

# HOBET: QCD to Effective Interaction and Effective Operators 

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## Two Paths to Connect QCD to the Effective Interaction

* HOBET (Harmonic-Oscillator-Based Effective Theory)
* Compute phase shifts in LQCD - Fit HOBET LECs
* See arXiv:1511.02262 (CalLat LQCD)+ arXiv: 1902.03543 (McElvain, Haxton)
* Compute nucleon scattering spectrum in LQCD - Fit HOBET LECs directly to spectrum + periodic boundary conditions.


## The Bloch-Horowitz Equation

* For practical calculation reasons we often want to work in a subspace of the full Hilbert space.
* P projects the subspace and $\mathrm{Q}=1$ $P$ gets the rest.
* The BH equation is the answer to the question: Does there exist an operator Heff that lives in P with the same eigenvalues and projected eigenvectors of the full H .

$$
\operatorname{Insert}(P+Q) \operatorname{in} H\left|\psi_{i}\right\rangle=E_{i}\left|\psi_{i}\right\rangle
$$

$$
\begin{aligned}
& P H P\left|\psi_{i}\right\rangle+P H Q\left|\psi_{i}\right\rangle=E_{i} P\left|\psi_{i}\right\rangle \\
& Q H P\left|\psi_{i}\right\rangle+Q H Q\left|\psi_{i}\right\rangle=E_{i} Q\left|\psi_{i}\right\rangle
\end{aligned}
$$

$$
Q\left|\psi_{i}\right\rangle=\frac{1}{E_{i}-Q H} Q H P\left|\psi_{i}\right\rangle
$$

$$
P\left(H+H \frac{1}{E_{i}-Q H} Q H\right) P\left|\psi_{i}\right\rangle=E_{i} P\left|\psi_{i}\right\rangle
$$

$$
H^{e f f}(E) P\left|\psi_{i}\right\rangle=P H \frac{1}{E_{i}-Q H} P\left|\psi_{i}\right\rangle=E_{i} P\left|\psi_{i}\right\rangle
$$

## BH Characteristics

* Eigenstates of $H^{\operatorname{eff}}(\mathrm{E})$ are projections with the same eigenvalues.
* All eigenstates that overlap P are included!
* True even if P projects a finite number of states.
* It is continuous in energy, including across $\mathrm{E}=0$. An effective theory based on the BH equation can be fit in the continuum and used to find bound states.
* Explicitly energy dependent: Must solve self consistently.
* Simple fixed point iteration converges rapidly.


## HO Effective Theory

*Why the HO basis?

* Discrete so we can use matrix techniques for solution.
* Good for confined wave function of nucleus
* With a consistent A-body quanta cutoff the center of mass is separable.
* In an HO ET with included space projector $\mathrm{P}(\Lambda, \mathrm{b})$, both UV and IR are excluded.
* Major Issue
* The kinetic energy operator T is a hopping operator, strongly connecting P \& Q (IR).


## HOBET Introduction

* HOBET is based on a reorganization of the BlochHorowitz equation by Haxton and Luu.

$$
H^{f f i \wedge}(E)=P\left[H \frac{E}{E-Q H}\right] P=P \frac{E}{E-T Q}\left[T+T \frac{Q}{E} T+V+V \frac{1}{E-Q H} Q V\right] \frac{E}{E-Q T} P
$$

* The reorganization isolates the impact of T for analytic calculation to all orders.
* The remaining part is replaced by a long range potential (like an OPEP) plus $\mathrm{V}_{\delta,}$ which is a short range expansion around it.


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## E/(E-QT) Transform of Edge States



- Acting on edge state with
$E / \hbar \omega=1 / 2$.
Recovers scattering wave function with phase shift.

* Acting on edge state with $E / \hbar \omega=-1 / 2$.

Recovers bound state exponential decay from gaussian falloff of HO state.

$$
G_{Q T} P=\frac{E}{E-Q T} P=\frac{E}{E-T}\left\{P \frac{E}{E-T} P\right\}^{-1}, \quad b_{i j}=\left\{P \frac{E}{E-T} P\right\}_{i j}^{-1}
$$

## Sum T to All Orders

* T contributions can be summed to all orders.

$$
\langle j| \frac{E}{E-T Q}\left[T-T \frac{Q}{E} T\right] \frac{E}{E-Q T}|i\rangle=E\left(\delta_{j i}-b_{j i}\right)
$$

* A surprisingly simple result.
* A non-perturbative sum of kinetic energy scattering is key to a convergent ET expansion of the remaining parts.


## The $\mathrm{V}_{\delta}$ Expansion

* $\mathrm{V}_{\delta}$ is described in terms of HO lowering operators.

$$
\begin{aligned}
\hat{C} & \text { lowers } L, \hat{A} \text { lowers nodal } n, \quad[\hat{C}, \hat{A}]=0 \\
V_{\delta}^{S} & =a_{L O}^{s} \delta(r)+a_{\text {NLO }}^{s}\left(\hat{A}^{\dagger} \delta(r)+\delta(r) \hat{A}\right)+\ldots \\
V_{\delta}^{\text {SD }} & =a_{\text {NLO }}^{\text {SD }}\left(\hat{C}^{+2} \delta(r)+\delta(r) \hat{C}^{2}\right)+a_{\text {NNLO }}^{22, S D}\left(\hat{C}^{+2} \delta(r) \hat{A}+\hat{A}^{\dagger} \delta(r) \hat{C}^{2}\right) \\
& +a_{\text {NNLO }}^{00, S D}\left(\hat{C}^{+2} \hat{A}^{+} \delta(r)+\delta(r) \hat{A} \hat{C}^{2}\right)+\ldots
\end{aligned}
$$

* This is slightly simplified by absorbing a constant related to coupling spins to angular momentum into the LECs.
* $\left[\hat{C}^{2}\right.$ is really $\hat{C}^{2} \equiv[\tilde{a} \otimes \tilde{a}]^{(2)} \odot\left[\vec{\sigma}_{1} \otimes \vec{\sigma}_{2}\right]^{(2)}$, coupling angular momentum to spins, with vector HO lowering op $\tilde{a}$ ]


## HOBET Lepage Plots





* Heff S-D matrix elements in $\Lambda=8$ are are directly calculated at $\mathrm{E}=-2.2245 \mathrm{MeV}$ from HO matrix elements in $\Lambda=400$ and LECs fit in a scheme independent way.
* For the middle plot, $\mathrm{O}(1)$ errors appear at $\mathrm{n}^{\prime}+\mathrm{n}=8$ making diagonal matrix elements with $\mathrm{n}=4$ unreliable. $\mathrm{n}=3$ at $4 \hbar \omega=60 \mathrm{MeV}$ is a reasonable breakdown scale.
* Convergence is good order by order, but the value of a good long range VIR is clear.


## Power Counting

* The expansion is for $V_{\delta}$, whose range $R$ is shorter than that of $V$.
* For a known $V_{\delta}$ the LECs are proportional to a non-local Talmi integral.
- The tail overlap shrinks rapidly with order.
* For a short range $V_{\delta}$ the expansion parameter is function of $b / R$.

$$
\begin{aligned}
& V_{\delta}\left(r^{\prime}, r\right)=V-V_{I R}+V \frac{1}{E-Q H} Q V \\
& L E C_{n^{\prime}, n} \propto \int r^{\prime 2} d r^{\prime} r^{2} d r r^{2(n-1)} e^{-r / 2} V_{Q} e^{-r / 2} r^{2(n-1)}
\end{aligned}
$$

## Energy Independence of LECs

* Enables fitting to data a range of energies.
* The upper blue dots are the result of solving for aLo at individual samples $\left(\mathrm{E}_{\mathrm{i}}, \delta_{\mathrm{i}}\right)$

* The lower gold dots represent an NLO fit to data from 1 and 10 MeV to determine $\mathrm{a}_{\text {NLO }}$ followed by refitting a $\mathrm{L}_{\mathrm{LO}}$ at each energy while holding $a_{\text {NLO }}$ constant.
* Conclusion: Energy dependence is adsorbed into higher order operators.


## Fitting LECs

* Principle: The BH equation is energy self consistent

$$
H_{e f f}^{\text {fill }}\left(E_{i}\right) P\left|\psi_{i}\right\rangle=E_{i} P\left|\psi_{i}\right\rangle
$$

* At fixed order we instead have a nearby eigenstate.

$$
H_{e f f}\left(E_{i}, L E C s\right) P\left|\psi_{i}^{\prime}\right\rangle=\varepsilon_{i} P\left|\psi_{i}^{\prime}\right\rangle
$$

* The mismatch must be due to LEC values.
- Repair by minimizing

$$
\sum_{i \in \text { samples }}\left(\varepsilon_{i}-E_{i}\right)^{2} / \sigma_{i}^{2}
$$

* The variance for the difference can be estimated from the sensitivity of $\varepsilon_{i}$ to next order LECs, automatically suppressing data outside the validity range for the current LEC order.


## Predicting the Deuteron

* Prediction of Deuteron WF from phase shift fit.
* ET Wave functions should match projections of numerical solutions with $\mathrm{Av}_{18}$ solid blue lines

- The matrix elements are continuous in energy across $\mathrm{E}=0$, one can fit $\mathrm{V}_{\delta}$ in the continuum and determine bound states.
* Using the same phase shift data we get
* With pionful $\mathrm{V}_{\text {IR }}=$ OPEP, at N3LO Ebinding $=-2.2278 \mathrm{MeV}$
* With pionless $\mathrm{V}_{\mathrm{IR}}=0$, at N3LO E E binding $=-2.0690 \mathrm{MeV}$


## Continuum Wave Functions

* ET Wave functions (long black dashes) should match projections of numerical solutions with $\mathrm{Av}_{18}$ (dotted colored lines)

* The energies chosen in the plot are deliberately chosen to be distinct from the $\left(\mathrm{E}_{\mathrm{i}}, \delta_{\mathrm{i}}\right)$ used in fitting the LECs.
* Phase shifts are recovered by solving for $\delta$ in

$$
H^{e f f}\left(E_{i}, L E C s, \delta\right) P|\psi\rangle=E_{i} P|\psi\rangle .
$$

## LECs $\rightarrow$ Phase Shifts

* Use fixed LECs at energy E, dial phase shift produce eigenvalue match to E .
* Even NLO 3P1 fit produces a good reproduction of phase shifts.
* A very small number of LECs reproduce phase shifts. P channel NLO has 1, other N3LO have 4.


## Connecting to LQCD

* Lüscher's method can be used to map the spectrum of two nucleons to phase shifts.
* Use traditional path: collect enough phase shift data in multiple channels and use it to fit the HOBET effective interaction.
* This is the first method of connecting QCD to HOBET.
- Sources of error
* Tail of interaction exceeding L/2.

* Divergences of the zeta function in higher order terms of Lüscher's formula.


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## HOBET in Periodic Volumes

* This is the second way to connect QCD to HOBET.
* Phase shifts as boundary conditions are replaced by periodic boundary conditions.
* Easier to construct in Cartesian HO basis.


Slice of 3D Cartesian State

* Key Observation: $\mathrm{V}_{\delta}$ is short range and isolated from the boundary conditions by Green's functions. It is the same object in infinite volume, or periodic volumes.
* We can use Cartesian-spherical brackets to relate $\mathrm{V}_{\delta}$ in both domains. The Cartesian $\mathrm{V}_{\delta}$ can be written in terms of the infinite volume spherical LECs!
* If $\mathrm{V}_{\text {IR }}$ is longer range than $\mathrm{L} / 2$, introduce images of $\mathrm{V}_{\text {IR }}$.
* This is a key advantage over Lüscher's method which requires a free propagation region outside the range of V , but inside the volume.


## Evaluate by Inserting Periodic Basis

Sum T to all orders: $\left\langle\vec{n}^{\prime}\right| \frac{E}{E-T Q}\left[T+T \frac{Q}{E} T\right] \frac{E}{E-Q T} P|\vec{n}\rangle=E\left(\delta_{\vec{n}^{\prime} \bar{n}}-b_{\vec{n} \vec{n}}\right)$

$$
b_{i j}=\left\{P \frac{E}{E-T} P\right\}_{i j}^{-1}
$$

* $\mathrm{V}_{\text {IR }}$ matrix elements are the most expensive part of $\mathrm{H}_{\text {eff }}$

$$
\left\langle\vec{n}^{\prime}\right| G_{T Q} V_{I R} G_{Q T}|\vec{n}\rangle=\sum_{\bar{m}^{\prime}, \vec{m}, \vec{s}, \bar{i}} b_{\vec{n}^{\prime}, s} \frac{E}{E-\lambda_{\vec{m}^{\prime}}}\left\langle\vec{s} \mid \vec{m}^{\prime}\right\rangle\left\langle\vec{m}^{\prime}\right| V_{I R}|\vec{m}\rangle\langle\vec{m} \mid \vec{t}\rangle \frac{E}{E-\lambda_{\vec{m}}} b_{i, \bar{n}}
$$

$\vec{m}, \vec{m}^{\prime}$ are discrete momentum states; s,t are HO states

* All pieces are precomputed, but sum is still very large.
* For $\vec{n}^{\prime}, \vec{n} \in P^{-} \quad G_{Q T}=1$, which can be used to check results.


## Testing Plan



## Induced Mixing

* Setup: spherical well potential in a periodic finite volume.
* The wave function is sampled on sphere outside potential and displayed as a radial displacement from a unit sphere.
* Higher order structure induced by periodic boundary conditions is obvious.
* All this mixing is isolated in E/(E-QT) Green's functions.



## Test Setup: Range(V)>L/2




$$
\begin{aligned}
& L=14.3 \mathrm{fm} \\
& m_{\pi} L=10
\end{aligned}
$$

* Periodic images of the potential make a contribution.
* Continuum extrapolation done on $\mathrm{N}^{\wedge} 3$ lattice with $\mathrm{N}=\{350,400,450\}$.
* Infinite volume bound state at -4.052 MeV.
* LECs are fit using states with $L=0$ overlap.

| Rep | MeV | $\mathrm{L}=0$ | $\mathrm{~L}=2$ | $\mathrm{~L}=4$ | $\mathrm{~L}=6$ |
| :--- | ---: | ---: | ---: | ---: | ---: |
| $A_{1}^{+}$ | -4.4428 | 0.5 | 0 | 0.866 | 0 |
| $A_{1}^{+}$ | 2.0314 | 0.155 | 0 | 0.988 | 0 |
| $E^{+}$ | 7.5995 | 0 | 0.424 | 0.361 | 0.830 |
| $E^{+}$ | 15.2980 | 0 | 0.474 | 0.393 | 0.788 |
| $A_{1}^{+}$ | 21.6167 | 0.326 | 0 | 0.265 | 0.908 |
| $E^{+}$ | 23.2423 | 0 | 0.468 | 0.597 | 0.651 |
| $A_{1}^{+}$ | 29.4041 | 0.521 | 0 | 0.853 | 0.023 |
| $E^{+}$ | 30.9457 | 0 | 0.567 | 0.428 | 0.704 |
| $A_{1}^{+}$ | 35.2449 | 0.655 | 0 | 0.189 | 0.732 |
| $E^{+}$ | 38.4043 | 0 | 0.882 | 0.176 | 0.437 |
| $A_{1}^{+}$ | 45.1402 | 0.526 | 0 | 0.576 | 0.625 |

## Phase Shift Setup

* Reference phase shifts for $\mathrm{L}=0$ and $\mathrm{L}=4$ are directly calculated from V.
* HOBET S-channel phase shifts are computed from the N3LO LECs that reproduce the spectrum. The phase shift is found by dialing the phase shift to produce energy self consistency.
* Lüscher's method phase shifts come from the formula

$$
k \cot \delta_{0}=\frac{2}{\sqrt{\pi} L} \mathcal{Z}_{0,0}\left(1 ; \tilde{k}^{2}\right)+\frac{12288 \pi^{7}}{7 L^{10}} \frac{\mathcal{Z}_{4,0}\left(1 ; \tilde{k}^{2}\right)^{2}}{k^{9} \cot \delta_{4}}+\mathcal{O}\left(\tan ^{2} \delta_{4}\right) \quad \text { Luu, Savage, }
$$

* An effective range expansion up to $\mathrm{k}^{6}$ is used to interpolate.
* For simplicity the second term is evaluated using the $\mathrm{L}=4$ phase shift directly determined from V .


## Phase Shift Results <br> $$
\begin{aligned} & L=14.3 \mathrm{fm} \\ & m_{\pi} L=10 \end{aligned}
$$

The V column should be considered the reference.

|  |  |  | Leading | Next Order <br> E MeV |
| :---: | :---: | :---: | :---: | :---: |
| V | HOBET | Lüscher | Lüscher |  |
| 1 | 142.023 | 141.931 | 142.552 | 142.751 |
| 2 | 128.972 | 128.860 | 129.571 | 129.823 |
| 4 | 113.602 | 113.464 | 114.205 | 114.403 |
| 8 | 96.919 | 96.752 | 97.575 | 97.3135 |
| 10 | 91.473 | 91.296 | 92.228 | 91.6403 |
| 15 | 81.672 | 81.480 | 82.852 | 81.3184 |
| 20 | 74.876 | 74.691 | 76.667 | 74.0936 |

* ET bound state found at -4.066 MeV v.s. -4.052 MeV (directly from V).
* HOBET errors are from PV solution + Momentum basis cutoff.
* Lüscher errors are from Range $(\mathrm{V})>\mathrm{L} / 2$ and magnification of errors by Zeta function poles.


## Effective Operators

* The Bloch Horowitz equation tell us how to renormalize an operator: $\quad \hat{O}_{i i}^{\text {eff. }}(E)=P \frac{E_{j}}{E_{j}-H Q} \hat{o} \frac{E_{i}}{E_{i}-Q H} P$
* $\quad i, j$ Label eigenstates of H .

$$
(P+Q) H\left|\psi_{i}\right\rangle=E_{i}\left|\psi_{i}\right\rangle
$$

* The Green's functions reconstruct $E_{i} P\left|\psi_{i}\right\rangle=\left(E_{i}-Q H\right)\left|\psi_{i}\right\rangle$ the full wave function from the projection.

$$
\left|\psi_{i}\right\rangle=\frac{E_{i}}{E_{i}-Q H} P\left|\psi_{i}\right\rangle
$$

* In bound states the boundary condition for $\mathrm{E} /(\mathrm{E}-\mathrm{QH})$ is an exponential decay outside the range of V .


## Operator Expansion

- Short range operators can also be matched to an expansion.

$$
\begin{aligned}
\hat{O}_{j i}^{\text {eff }} & =P \frac{E_{j}}{E_{j}-H Q} \hat{O} \frac{E_{i}}{E_{i}-Q H} P \\
& =P \frac{E_{j}}{E_{j}-T Q}\left[\hat{O}+V Q \frac{E_{j}}{E_{j}-H Q} \hat{O}+\hat{O} \frac{E_{i}}{E_{i}-Q H} Q V+V Q \frac{E_{j}}{E_{j}-H Q} \hat{O} \frac{E_{i}}{E_{i}-Q H} Q V\right] \frac{E_{i}}{E_{i}-Q T} P \\
& \rightarrow P \frac{E_{j}}{E_{j}-T Q}\left[\hat{O}+\hat{O}_{\delta}\right] \frac{E_{i}}{E_{i}-Q T} P
\end{aligned}
$$

* $\mathrm{O}_{\delta}$ has an expansion much like $\mathrm{V}_{\delta}$ with an expansion in harmonic oscillator quanta.
* Key Point: The LECs $\hat{1}_{\text {of }}$ the expansion can be fit to a set of LQCD measurements. The boundary conditions are then replaced in $\mathrm{E} /(\mathrm{E}-$ QT) with the infinite volume boundary conditions (phase shifts) to give the effective operator in infinite volume.


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& =P \frac{E_{j}}{E_{j}-T Q}\left[\hat{O}+V Q \frac{E_{j}}{E_{j}-H Q} \hat{O}+\hat{O} \frac{E_{i}}{E_{i}-Q H} Q V+V Q \frac{E_{j}}{E_{j}-H Q} \hat{O} \frac{E_{i}}{E_{i}-Q H} Q V\right] \frac{E_{i}}{E_{i}-Q T} P \\
& \rightarrow P \frac{E_{j}}{E_{j}-T Q}\left[\hat{O}+\hat{O}_{i}\right] \frac{E_{i}}{E_{i}-Q T} P
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* $\mathrm{O}_{\delta}$ has an expansion much like $\mathrm{V}_{\delta}$ with an expansion in harmonic oscillator quanta.
- Key Point: The LECs ${ }^{\hat{1}}$ of the expansion can be fit to a set of LQCD measurements. The boundary conditions are then replaced in E/(EQT) with the infinite volume boundary conditions (phase shifts) to give the effective operator in infinite volume.


## Application to $0 v \beta \beta$ Operators

Example: $\quad V_{i}^{n n \rightarrow p p}=-O_{i} \frac{g_{A}^{2}}{4 F_{\pi}^{2}} \tau_{1}^{+} \tau_{2}^{+} \frac{\sigma_{1} \cdot \mathbf{q} \sigma_{2} \cdot \mathbf{q}}{\left(|q|^{2}+m_{\pi}^{2}\right)^{2}}$
Nicholson et al. Phys. Rev. Lett. 121, 172501 (2018)

Running of 1S-1S Matrix Element, $E=-1.961 \mathrm{MeV}$

* Boxed part * $10^{6}$.
* HO Length scale $b=1.7 \mathrm{fm}$
* Start in $\Lambda_{\infty}=80$ and integrate out shell by shell.
* Note jump when 15 becomes an edge state at $\Lambda_{\mathrm{P}}=0$.



## Effective Operators in A-Body Context

* The E in Ôeff is the A-body E.
* Translation invariance requires a total $\Lambda$ cutoff. If spectators are
 excited - red dots, then $\Lambda_{12}$ must be reduced.
* We add a spectator quanta index to the standard density matrix. The interacting particles are then in a 2 particle P space defined by $\Lambda_{12}=\Lambda-\Lambda$ s.
- Matched with this we produce ${ }_{i j}^{e f \cdot \Lambda_{12}}$ for $\Lambda_{12}=0 \ldots \Lambda$.


## Implementation with BIGSTICK

* We (Evan Rule erule@berkeley.edu) are constructing a 2body spectator dependent density matrix for BIGSTICK.
- We will use a realistic potential for H in Ôeff.
* Given universality with respect to the starting potential, we hope for the same with $\hat{O}_{i j}^{\text {eff }, \Lambda_{1}}$.
* We will test with operators associated with experiments.
* Last we will evaluate various $0 \imath \beta \beta$ operators.

$$
\hat{o}=\tau_{1}^{+} \tau_{2}^{+} \frac{\sigma_{1} \cdot \mathbf{q} \sigma_{2} \cdot \mathbf{q}}{\left(|q|^{2}+m_{\pi}^{2}\right)^{2}}
$$

## Summary

* HOBET can be connected to QCD via LQCD observables, or an LQCD nucleon scattering spectrum in finite volume.
* Operators have an expansion, with LECs isolated from boundary conditions by Green's functions and can be fit to LQCD measurements.
* We have made progress on operator renormalization and evaluation in an A-body context. We hope to have results for $0 v \beta \beta$ soon via a hybrid approach with a standard shell model.
* Longer term we are continuing on a path to a HOBET based shell model code.


## End

