a general formalism using unitary transformations that (a) preserve desired properties (on-shell matrix elements, eigenvalues) and (b) can be fitted to data.

Preliminary application to a cold atomic gas at unitarity is promising.



Next step: apply to nuclear systems (more complicated, multi-channel; not only binding energies, but also spin-orbit splitting usually attributed to 3-body forces)

## Center of mass without exact factorization

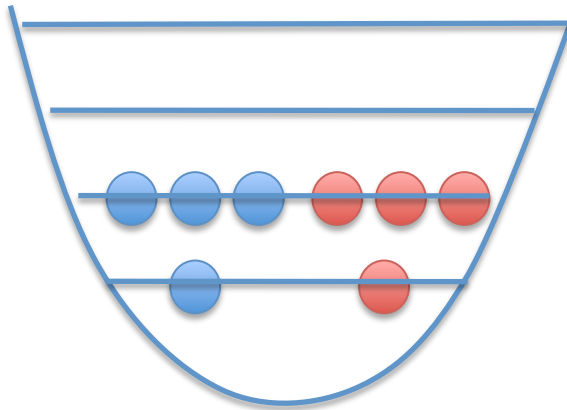
Center-of-mass is an important contamination in nuclear structure calculations.

A theorem (Palumbo, later Lawson) showed that in a h.o. basis, a specific truncation (the  $Nh\Omega$  truncation) guarantees a system with a translationally invariant interaction can decouple relative from c.m. motion.

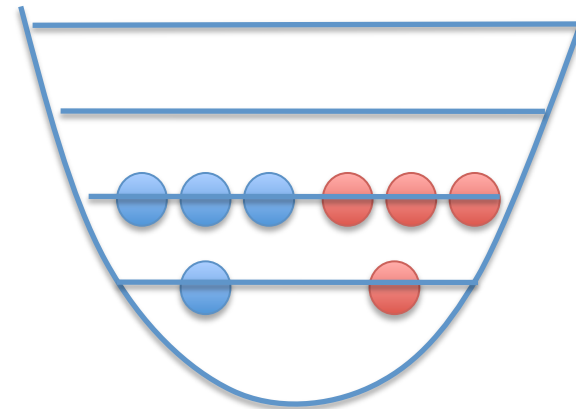
However a more “natural” truncation is by maximal orbits: this is natural in Hartree-Fock, coupled-cluster, etc.

## A tale of two truncations

orbit truncation: all  
excitations



$Nh\Omega$  (or energy) truncation: only  
those excitations in noninteracting  
h.o. with energy  $\leq Nh\Omega$



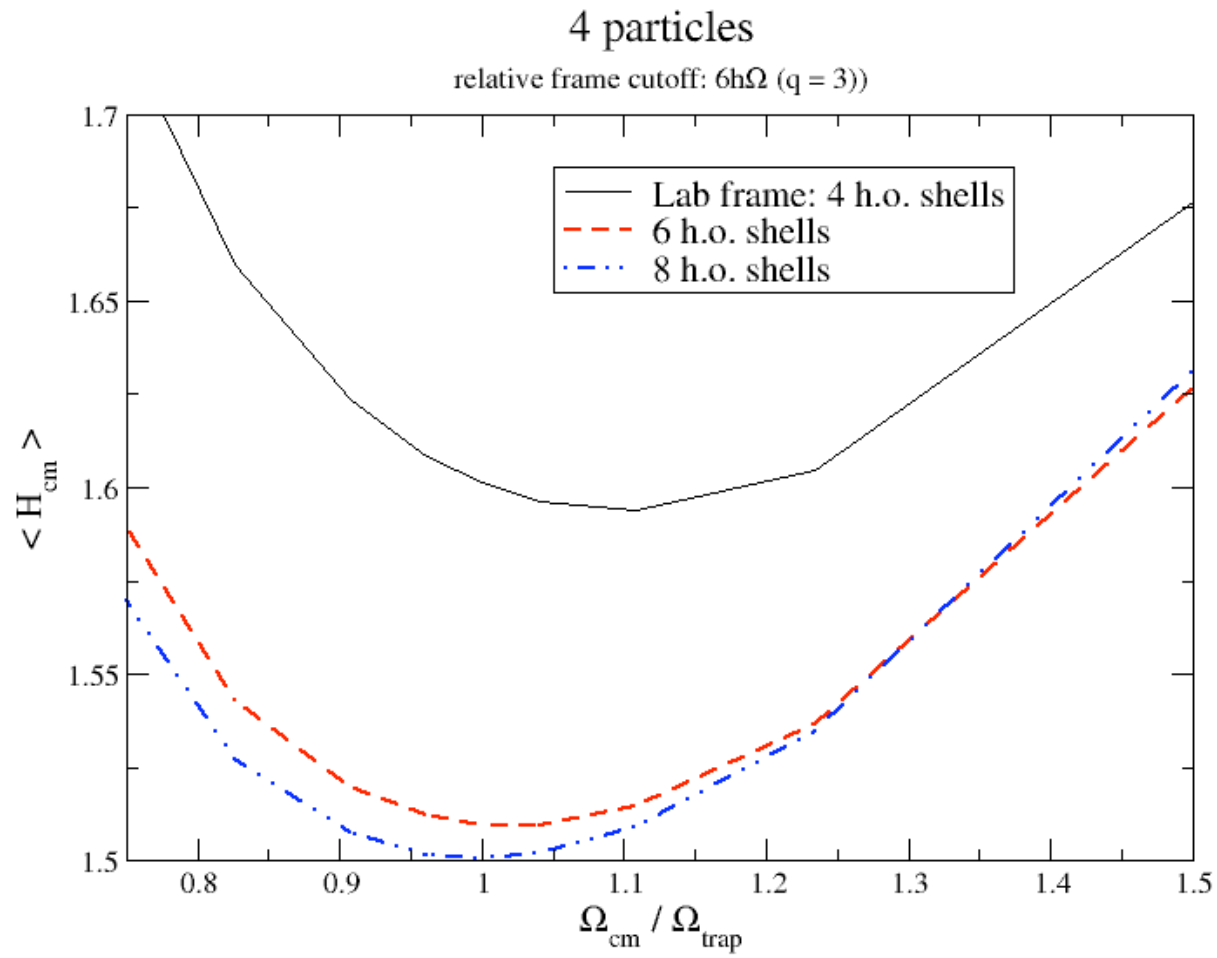
## But is the orbit truncation bad?

Hagen, Papenbrock, and Dean: in CC, look at  $\langle H_{\text{cm}} \rangle$

$H_{\text{cm}}$  is minimized, only with h.o. frequency different from the basis

Roth, Gour, and Piecuch: in importance-truncated CI, look also at perturbations by adding  $\beta H_{\text{cm}}$ ; considerable contamination in orbital truncation

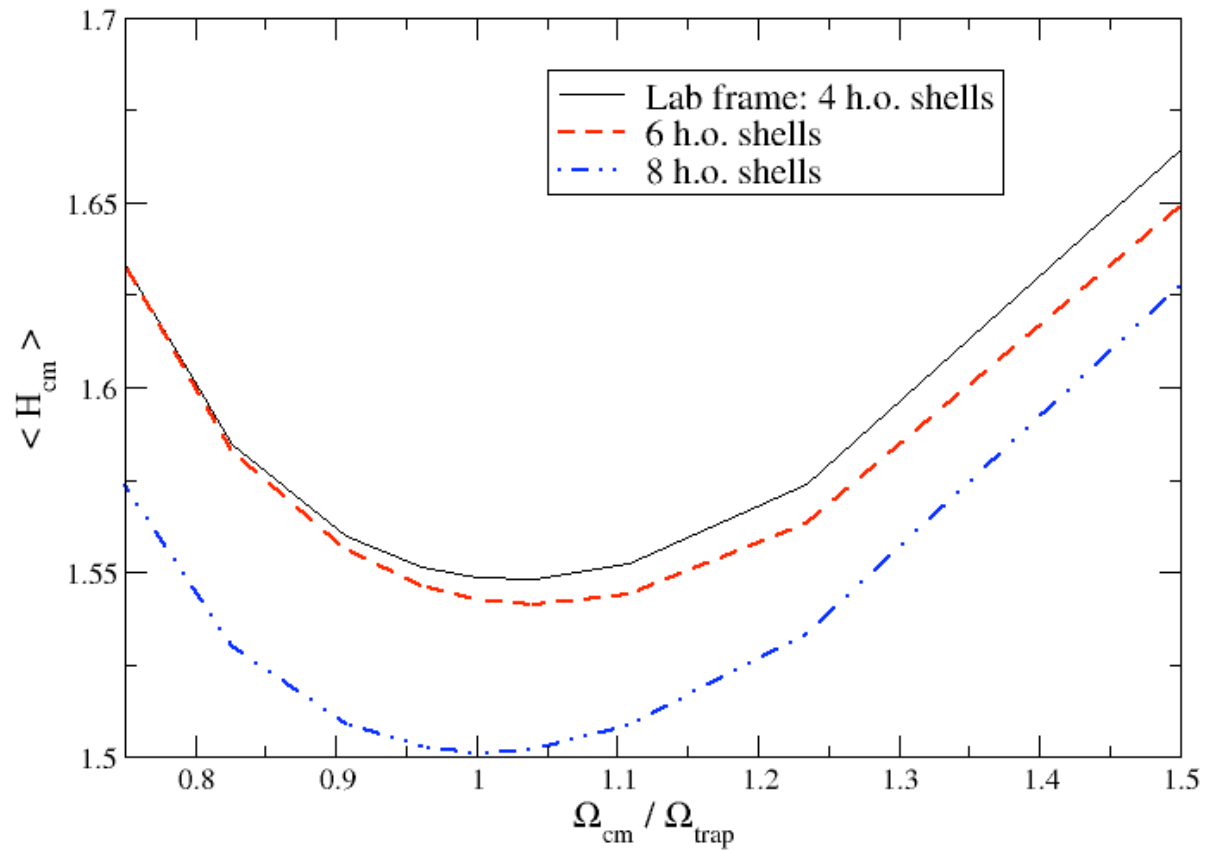
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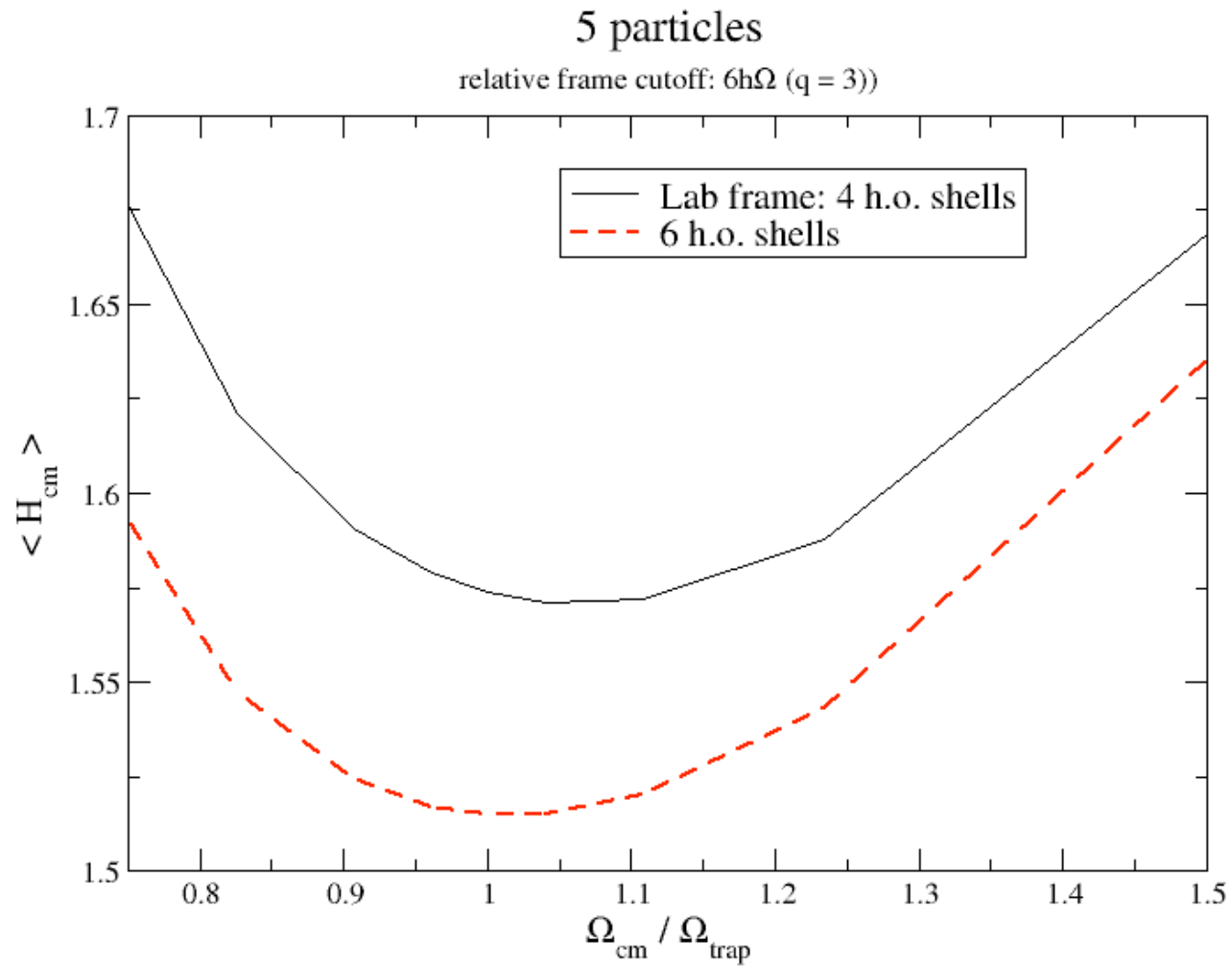
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4 particles -- 1st excited state

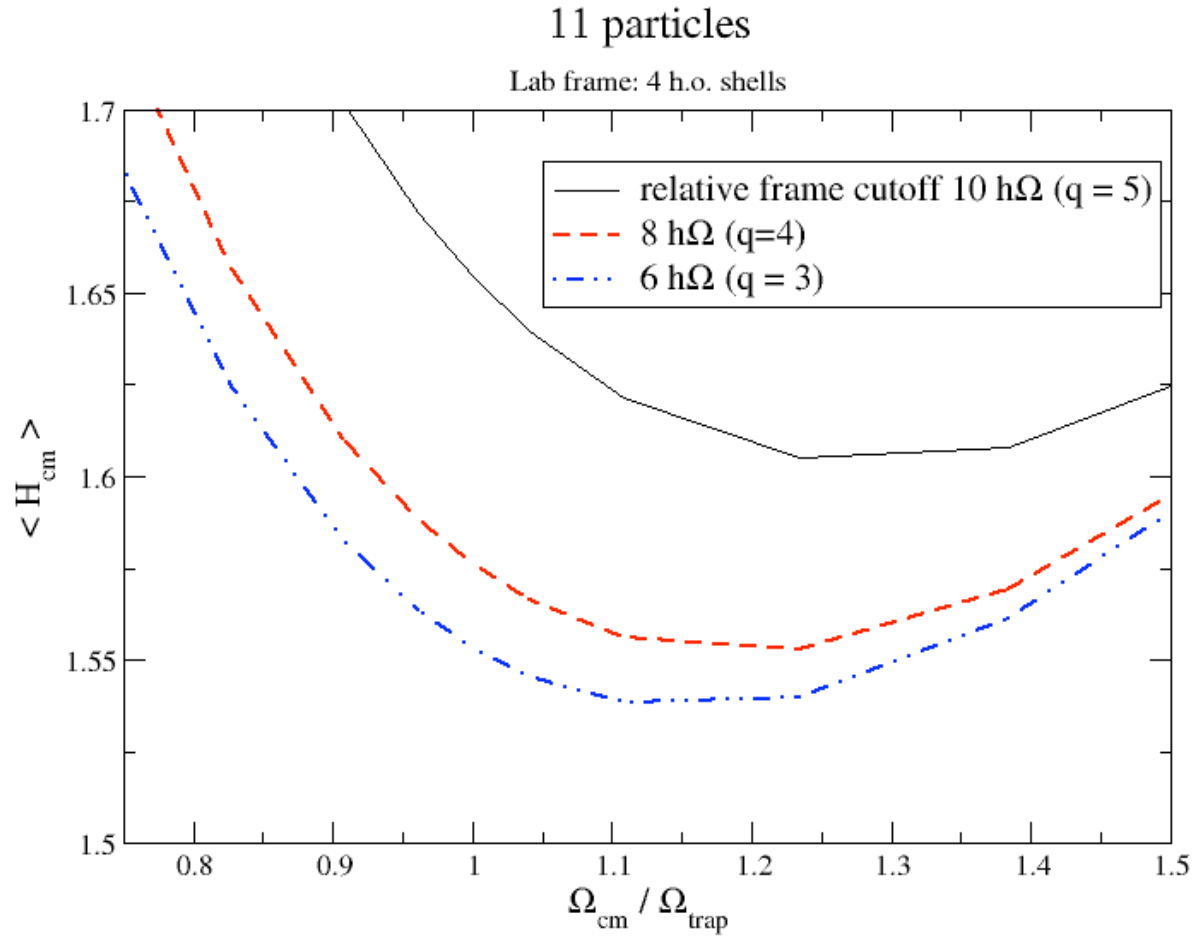
relative frame cutoff:  $6\hbar\Omega$  ( $q = 3$ )



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Need to look at sensitivity to adding  $\beta H_{cm}$ ;

Work is under way for nuclei