

7/11/2013

Th2a-1

Plan for Renormalization Group 2:

- ① Recap of some important points from the slides shown yesterday (using W2a-2 and 2b plus the exercises)
- ② Local projection (visualizing potential changes in coordinate space) and flow to universal potentials (where does unresolved physics go?)
- ③ Alternative generators - Wegner and block diagonal (N_{low})
- ④ Operators and many-body contributions
- ⑤ Perturbativeness - Weinberg eigenvalues
- ⑥ Computational aspects

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- Recap of important points from the slides

• Nuclei would be at low resolution based on Fermi momentum in large nucleus,

• recall from exercises $\rho = \frac{2}{3\pi^2} k_F^3$ (for protons and neutrons)
 $\Rightarrow k_F = (3\pi^2 \rho)^{1/3}$ (see next page)

and density of heavy nuclei about constant $\Rightarrow k_F \approx 1.1 - 1.3 \text{ fm}^{-1}$

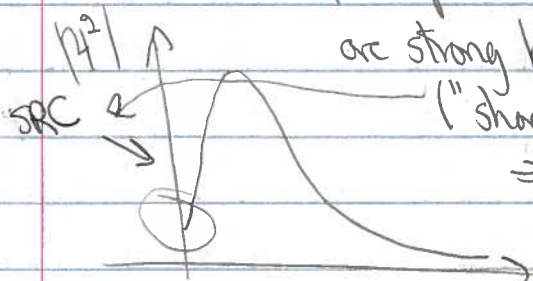
• So typical relative momentum of $\approx 1 \text{ fm}^{-1}$ ($\approx 200 \text{ MeV}$) in a large nucleus. Even less in light nuclei

• But if the potential has a repulsive core, then there

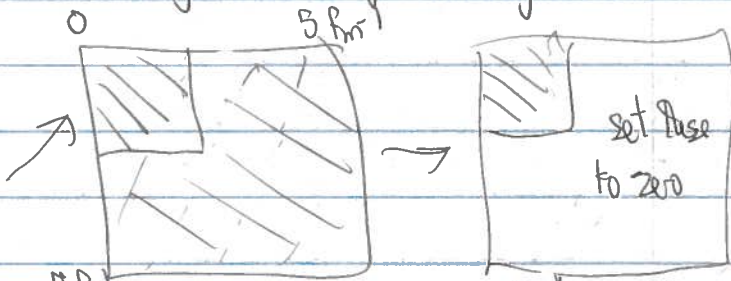
are strong high momentum components ("short-range correlations")

\Rightarrow • slow-down convergence of many-body nuclei.

• eg. matrices get too big,



• low pass filter fails even for low energy.



Why? Because of

quantum mechanics,

$$T = V + V \frac{1}{E - H_0} V + \dots$$

If strong off-diagonal comp \Rightarrow can't ignore.

$$\Rightarrow \langle k | T | k \rangle = \langle k | V | k \rangle + \frac{2}{\pi} \int_0^{k_F} dk' \frac{\langle k | V | k' \rangle \langle k' | V | k \rangle}{(k^2 - k'^2)/m} + \dots$$

\uparrow low momentum

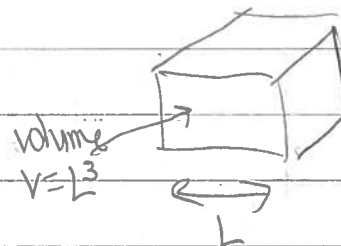
Solution? Unitary transformation to decouple! Use RG to do it,

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WDA-2b

[Aside: deriving $g = \frac{2}{3\pi^2} k_F^3$...

• For a non-interacting Fermi gas, imagine putting N particles in a box of side $L \Rightarrow g = \frac{N}{V}$



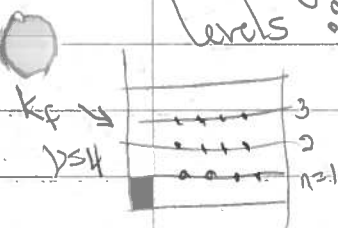
• when in doubt for uniform system, put in box and take $V \rightarrow \infty$ limit at end

• Let ν be the spin-isospin degeneracy:

$\nu = 2$ for neutrons only (spin up, spin down)

$\nu = 4$ for symmetry matter ($\uparrow, \downarrow, n, p$)

• Apply periodic boundary conditions \Rightarrow discrete momentum levels:



Then $N = \sum_{k \leq k_F} \nu \cdot 1$

But pbc: $e^{ik(x+L)} = e^{ikx}$ in each dimension

$\Rightarrow k_n L = 2\pi n$ $n=1, 2, 3, \dots$ are allowed

$\Rightarrow n = \frac{kL}{2\pi}$ or $\Delta n = 1 = \frac{L}{2\pi} \Delta k$ in each dimension

in large V limit $\Rightarrow \sum_n \rightarrow \left(\frac{L}{2\pi}\right)^3 \int d^3k = \frac{V}{(2\pi)^3} \int d^3k$

$\Rightarrow N = \frac{V}{(2\pi)^3} \int_0^{k_F} d^3k \nu =$
 \leftarrow volume of sphere in large V limit

or $\frac{N}{V} = g = \frac{1}{8\pi^3} \cdot \frac{4}{3}\pi k_F^3 \cdot \nu = \frac{\nu k_F^3}{6\pi^2}$

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Wednesday exercises review

1(c) Why would we want to repeat nuclear structure calculations for different values of the SRG λ (or s)?

do 2, 3,
first
then return

- Observables are supposed to be unchanged
 \Rightarrow test if a quantity is an observable (example, clear demonstration that D-state probability in the deuteron is not)
- \Rightarrow determine the scale dependence of a quantity
- Test for errors
- Test approximations
 - We will see this particularly in considering many-body potentials and other operators,

2. General equation is $\frac{dH_s}{ds} = [\eta(s), H_s] = [[G_s, H_s], H_s]$

- T_{rel} (or T) doesn't change by construction
- What if we used $T = T_{com} + T_{rel}$ for G_s instead of T_{rel} ?

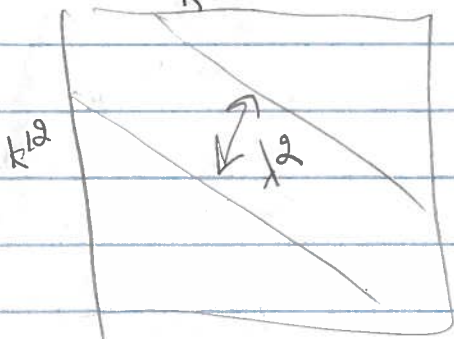
(answer: $[T_{com}, H_s] = 0$, so no difference!)

• other choices for G_s ?

3. $\langle k | \frac{dV_s}{ds} | k' \rangle \equiv \frac{dV_s(k, k')}{ds} = -\frac{1}{(k-k')^2} V_s(k, k') + \frac{2}{\pi} \int_0^{\infty} q^2 dq \frac{(k^2 - k'^2 - q^2)}{(k^2 - q^2)(k'^2 - q^2)} V_s(k, q) V_s(q, k')$

• If $-\frac{1}{(k-k')^2} V_s(k, k')$ dominates then $V_s(k, k') \approx V_{so}(k, k')$

$k \neq k'$
(must be sufficiently off diagonal)



• Look at slides

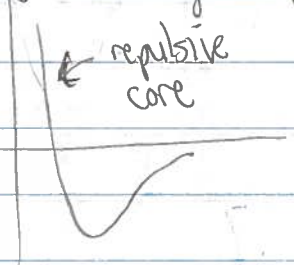
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Visualizing the softening of NN interactions

- In momentum space we associate softening of a potential with the decreased coupling between high and low momentum:

$$\langle k_{high} | V | k_{low} \rangle \rightarrow 0$$

But what does this do to our picture of potentials having strong short-range repulsion?



Visualizing is not so easy, because the potential becomes non-local, so it is a functional of r and r'

- note that the $-(k^2 - E^2)^{-2} V_0(k, k')$ term in the SRG equation (not partial wave projected) can be written using

$$k^2 - E^2 = (k+k')(k-k') \rightarrow \vec{p} \vec{q}$$

as an explicit function of total momentum $\vec{p} = (k+k')$ and not just momentum transfer $\vec{q} = k-k' \Rightarrow$ nonlocal

Why do partial waves not mix? (except for already coupled \Rightarrow insert expansion, only k^2, k'^2, p^2 dependence from SRG)

Plan: use a local projection

- The high-momentum tails of low-energy wavefunctions are suppressed by RG evolution which implies the wavefunction variation over short distances is small. So in the non-local Schrödinger equation:

$$-\frac{1}{2\mu} \nabla^2 \psi(\vec{r}) + \int d^3r' V(\vec{r}, \vec{r}') \psi(\vec{r}') = E \psi(\vec{r})$$

← treat as constant over range of V nonlocality

$$\rightarrow -\frac{1}{2\mu} \nabla^2 \psi(\vec{r}) + \psi(\vec{r}) \int d^3r' V(\vec{r}, \vec{r}') \approx E \psi(\vec{r})$$

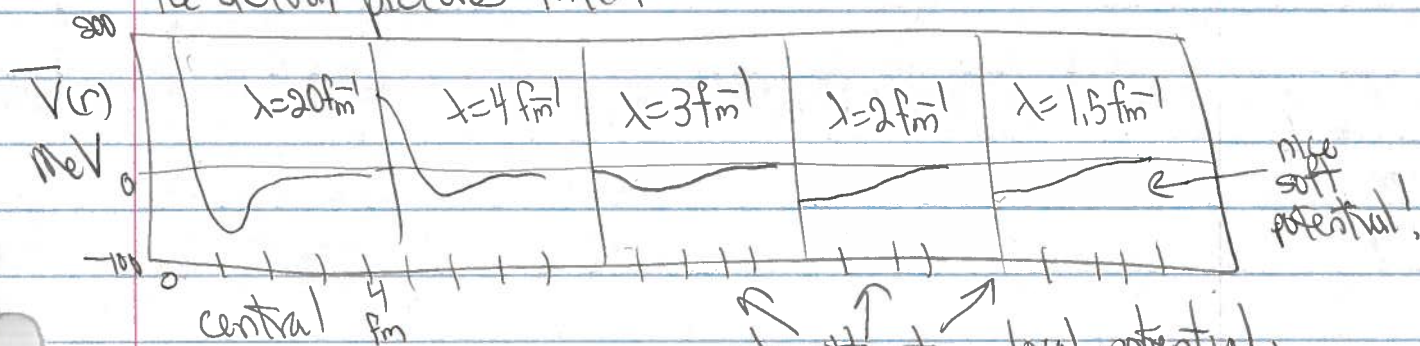
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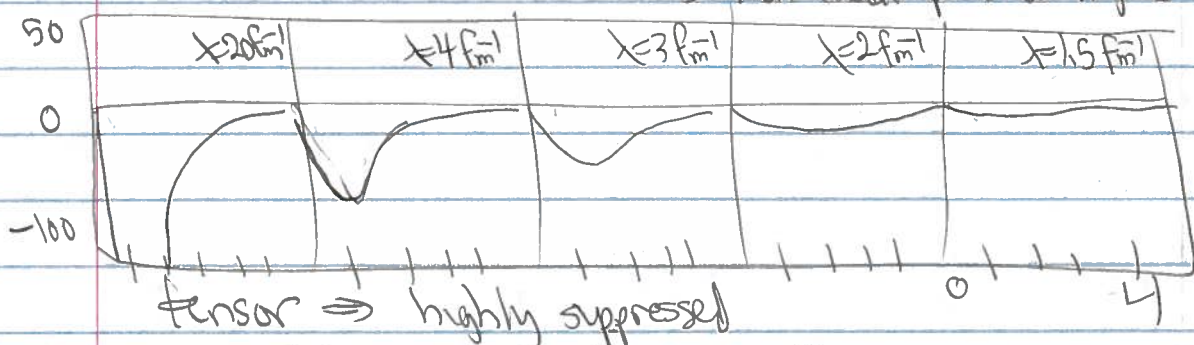
⇒ define $\bar{V}_\lambda(r) = \int d^3r' V_\lambda(\vec{r}, \vec{r}')$ as the local projection,

• Kyle Wendt has developed this idea further, to apply beyond S-waves (which is all that survives the angular integral).

• We'll sketch the result for the AV18 potential and look at the actual pictures later.



purely attractive local potential,
so phase shift must fail to change sign
⇒ non-local part at higher momentum



- D-state probability changes greatly
- but asymptotic D-S ratio unchanged!
- what about quadrupole moment?

• Different potentials evolve to same in both momentum rep. (at momentum below λ and in local projection),

• Where do you expect high energy contributions to go? c.f. $\lambda_0 \rightarrow \lambda_0 + \lambda_0$ $\rightarrow \lambda_0 - \lambda_0 + \lambda_0$
⇒ see slides (same thing here!)

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- Wegner choice for the flow-equation generator
 \Rightarrow use the diagonal of H_S in whatever basis you are in. E.g., $H_d = T_{rel} + V_S(k, k)$
 \nwarrow diagonal

- Let's consider the general case with $H_{ii} \equiv e_i$ where we are labeling the basis elements $|i\rangle, |j\rangle, \dots$
 - Note that these could be plane waves, harmonic oscillators, 2 particle or more, ...

$$\langle i | \frac{dH_S}{ds} | j \rangle = \frac{dH_{ij}}{ds} = \langle i | [[H_d, H_S], H_S] | j \rangle$$

$$= \langle i | H_d H_S H_S - H_S H_d H_S - H_S H_d H_S + H_S H_S H_d | j \rangle$$

\nwarrow \nwarrow \nwarrow \nwarrow
 insert $\sum_k |k\rangle \langle k|$ and we $H_d |j\rangle = e_j |j\rangle$, etc.

$\Rightarrow \frac{dH_{ij}}{ds} = \sum_k (e_i - e_k - e_k + e_j) H_{ik} H_{kj}$ ← simple matrix multiplication

$i=j \Rightarrow \frac{dH_{ii}}{ds} = 2 \sum_k (e_i - e_k) |H_{ik}|^2$

We want to ask: what can we say about $\frac{d}{ds} \sum_{i \neq j} |H_{ij}|^2$?

• This is the sum of the squares of the off-diagonal parts. Does it decrease?

• The full sum is $\sum_{ij} |H_{ij}|^2 = \sum_{ij} H_{ij} H_{ji} = \text{Tr } H_S^2 = \text{constant!}$ The trace is invariant.

So $\frac{d}{ds} \sum_{i \neq j} |H_{ij}|^2 = - \frac{d}{ds} \sum_i |H_{ii}|^2 = -2 \sum_i \underbrace{H_{ii}}_{e_i} \underbrace{\left(\frac{dH_{ii}}{ds} \right)}_{\text{from eqn. 1}} = -4 \sum_{i \neq k} e_i (e_i - e_k)$

\Rightarrow except for degeneracies, off-diagonal elements decrease. $= -2 \sum_{i \neq k} (e_i - e_k)^2 |H_{ik}|^2 \leq 0!$

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Th 2a'6

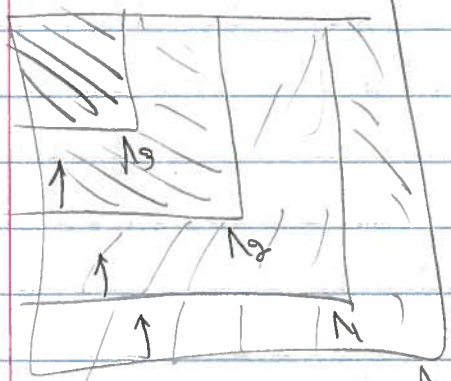
• The use of T_{rel} instead of $H_0 = T_{rel} + H_{ij}$ is ok for nuclear physics, at least in the momentum basis, because $T_{rel} \gg (V_s)_{ij}$ so $H_0 \approx T_{rel}$

25
Schulz 2

- It can fail, though, B \Rightarrow see Wendt et al, with large Λ leading order forces,
 \Rightarrow good example of decoupling.

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• $V_{low k}$ RG equation - Bogner, Kuo, Schwenk (2001)



• based on requiring the half-off shell T matrix to be invariant with a change in cutoff on the sum over intermediate states,

$$\left[\frac{d}{d\Lambda} \left(\begin{array}{c} \text{Diagram 1} \\ \text{Diagram 2} \\ \text{Diagram 3} \end{array} \right) \right] = 0$$

initial $T(k', k; k^2) = V_{low k}^{(k', k)} + \frac{2}{\pi} \mathcal{P} \int_0^{\Lambda_0} \frac{V_{low k}(k', p) T(p, k; k^2)}{k^2 - p^2} p^2 dp$

for all $k, k' < \Lambda$ } with cutoff $= V_{low k}^\Lambda(k', k) + \frac{2}{\pi} \mathcal{P} \int_0^\Lambda \frac{V_{low k}^\Lambda(k', p) T(p, k; k^2)}{k^2 - p^2} p^2 dp$

principal value so real

half-on-shell because k, k^2 but $p \neq k$

Take $\frac{dT}{d\Lambda} = 0 \Rightarrow \frac{d}{d\Lambda} V_{low k}^\Lambda(k', k) = \frac{2}{\pi} \frac{V_{low k}^\Lambda(k, \Lambda) T(\Lambda, k; \Lambda^2)}{1 - (k/\Lambda)^2}$

derivation is not immediate, (see Bogner et al.)

cf, partial wave SRG equation (with $G_S = T_{rel}$)

$$\frac{d}{d\lambda} V_\lambda(k, k') = \left(-\frac{4}{\lambda^5} \right) (k^2 - k'^2)^2 V_\lambda(k, k') + \frac{2}{\pi} \int_0^\infty \frac{(k^2 + k'^2 - 2q^2)}{(k^2 - q^2)(k'^2 - q^2)} V_\lambda(k, q) V_\lambda(q, k') q^2 dq$$

from $\frac{d}{d\lambda} = \frac{d}{d\Lambda} \left(\frac{d\Lambda}{d\lambda} \right)$
and $s = \lambda^4$

• Compare rhs: T matrix for $V_{low k}^\Lambda$ but just potential for SRG

\Rightarrow SRG much easier for $A > 2$ (otherwise need T matrix in all channels).

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Can we get a $V_{low k}$ -like potential from the SRG flow equation by an appropriate choice of G_s ? Yes!

Use $\frac{dH_s}{ds} = [[G_s, H_s], H_s]$ with $G_s = \begin{pmatrix} PH_sP & 0 \\ 0 & QH_sQ \end{pmatrix}$

Choose Λ and then G_s is the running Hamiltonian with the off-diagonal blocks defined by Λ set equal to zero.

P and Q are projection operators, $P+Q=1$.

Proof that this does what we want [Gubankova et al.]

A measure of off-diagonal coupling is QH_sP
 so this is the part that does the coupling, \uparrow upper \leftarrow lower

$\Rightarrow \sum_n \langle \gamma_n | (QH_sP)^\dagger (QH_sP) | \gamma_n \rangle = \text{Tr} [PH_sQH_sP] \geq 0$ (since $Q^2=Q, P^2=P$)
 $[G_s, H_s] = [PH_sP + QH_sQ, H_s]$

Now check how this changes with s using $\frac{d}{ds} H_s = [M_s, H_s]$

$\frac{d}{ds} \text{Tr} [PH_sQH_sP] = \text{Tr} [PM_sQ(QH_sQH_sP - QH_sPH_sP)]$
 $+ \text{Tr} [(PH_sPH_sQ - PH_sQH_sQ)Q M_s P]$

you are invited to prove it, $\Rightarrow = -2 \text{Tr} [(QM_sP)^\dagger (QM_sP)] \leq 0$

\Rightarrow The off-diagonal QH_sP block will decrease (or not increase) as s increases.

Two examples: $G_s = \text{Tr} \begin{pmatrix} \diagdown & \\ & \diagup \end{pmatrix} \Rightarrow$ goes to $\begin{pmatrix} \diagdown & \\ & \diagup \end{pmatrix}$

$G_s = \begin{pmatrix} \text{||||} & \\ & \text{||||} \end{pmatrix} \Rightarrow$ goes to $\begin{pmatrix} \diagdown & \\ & \text{||||} \end{pmatrix}$

Does it always evolve to the pattern of G_s ? See pictures!

7/11/2003

Thda-9

SRG Flow of Operators

- We'll have more to say about operators in a future lecture; just some basic ideas here.

- When we transform $H_s = U(s)H_{s_0}U^\dagger(s)$, the wave functions also get transformed: $| \psi_n^s \rangle = U(s) | \psi_n^{s_0} \rangle$, so that energies are unchanged $E_n = \langle \psi_n^{s_0} | H_{s_0} | \psi_n^{s_0} \rangle = \langle \psi_n^s | H_s | \psi_n^s \rangle$

- So this means that any operator O must be transformed:

$$O_s = U(s)O U^\dagger(s)$$

- We can calculate this directly by constructing $U(s)$ and $U^\dagger(s)$. We do this by first evolving H to H_s , then finding all the eigenstates $| \psi_n^s \rangle$ of H and $| \psi_n^{s_0} \rangle$ of H_{s_0} .

- Then we have $U(s) = \sum_n | \psi_n^s \rangle \langle \psi_n^{s_0} |$

- in a basis like momentum space, this would give us the matrix element $\langle k | U(s) | k' \rangle = \sum_n \langle k | \psi_n^s \rangle \langle \psi_n^{s_0} | k' \rangle$

- just an outer product,

- This works fine in practice but there are two other ways

i) Evolve O_s with its own flow equation

ii) evolve $U(s)$ " " " " " and then use $O_s = U(s)O U^\dagger(s)$

What are the equations? \Rightarrow you do that for exercises!

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Ph20-10

How do we know that SRG evolution of operators (including the Hamiltonian) must generate many-body terms?

• In exercises: fairy about physics, here: formal discussion.

• Consider 2nd quantization. This is defined with two ingredients:

① a single-particle basis (eg. plane waves in a box or HO wfs)

② a reference state that serves as the "vacuum".

examples

- could be the actual vacuum

- or a filled core (Fermi sea or a closed shell)

• Kinetic energy: $T = \sum_i \frac{p_i^2}{2m} a_i^\dagger a_i$

• Two-body potential: $\frac{1}{4} \sum_{ijkl} V_{ijkl}^{(2)} a_i^\dagger a_j^\dagger a_l a_k$

• 3-body potential: $\frac{1}{36} \sum_{ijklmn} V_{ijklmn}^{(3)} a_i^\dagger a_j^\dagger a_k^\dagger a_n a_m a_l$

• These operators have anti-commutation relations:

$$\{a_i, a_j^\dagger\} = a_i a_j^\dagger + a_j^\dagger a_i = \delta_{ij}, \quad \{a_i, a_j\} = \{a_i^\dagger, a_j^\dagger\} = 0$$

Claim, $\frac{dV_S}{ds} = \left[\underbrace{\sum_{Q \neq T} a^\dagger a}_{1\text{-body}}, \underbrace{\sum a^\dagger a^\dagger a a}_{2\text{-body}}, \underbrace{\sum a^\dagger a^\dagger a a}_{2\text{-body}} \right] = \dots + \underbrace{\sum a^\dagger a^\dagger a a}_{3\text{-body}}$

• And this is just one time step!

• \Rightarrow A-body operators generated

• Is this a problem?

• depends: we need to be able to truncate \Rightarrow need hierarchy

• also need to be able to calculate with minimal (usually 3-body)

In-medium

SRG \Rightarrow

• Alternative: Pick a different reference state \Rightarrow reshuffles what is many-body!

show this in the exercises

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SRG technology is to evolve 3-body forces

Three methods not exist:

slides by Angelo Calci

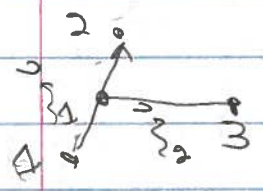
- i) evolve in a discrete harmonic oscillator basis
 - ⇒ applied to No-Core Shell Model (tomorrow)
- ii) evolve in a partial-wave momentum basis
 - ⇒ separate evolution of 2 and 3 body parts
 - ⇒ applied to neutron matter (next week)
- iii) evolve in a hyperspherical basis
 - ⇒ good features, visualization

Eric Jurgenson
many developments in Bonnstadt, P. Navrotili
Kai Hebel
instructor next week!

more later on these comparisons

Recent: 4-body evolution (see Angelo Calci talk from Trento)

Oscillator evolution is in a 3-body Jacobi basis



- generalization of center-of-mass and relative
 - $\vec{r}_0 = \frac{1}{\sqrt{3}} [\vec{r}_1 + \vec{r}_2 + \vec{r}_3]$ ← potential doesn't depend
 - $\vec{r}_1 = \sqrt{\frac{2}{3}} [\vec{r}_1 - \vec{r}_2]$ ← relative between 1 and 2.
 - $\vec{r}_2 = \sqrt{\frac{2}{3}} [\frac{1}{2}(\vec{r}_1 + \vec{r}_2) - \vec{r}_3]$ ← relative between 3 and com of 1 and 2

hard part!

must antisymmetrize HO basis: $|\alpha\rangle = |(N_1 L_1 S_1) S_1 (N_2 L_2 S_2) S_2 (S M_3) (T_1 T_2) T M_T\rangle$

momentum space evolution $|p q \alpha\rangle_i \equiv |p_i q_i; [(LS) S (L S)_i] | S_2 (T_1) T M_T\rangle$

