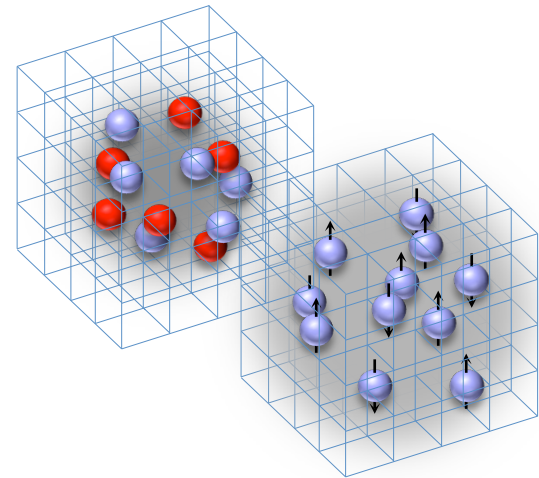


# New Developments in Lattice Effective Field Theory

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Intersections of Particle and Nuclear Physics  
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# Outline

Lattice effective field theory

Brief review of methods

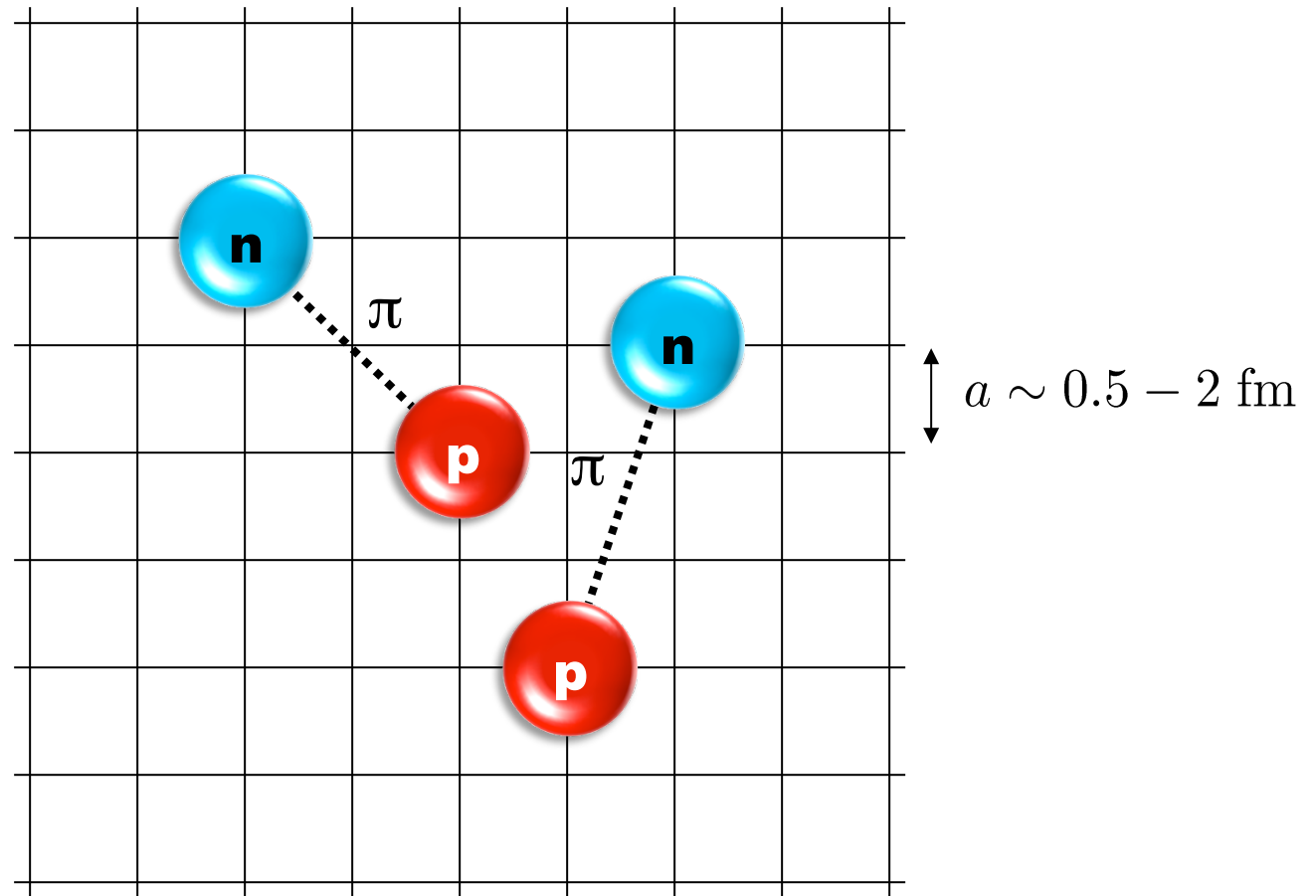
Thermodynamics and structure

Eigenvector continuation

Superfluid pair correlations

Summary and outlook

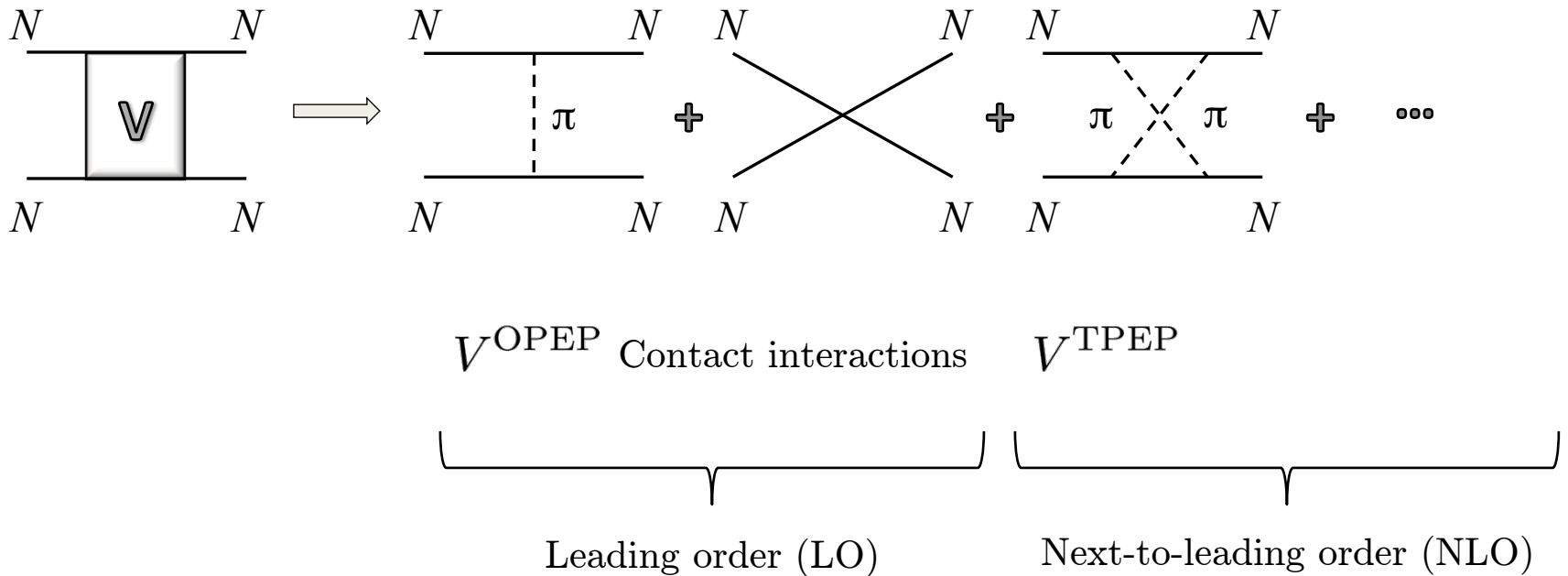
## Lattice effective field theory



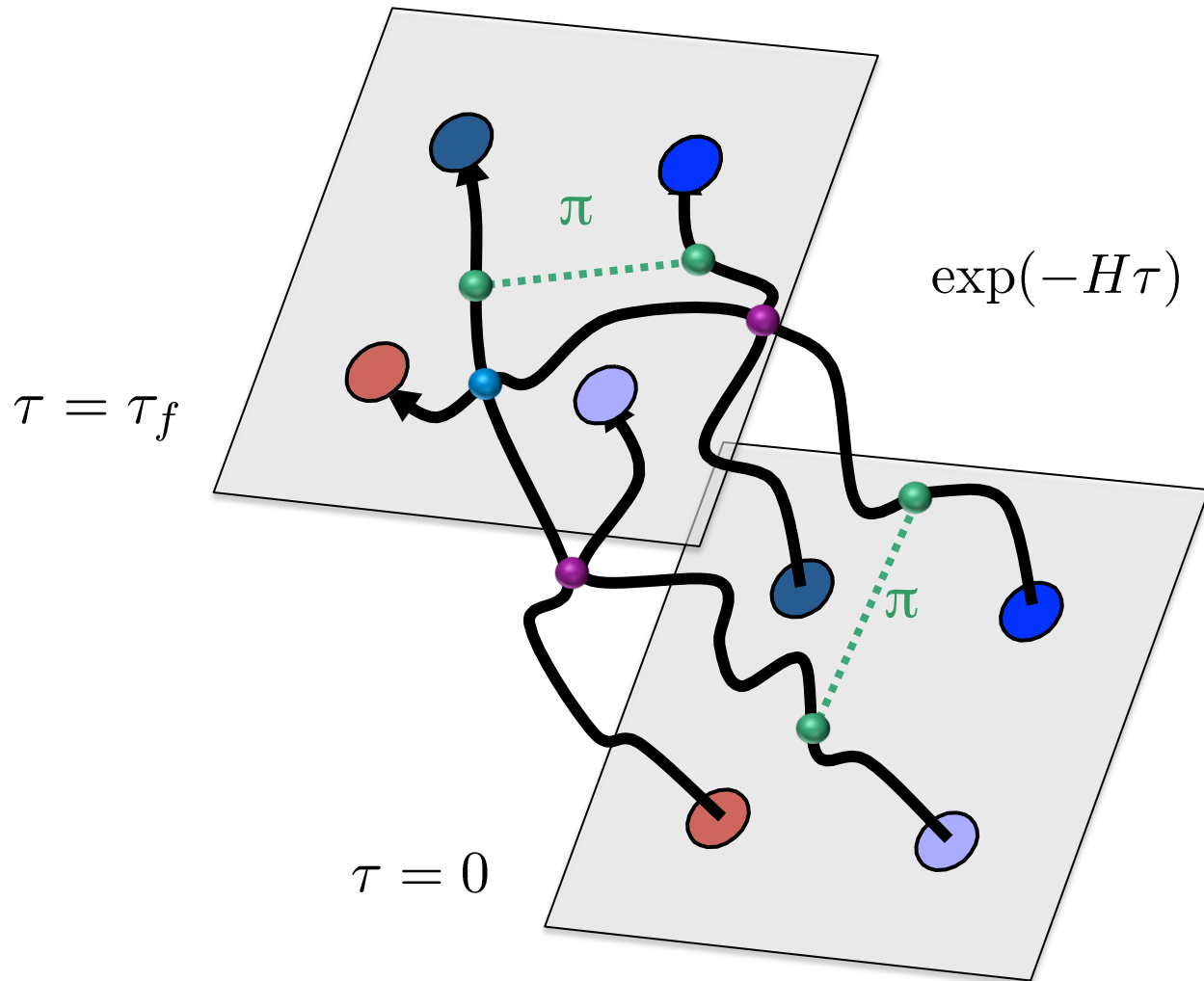
Review: D.L, Prog. Part. Nucl. Phys. 63 117-154 (2009)  
TALENT summer school lectures: [qmc2016.wordpress.ncsu.edu](http://qmc2016.wordpress.ncsu.edu)

# Chiral effective field theory

Construct the effective potential order by order



# Euclidean time projection

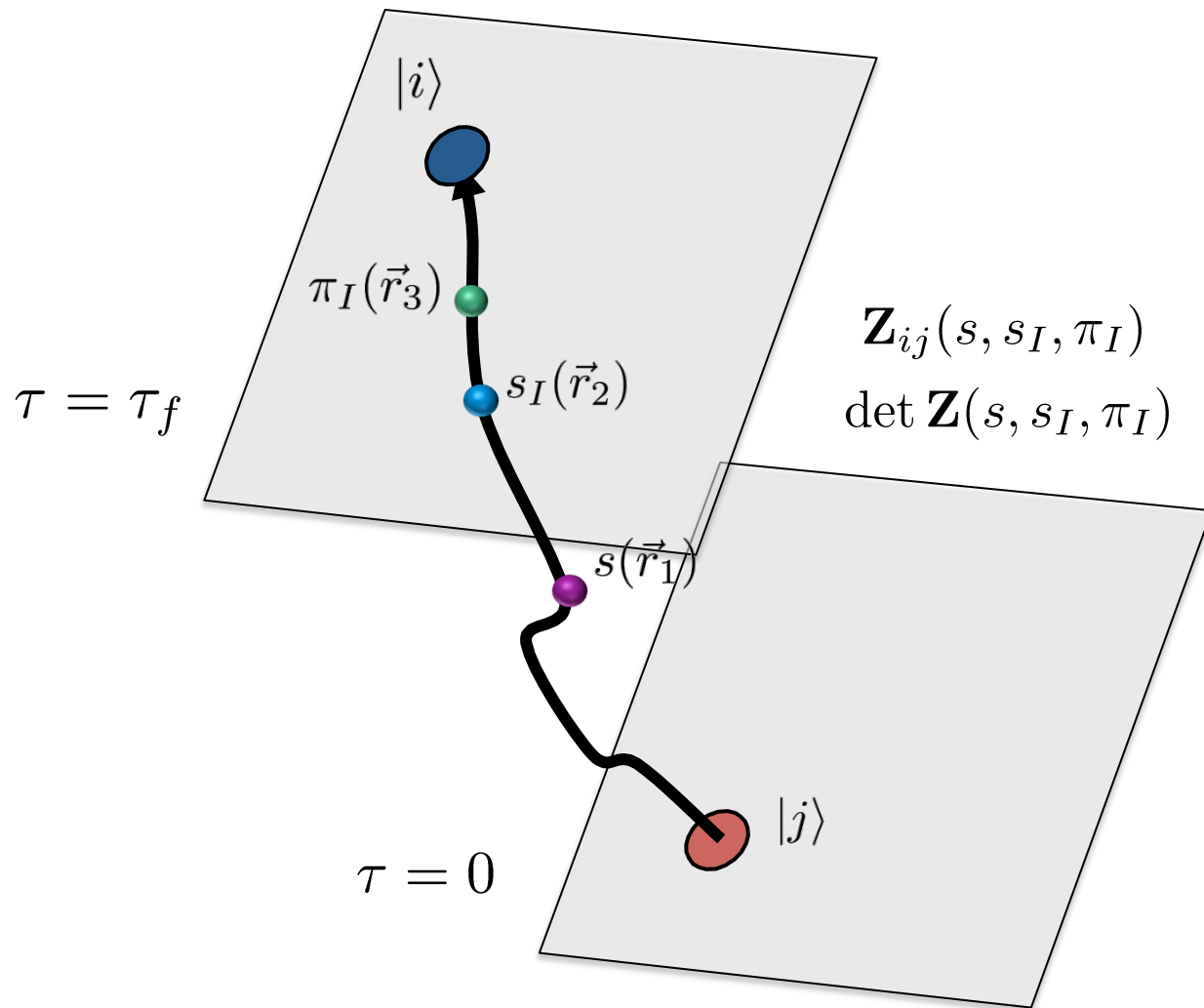


## Auxiliary field method

We can write exponentials of the interaction using a Gaussian integral identity

$$\begin{aligned} & \exp \left[ -\frac{C}{2} (N^\dagger N)^2 \right] \quad \times \quad (N^\dagger N)^2 \\ &= \sqrt{\frac{1}{2\pi}} \int_{-\infty}^{\infty} ds \exp \left[ -\frac{1}{2} s^2 + \sqrt{-C} s (N^\dagger N) \right] \quad \rangle \quad s N^\dagger N \end{aligned}$$

We remove the interaction between nucleons and replace it with the interactions of each nucleon with a background field.



## Pinhole algorithm

Consider the density operator for nucleon with spin  $i$  and isospin  $j$

$$\rho_{i,j}(\mathbf{n}) = a_{i,j}^\dagger(\mathbf{n})a_{i,j}(\mathbf{n})$$

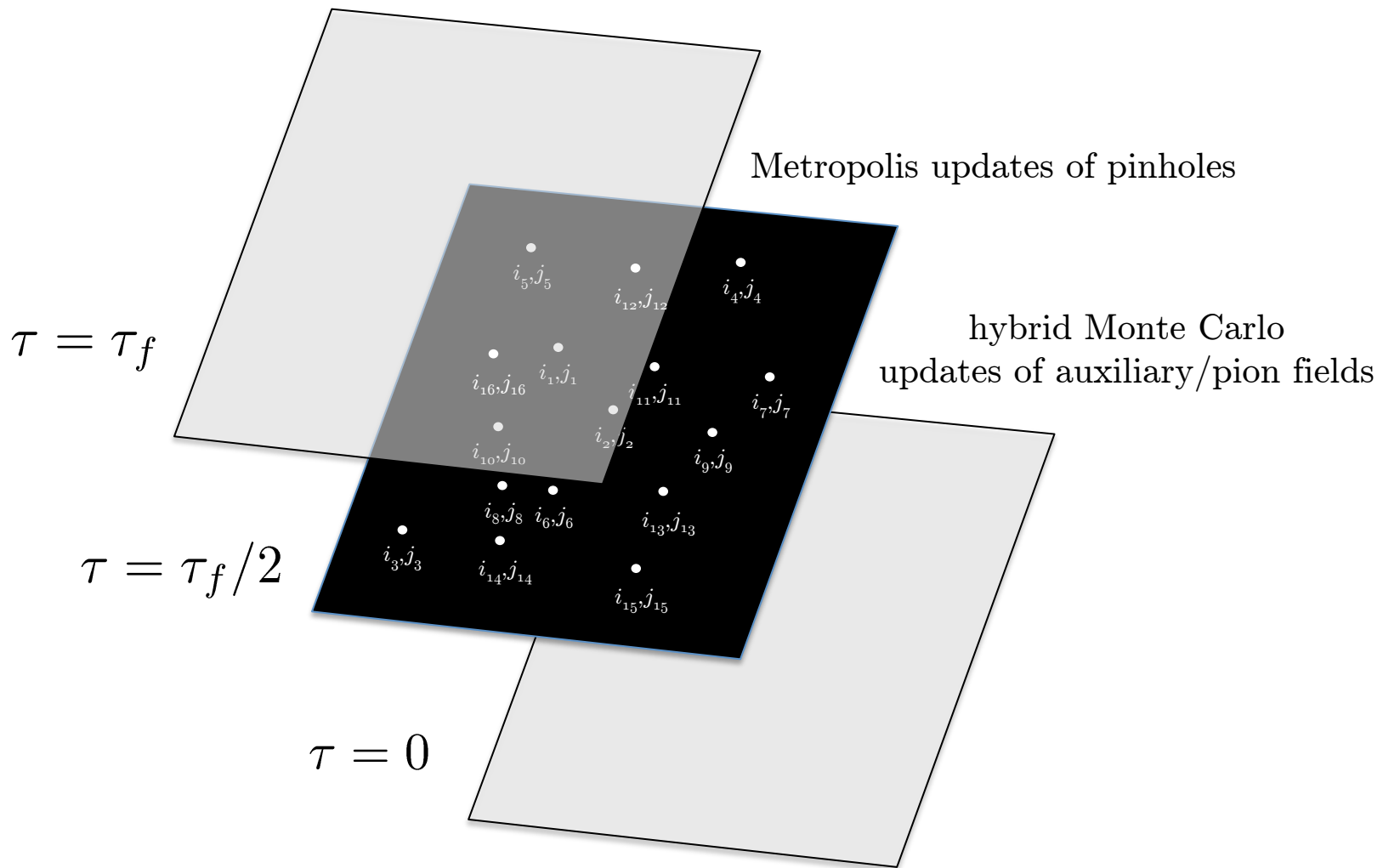
We construct the normal-ordered  $A$ -body density operator

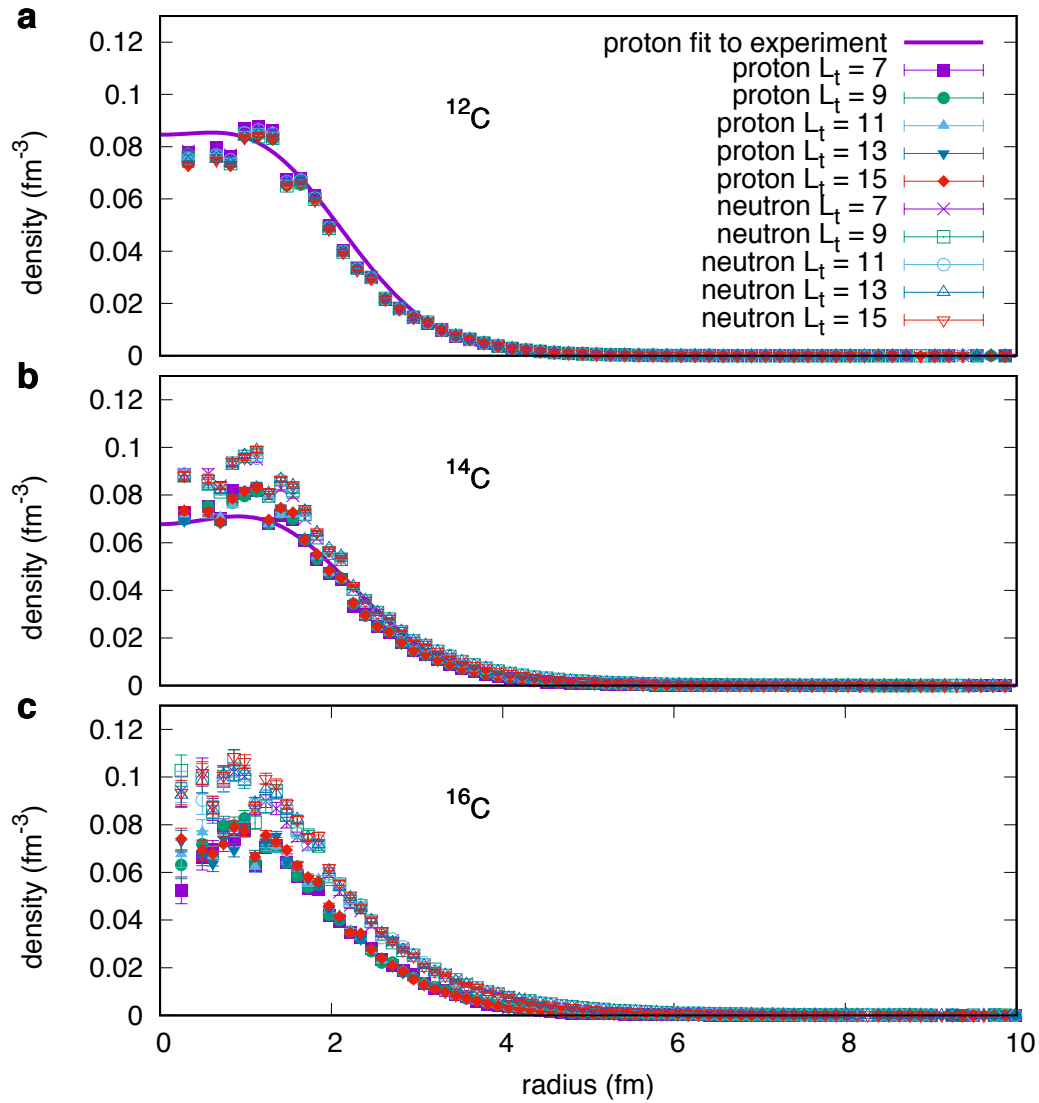
$$\rho_{i_1,j_1,\dots,i_A,j_A}(\mathbf{n}_1,\dots,\mathbf{n}_A) = : \rho_{i_1,j_1}(\mathbf{n}_1) \cdots \rho_{i_A,j_A}(\mathbf{n}_A) :$$

In the simulations we do Monte Carlo sampling of the amplitude

$$A_{i_1,j_1,\dots,i_A,j_A}(\mathbf{n}_1,\dots,\mathbf{n}_A,t) = \langle \Psi_I | e^{-Ht/2} \rho_{i_1,j_1,\dots,i_A,j_A}(\mathbf{n}_1,\dots,\mathbf{n}_A) e^{-Ht/2} | \Psi_I \rangle$$







Elhatisari, Epelbaum, Krebs, Lähde, D.L., Li, Lu, Meißner, Rupak, PRL 119, 222505 (2017)

## Challenge

In order to compute thermodynamic properties of finite nuclei, nuclear matter, and neutron matter, we need to compute the partition function

$$\text{Tr exp}(-\beta H)$$

The standard method for computing the partition function involves calculating determinants of matrices of size  $4V \times 4V$ , where  $V$  is the number of lattice points filling the spatial volume. Since  $V$  is usually several hundred or several thousand, these calculations are very expensive.

## Pinhole trace algorithm

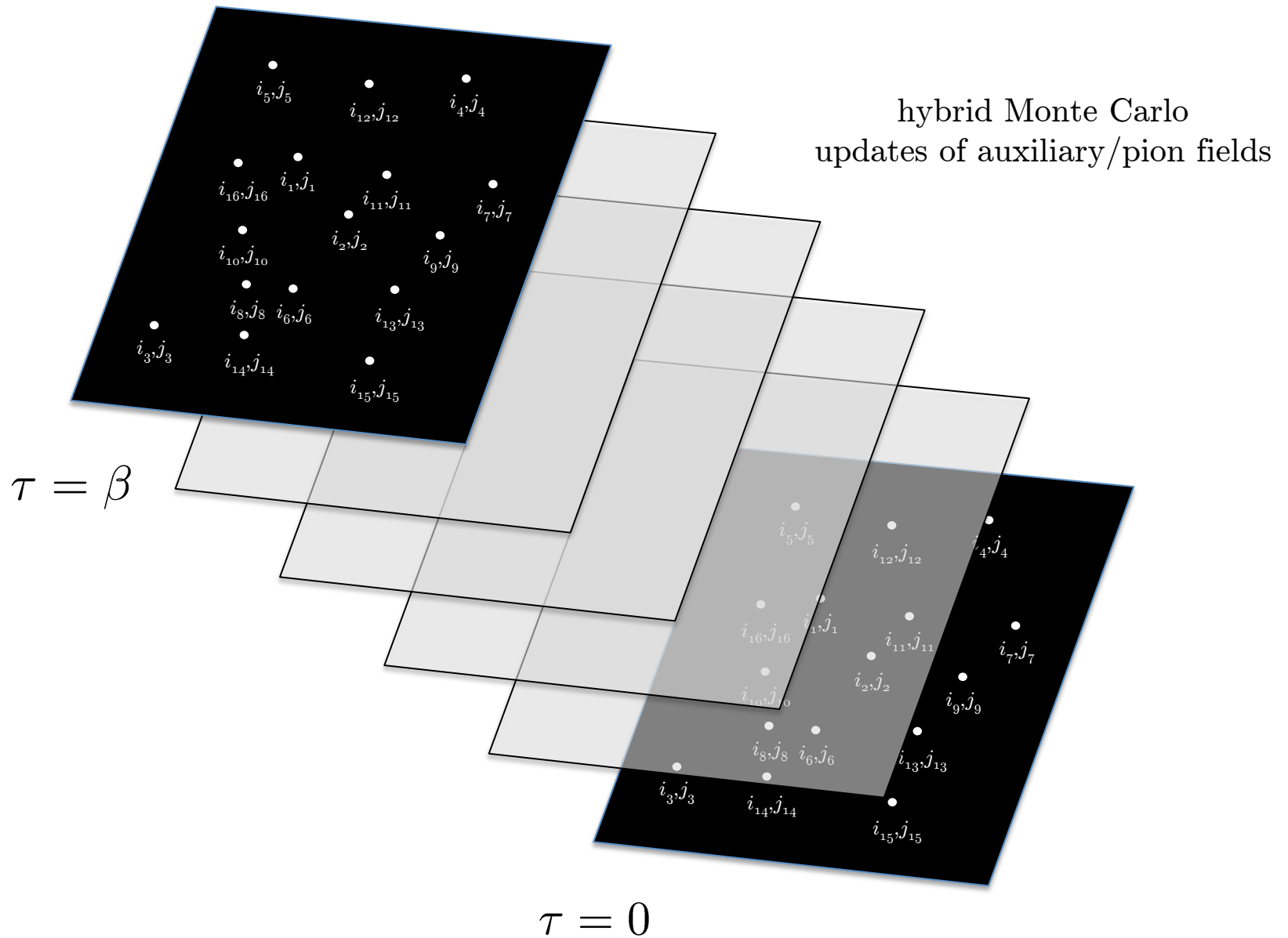
We have developed an alternative method using pinholes that calculates determinants of matrices of size  $A \times A$ , where  $A$  is the number of nucleons. The method does not suffer from severe sign oscillations.

We compute the quantum mechanical trace over  $A$ -nucleon states by summing over pinholes (position eigenstates) for the initial and final states

$$\begin{aligned} & \text{Tr } O \\ &= \frac{1}{A!} \sum_{i_1 \cdots i_A, j_1 \cdots j_A, \mathbf{n}_1 \cdots \mathbf{n}_A} \langle 0 | a_{i_A, j_A}(\mathbf{n}_A) \cdots a_{i_1, j_1}(\mathbf{n}_1) O a_{i_1, j_1}^\dagger(\mathbf{n}_1) \cdots a_{i_A, j_A}^\dagger(\mathbf{n}_A) | 0 \rangle \end{aligned}$$

This can be used to calculate the partition function in the canonical ensemble.

# Metropolis updates of pinholes



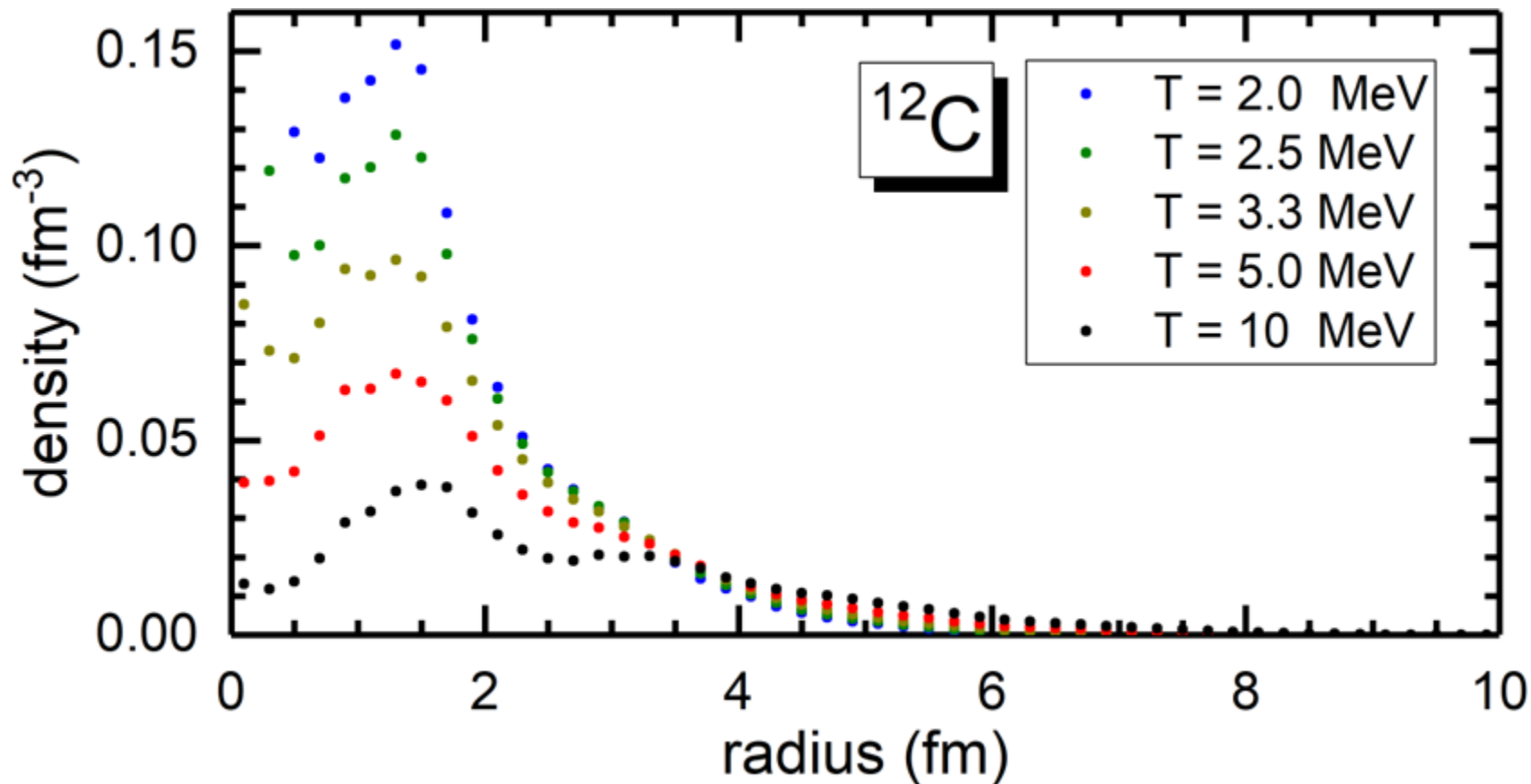


Figure courtesy of Bingnan Lu

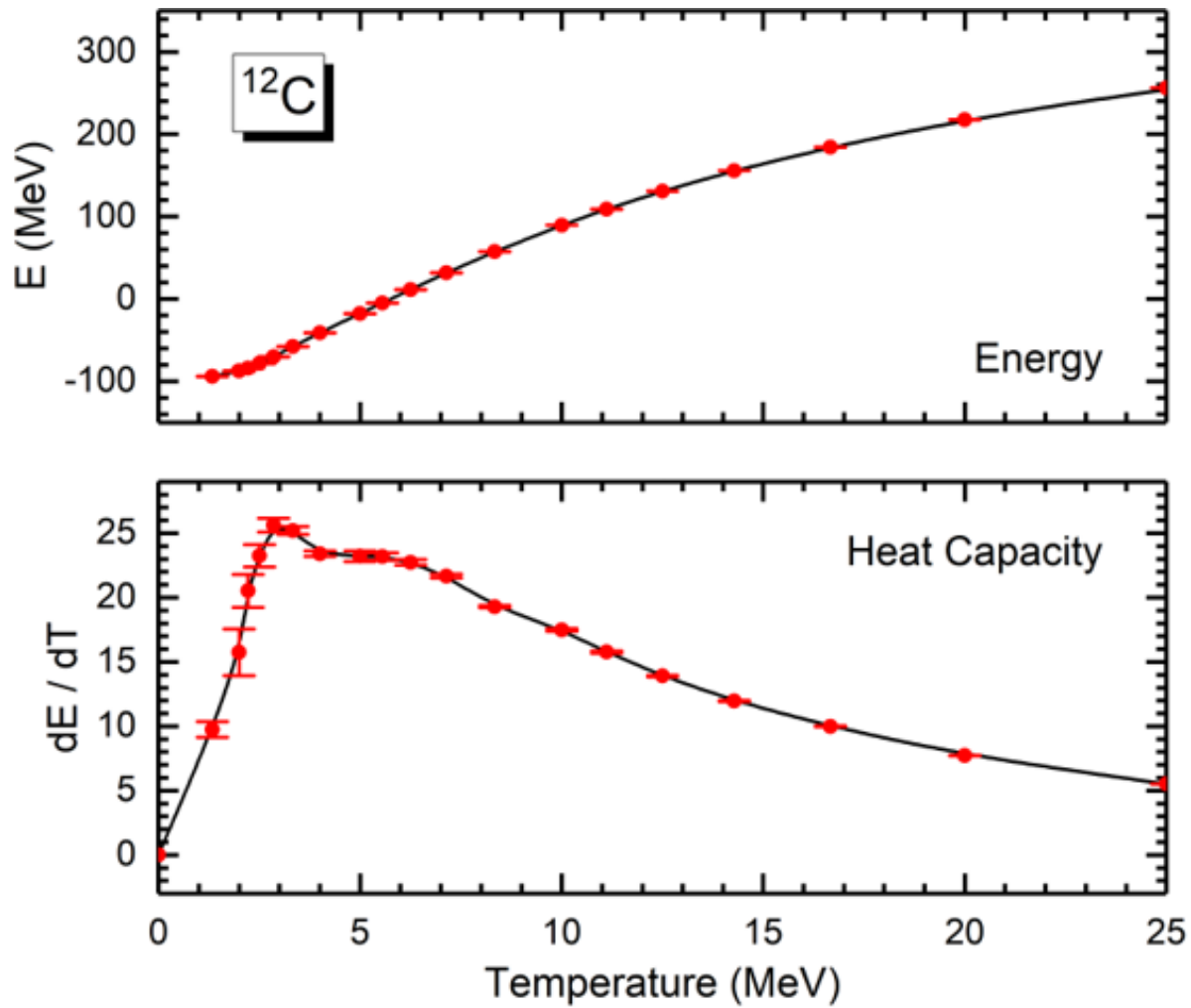


Figure courtesy of Bingnan Lu

## Eigenvector continuation

We demonstrate that when a control parameter in the Hamiltonian matrix is varied smoothly, the extremal eigenvectors do not explore the large dimensionality of the linear space. Instead they trace out trajectories with significant displacements in only a small number of linearly-independent directions.

We prove this empirical observation using analytic function theory and the principles of analytic continuation.

Since the eigenvector trajectory is a low-dimensional manifold embedded in a very large space, we can find the desired eigenvector using methods similar to image recognition in machine learning.

D. Frame, R. He, I. Ipsen, Da. Lee, De. Lee, E. Rrapaj, arXiv:1711.07090



Consider a one-parameter family of Hamiltonian matrices of the form

$$H(c) = H_0 + cH_1$$

where  $H_0$  and  $H_1$  are Hermitian. Let the eigenvalues and eigenvectors be

$$H(c)|\psi_j(c)\rangle = E_j(c)|\psi_j(c)\rangle$$

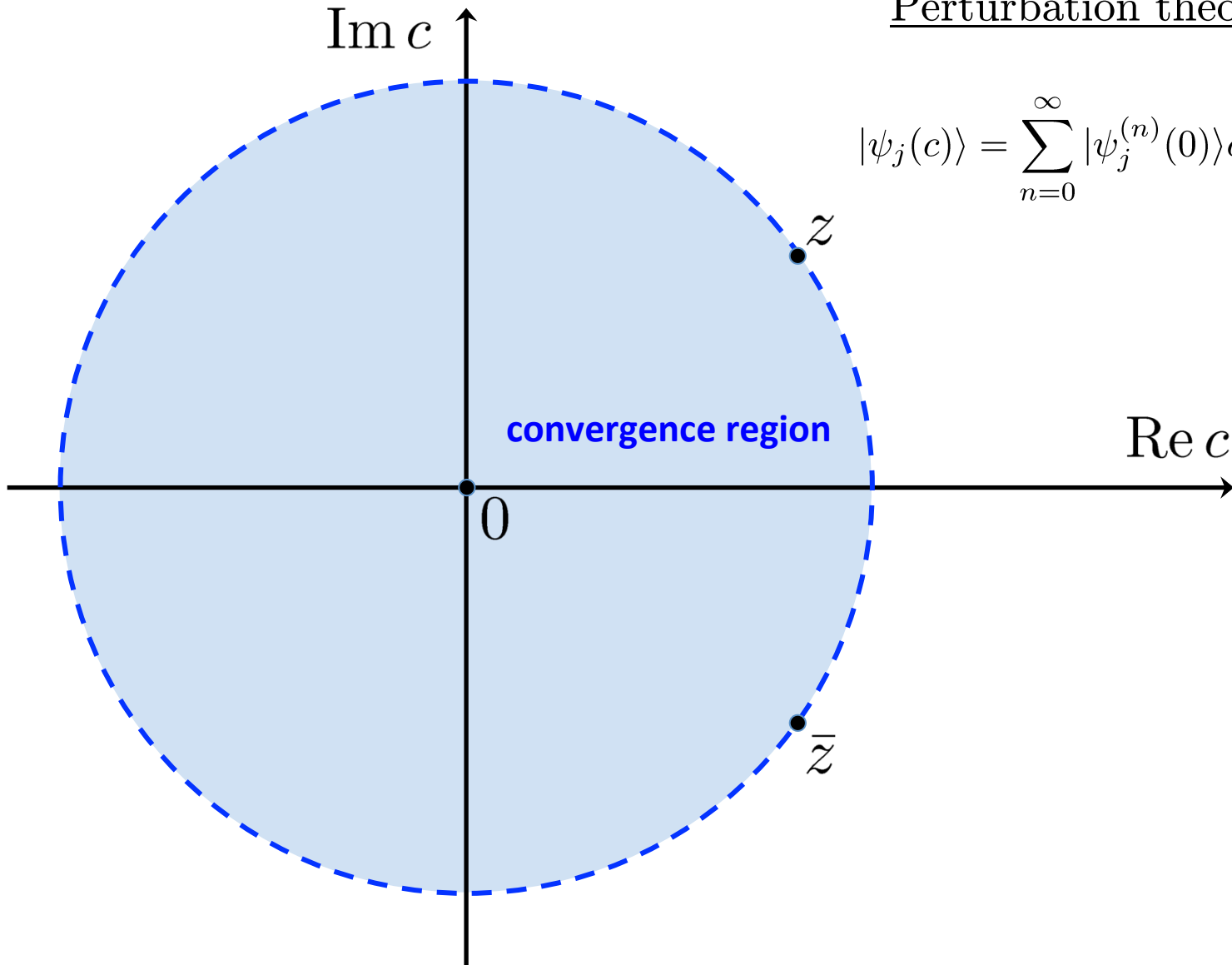
We can perform series expansions around the point  $c = 0$ .

$$E_j(c) = \sum_{n=0}^{\infty} E_j^{(n)}(0)c^n/n!$$
$$|\psi_j(c)\rangle = \sum_{n=0}^{\infty} |\psi_j^{(n)}(0)\rangle c^n/n!$$

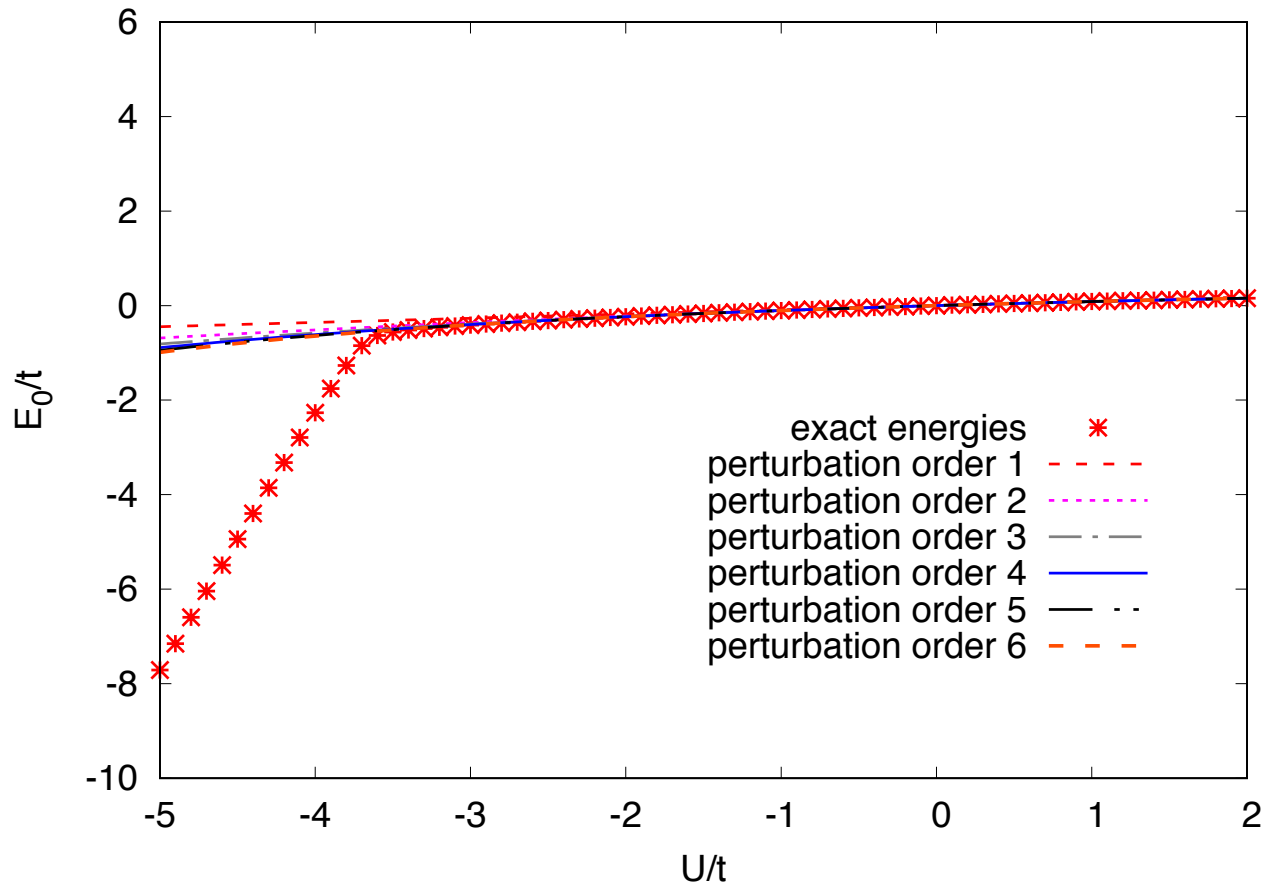
This is the strategy of perturbation theory. We can compute each term in the series when the eigenvalues and eigenvectors of  $H_0$  are known or computable.

## Perturbation theory

$$|\psi_j(c)\rangle = \sum_{n=0}^{\infty} |\psi_j^{(n)}(0)\rangle c^n / n!$$

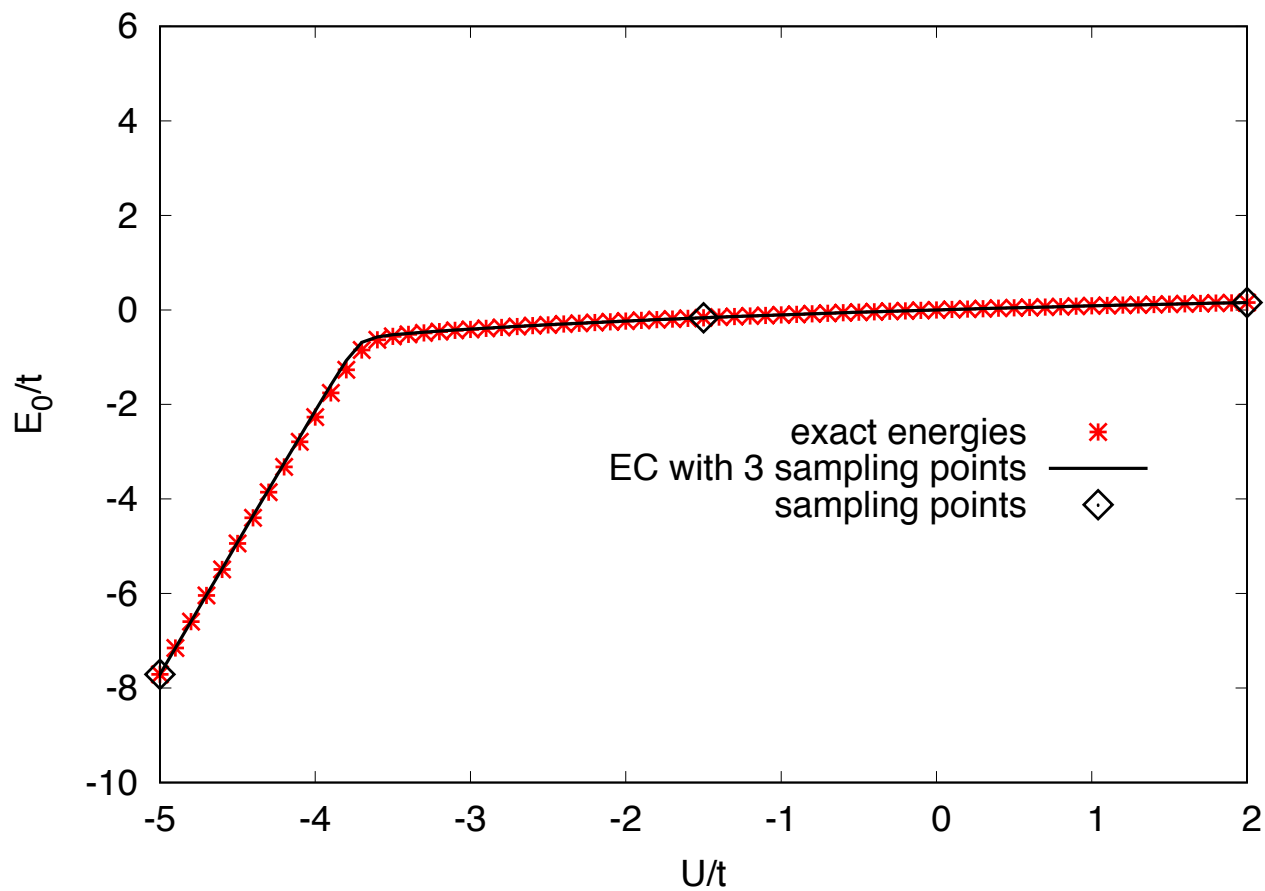


# Bose-Hubbard model

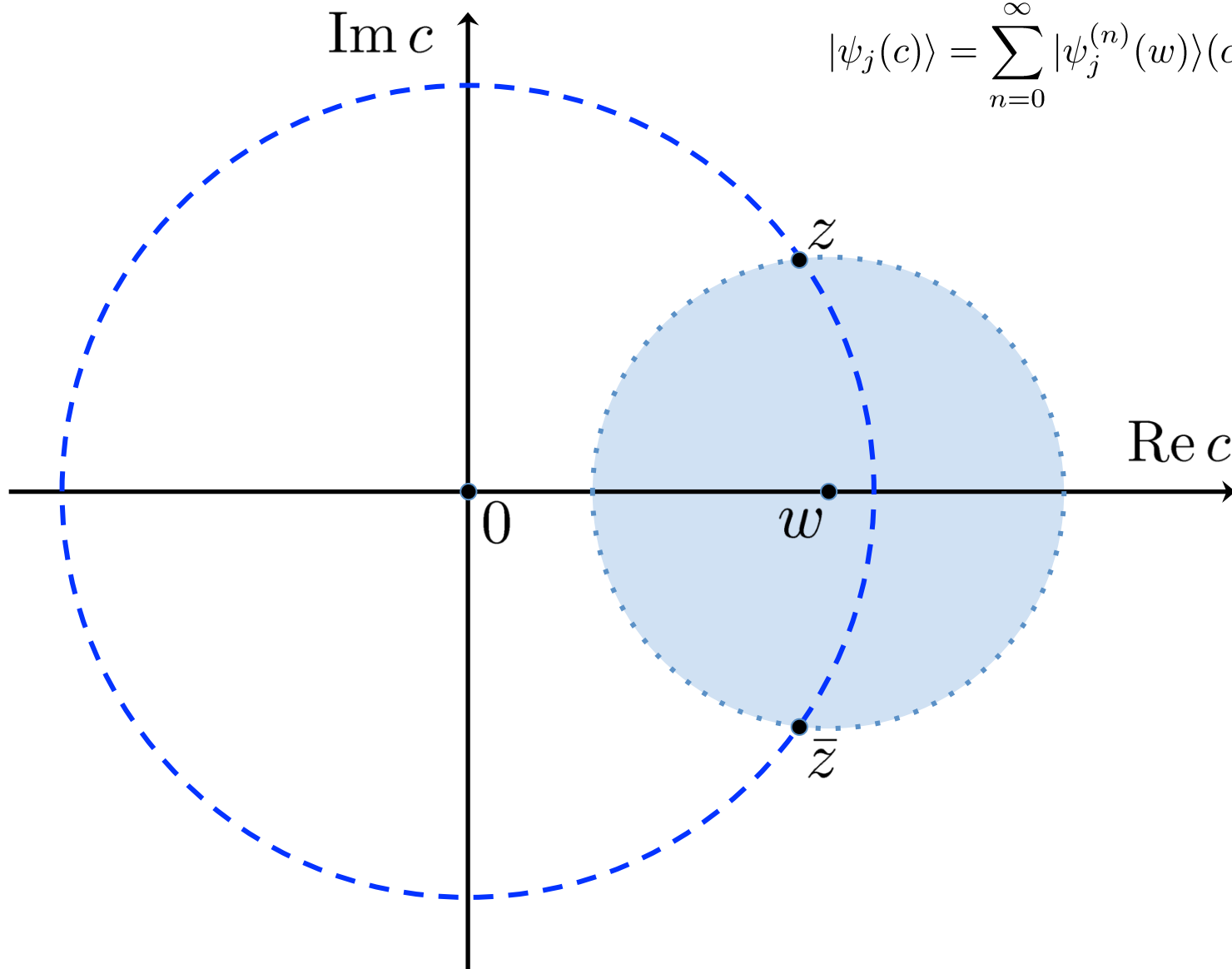


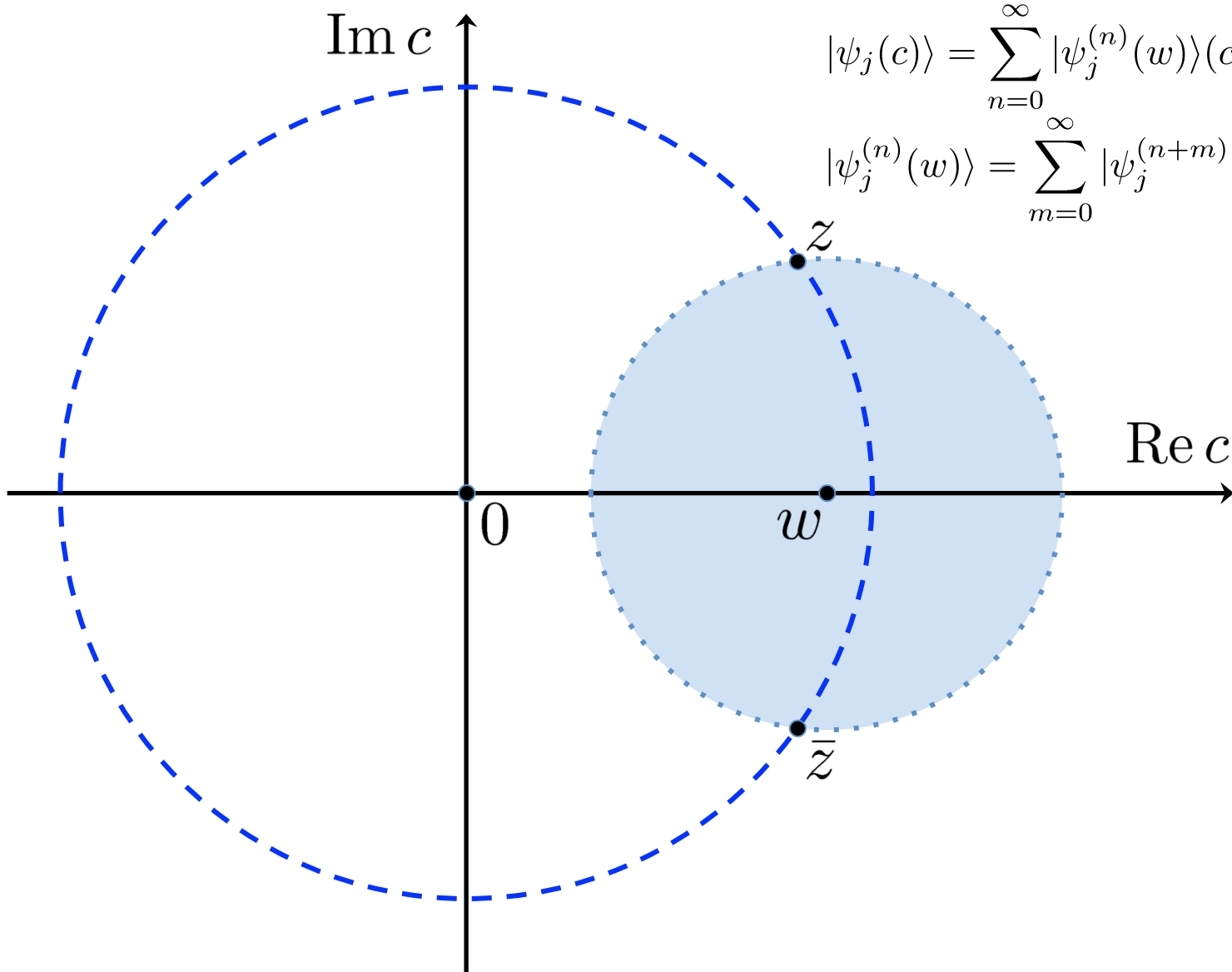
Perturbation theory fails at strong attractive coupling

Restrict the linear space to the span of three vectors



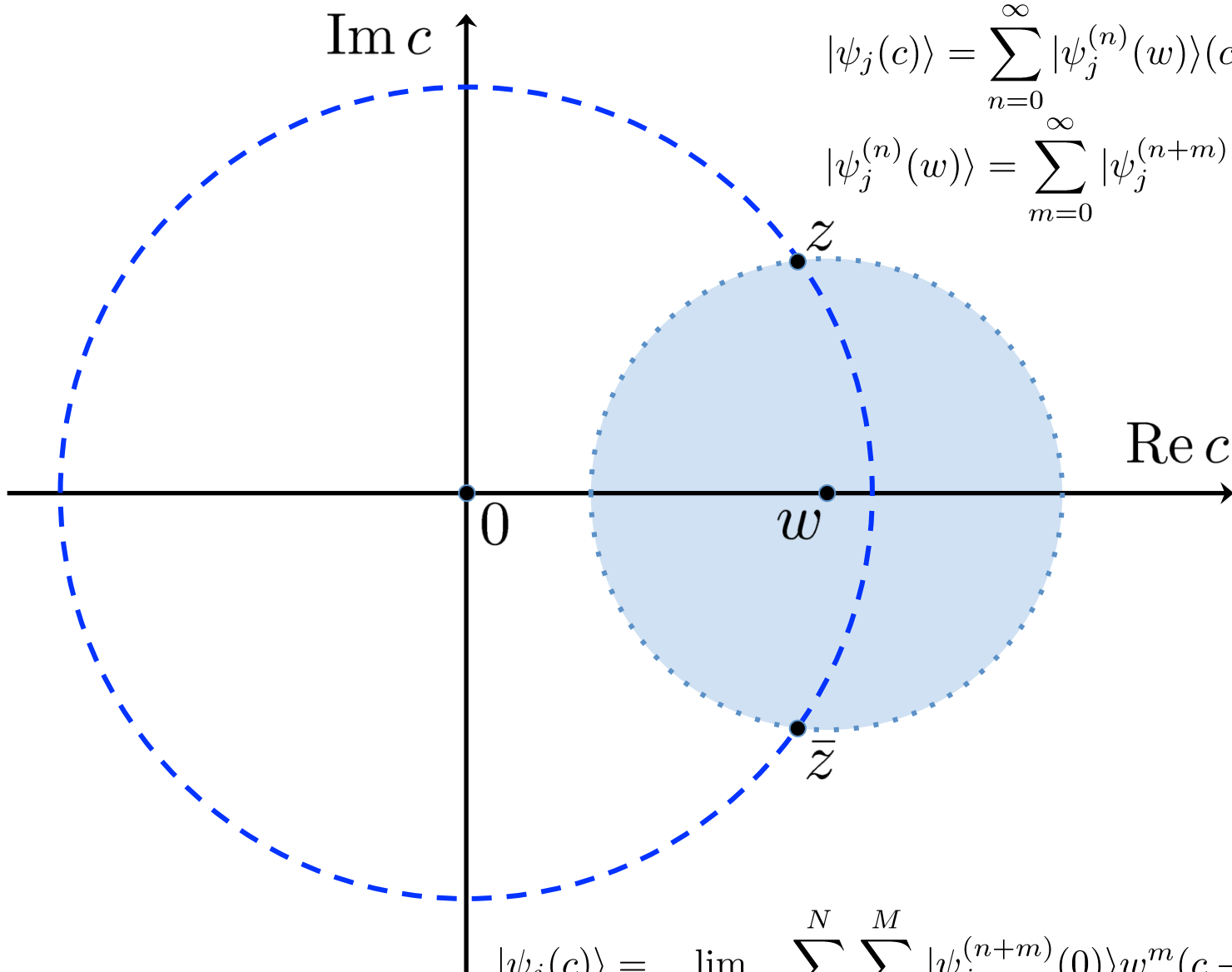
$$|\psi_j(c)\rangle = \sum_{n=0}^{\infty} |\psi_j^{(n)}(w)\rangle (c-w)^n / n!$$





$$|\psi_j(c)\rangle = \sum_{n=0}^{\infty} |\psi_j^{(n)}(w)\rangle (c-w)^n / n!$$

$$|\psi_j^{(n)}(w)\rangle = \sum_{m=0}^{\infty} |\psi_j^{(n+m)}(0)\rangle w^m / m!$$



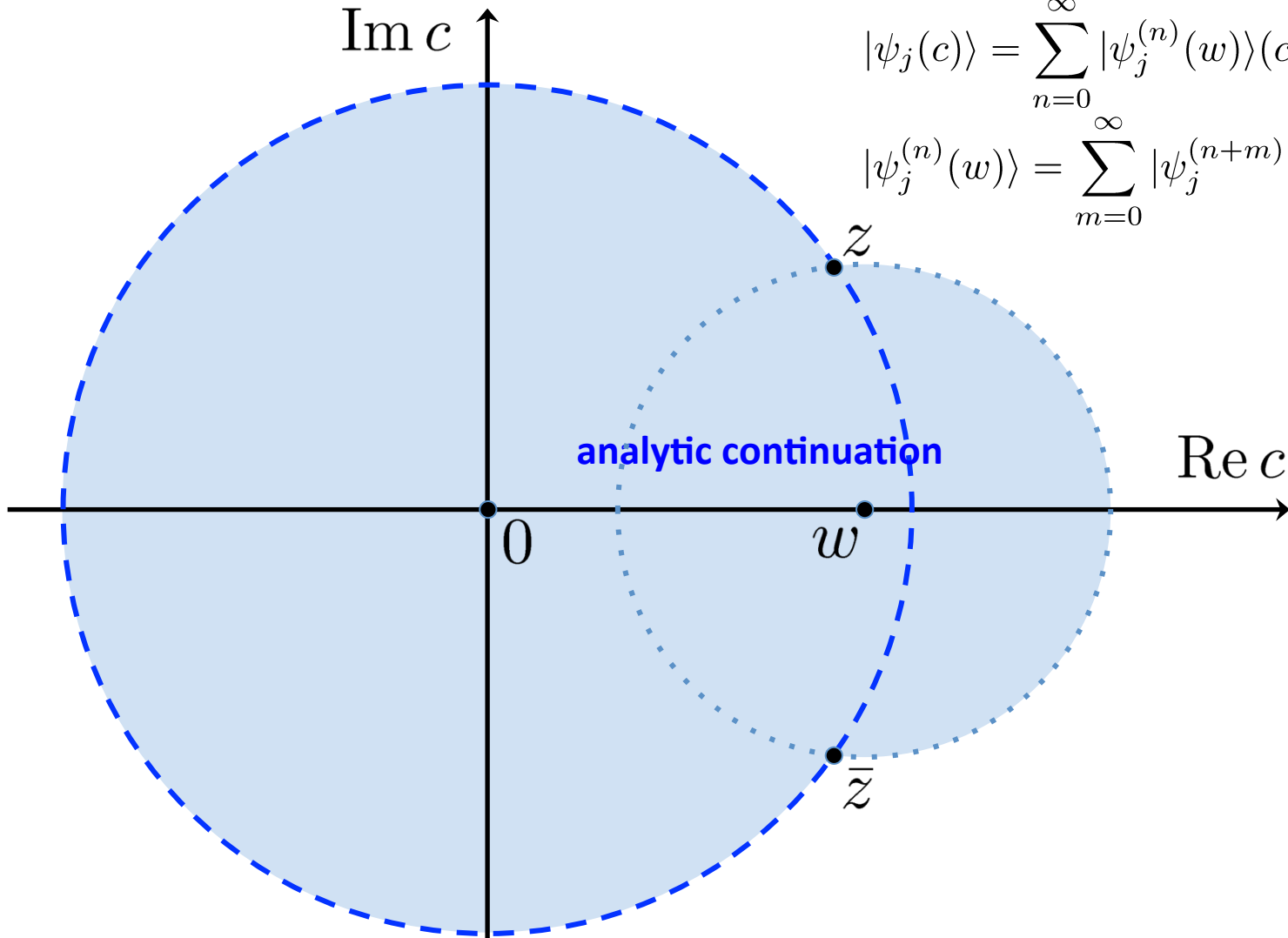
$$|\psi_j(c)\rangle = \sum_{n=0}^{\infty} |\psi_j^{(n)}(w)\rangle (c-w)^n / n!$$

$$|\psi_j^{(n)}(w)\rangle = \sum_{m=0}^{\infty} |\psi_j^{(n+m)}(0)\rangle w^m / m!$$

$$|\psi_j(c)\rangle = \lim_{N, M \rightarrow \infty} \sum_{n=0}^N \sum_{m=0}^M |\psi_j^{(n+m)}(0)\rangle w^m (c-w)^n / (m! n!)$$

$$|\psi_j(c)\rangle = \sum_{n=0}^{\infty} |\psi_j^{(n)}(w)\rangle (c-w)^n / n!$$

$$|\psi_j^{(n)}(w)\rangle = \sum_{m=0}^{\infty} |\psi_j^{(n+m)}(0)\rangle w^m / m!$$



$$|\psi_j(c)\rangle = \sum_{n=0}^{\infty} |\psi_j^{(n)}(0)\rangle c^n / n!$$

$$|\psi_j(c)\rangle = \lim_{N, M \rightarrow \infty} \sum_{n=0}^N \sum_{m=0}^M |\psi_j^{(n+m)}(0)\rangle w^m (c-w)^n / (m! n!)$$



The eigenvector can be well approximated as a linear combination of a few vectors, using either the original series expansion

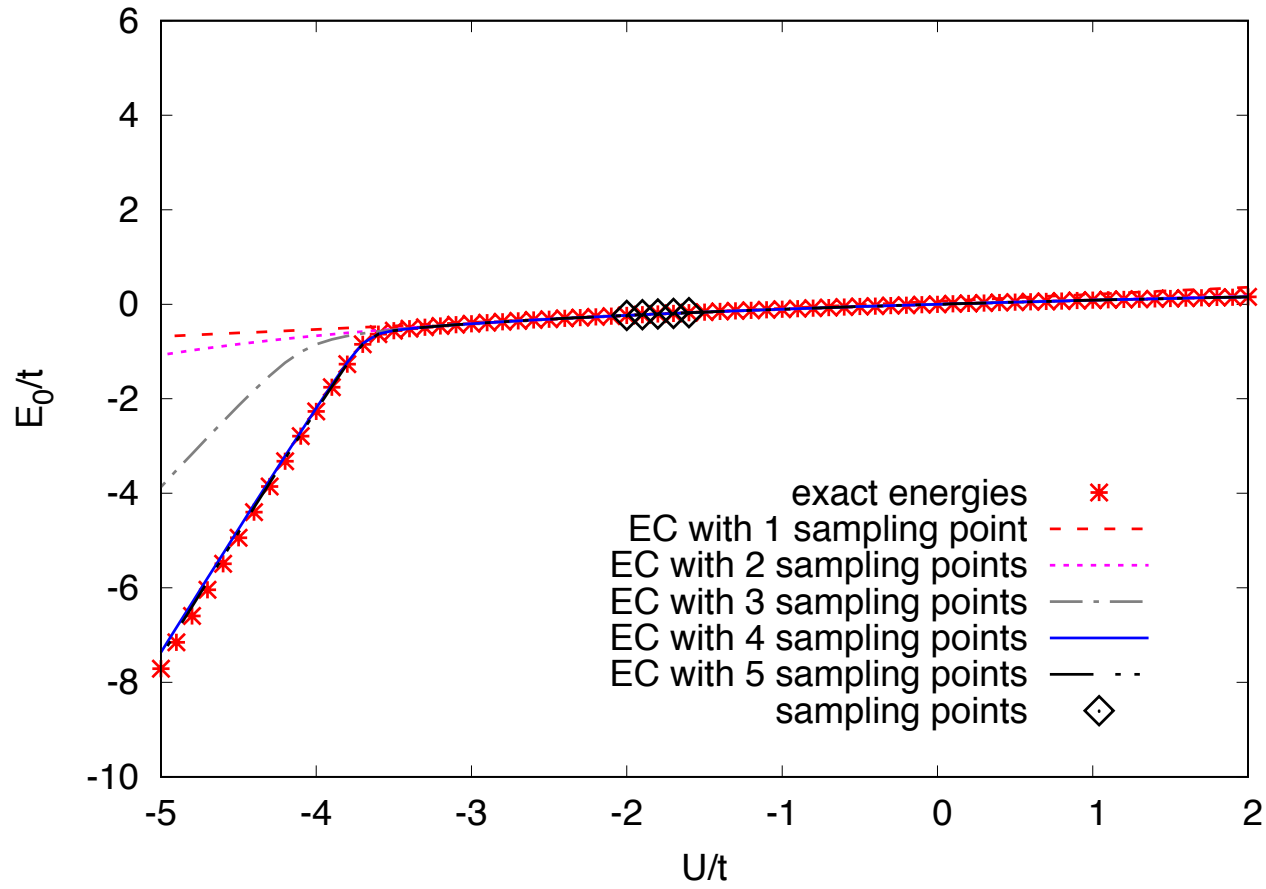
$$|\psi_j(c)\rangle = \sum_{n=0}^{\infty} |\psi_j^{(n)}(0)\rangle c^n / n!$$

or the rearranged multi-series expansion we obtained through analytic continuation

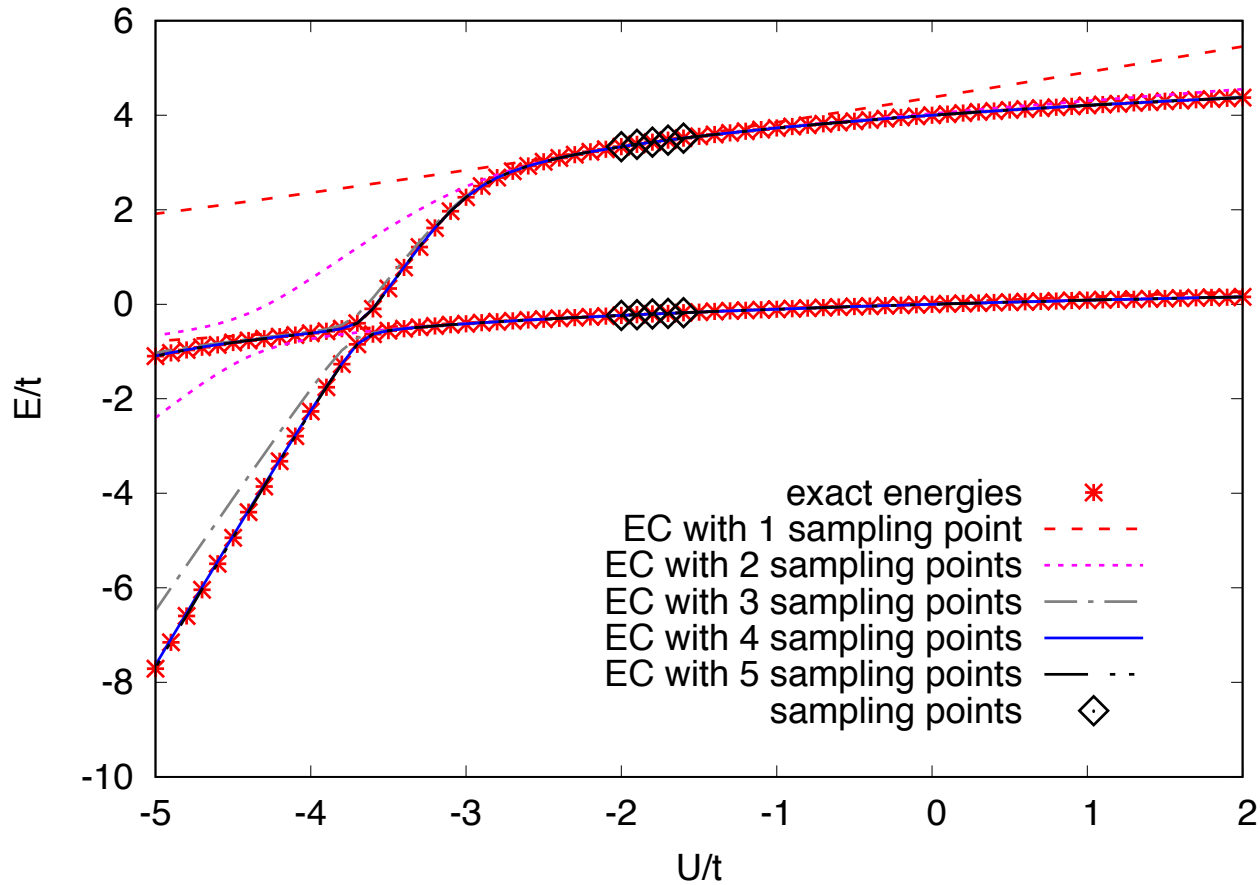
$$|\psi_j(c)\rangle = \lim_{N, M \rightarrow \infty} \sum_{n=0}^N \sum_{m=0}^M |\psi_j^{(n+m)}(0)\rangle w^m (c-w)^n / (m! n!)$$

As  $c$  is varied the eigenvector does not explore the large dimensionality of the linear space, but is instead well approximated by a low-dimension manifold.

We can “learn” the eigenvector trajectory in one region and perform eigenvector continuation to another region



Applying eigenvector continuation to more than one eigenvector at a time accelerates convergence near avoided level crossings.



## Superfluidity and pairing correlations

The two-body density matrix is defined as

$$\rho_2(\vec{r}'_1, \vec{r}'_2, \vec{r}_1, \vec{r}_2) = \left\langle \psi_{\downarrow}^{\dagger}(\vec{r}'_2) \psi_{\uparrow}^{\dagger}(\vec{r}'_1) \psi_{\uparrow}(\vec{r}_1) \psi_{\downarrow}(\vec{r}_2) \right\rangle$$

Long-range correlations in the two-body density matrix is a signature for pair superfluidity:

$$\begin{aligned} \rho_2(\vec{r}'_1, \vec{r}'_2, \vec{r}_1, \vec{r}_2) &\rightarrow \alpha N/2 \cdot \phi^*(|\vec{r}'_1 - \vec{r}'_2|) \phi(|\vec{r}_1 - \vec{r}_2|) \\ |\vec{r}_1 - \vec{r}'_1|, |\vec{r}_2 - \vec{r}'_2| &\rightarrow \infty \end{aligned}$$

Yang, RMP 34, 694 (1962)

In progress: Two-body density matrix in neutron matter

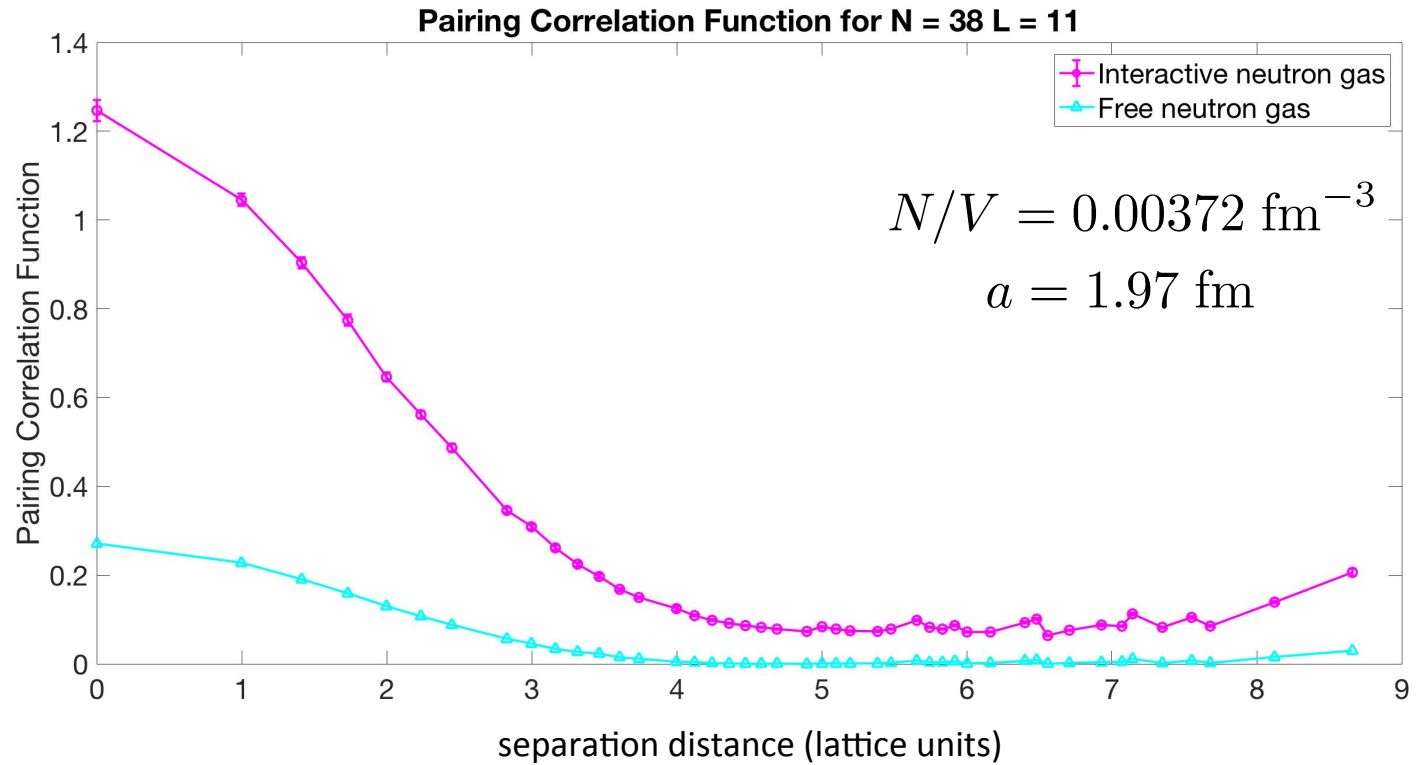


Figure courtesy of Rongzheng He

## Summary and Outlook

These are exciting times for the *ab initio* nuclear theory community. In lattice EFT, we have new projects in motion which are pushing the current frontiers.

Currently working to improve our understanding of the detailed connection between bare nuclear forces and nuclear structure for light and medium-mass nuclei.

Applying the adiabatic projection method to low-energy nucleon-nucleus and alpha-nucleus scattering and reactions.

Using the pinhole algorithm to study the detailed structure of nuclei and thermodynamics of finite nuclei, nuclear matter, and neutron matter.

Implementing eigenvector continuation to treat all higher-order interactions in chiral effective field theory.

Calculating the two-body density matrix to measure pairing correlations in neutron matter and finite nuclei.