Mixing of flavor-singlet light meson, charmonium and gluonic operators with optimal distillation profiles

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Motivation

Why is this mixing important?

- ? Relevant for calculating the low-lying flavor-singlet meson spectrum.
- ? States of theoretical and experimental interest are in this channel, e.g (recent) glueball candidates M. Ablikim *et al.* (BESIII), Phys. Rev. Lett., 132 (2024), exotics, hybrids, etc...
- ? Understanding the mixing dynamics between light meson, charmonium and gluonic operators can give some insights on the composition of these states of interest.

Challenges:

- ! Signal-to-noise problem, particularly in *disconnected* correlations.
- ! How to saturate the spectrum? "Physical"-looking operators, multi-particle operators, etc...
- ! Presence of dynamical quarks degrades the efficacy of gluonic operators: what is a glueball in this setup?



Mixing of flavor singlets

We work with $N_f = 3 + 1 \text{ QCD} \rightarrow \text{SU(3)}$ light flavor symmetry.

- Flavor singlet: f_0 , η_c , glueballs, singlet 2-pion states, etc...
- Flavor octet: π , a_0 , octet 2-pion states, etc...

Can we study charmonia, light mesons and glueballs separately? **No**. All flavor singlets are in the **same** lattice symmetry channel! **Mixing**!

$$\begin{split} \langle \chi_{c0} | \mathcal{O}_{\bar{c}c}^{\dagger} | \Omega \rangle &\neq 0 \rightarrow \text{This is expected!} \\ \langle f_0 | \mathcal{O}_{\bar{c}c}^{\dagger} | \Omega \rangle &\neq 0 \rightarrow \text{This is the issue!} \\ C_{\bar{c}c}(t) \stackrel{t \rightarrow \infty}{\approx} | \langle f_0 | \mathcal{O}_{\bar{c}c}^{\dagger} | \Omega \rangle |^2 e^{-E_{f_0} t} \neq | \langle \chi_{c0} | \mathcal{O}_{\bar{c}c}^{\dagger} | \Omega \rangle |^2 e^{-E_{\chi_{c0}} t} \end{split}$$

Energies obtained from charmonium operators **do not** necessarily correspond to charmonium energy eigenstates.

This mixing can be **artificially** removed to some extent by **ignoring charm annihilation effects** but doing so is **not** strictly correct.

Can we gain some insight on the nature of the different states?



Flavor SU(3) states are labeled by $|D, Y, I, I_3\rangle$ and we are interested in $|1, 0, 0, 0\rangle$:

• Meson interpolators: $\mathcal{O}_c = \bar{c}c$, $\mathcal{O}_l = \frac{1}{\sqrt{3}} \left(\bar{u}u + \bar{d}d + \bar{s}s \right)$

- Gluonic interpolators: $\mathcal{O}_g = \bigcup_{i=1}^{n}$, \mathcal{T}_i , etc...
- 2-pion interpolators: O_{2π} built from
 8 ⊗ 8 = 1 ⊕ 8 ⊕ 8' ⊕ 10 ⊕ 10 ⊕ 27 via SU(3) Clebsch-Gordan
 coefficients. P. McNamee & F. Chillton, Rev. Mod. Phys 36 (1964)

We build a mixing correlation matrix

$$\begin{pmatrix} \langle \mathcal{O}_{l}(t)\bar{\mathcal{O}}_{l}(0) \rangle & \langle \mathcal{O}_{l}(t)\bar{\mathcal{O}}_{c}(0) \rangle & \langle \mathcal{O}_{l}(t)\bar{\mathcal{O}}_{2\pi}(0) \rangle & \langle \mathcal{O}_{l}(t)\bar{\mathcal{O}}_{g}(0) \rangle \\ * & \langle \mathcal{O}_{c}(t)\bar{\mathcal{O}}_{c}(0) \rangle & \langle \mathcal{O}_{c}(t)\bar{\mathcal{O}}_{2\pi}(0) \rangle & \langle \mathcal{O}_{c}(t)\bar{\mathcal{O}}_{g}(0) \rangle \\ * & * & \langle \mathcal{O}_{2\pi}(t)\bar{\mathcal{O}}_{2\pi}(0) \rangle & \langle \mathcal{O}_{2\pi}(t)\bar{\mathcal{O}}_{g}(0) \rangle \\ * & * & * & \langle \mathcal{O}_{g}(t)\bar{\mathcal{O}}_{g}(0) \rangle \end{pmatrix} \end{cases}$$

and from the GEVP we obtain an optimal interpolator

 $\tilde{\mathcal{O}}^{(n)} = \mathbf{a_1^n}\mathcal{O}_c + \mathbf{a_2^n}\mathcal{O}_l + \mathbf{a_3^n}\mathcal{O}_q + \mathbf{a_4^n}\mathcal{O}_{2\pi}, \ a_i^n \in \mathbb{R}$

which maximizes $\langle n | \hat{\tilde{\mathcal{O}}}^{(n)\dagger} | \Omega \rangle$.



	\mathcal{O}_l	\mathcal{O}_{c}	$O_{2\pi}$	C	\mathcal{D}_{g}
\mathcal{O}_l		\$		\$	⊊ €¢
\mathcal{O}_c	-		00	0	S €S€°
$\mathcal{O}_{2\pi}$	-	-		0	⊊ &
\mathcal{O}_{g}	-	-	-	~ <u>&</u>	Se so

Disconnected correlations are **necessary** for flavor-singlets but have **large** statistical errors.

- **Explicit mixing:** including flavor-mixing off-diagonals.
- Implicit mixing: including disconnected contributions in diagonals.

Optimal distillation profiles

Distillation M. Peardon et al. Phys. Rev. D 80, 054506 (2009)

- Project quark fields onto low-dimensional subspace of smooth, gauge-covariant fields → Smearing.
- ► $\psi(t) \rightarrow V[t]V[t]^{\dagger}\psi(t)$ with V[t] the low-modes of the 3D gauge-covariant Laplacian operator.
- Perambulators: $\tau[t_1, t_2] = V[t_1]^{\dagger} D^{-1} V[t_2]$
- Elementals: $\Phi[t] = V[t]^{\dagger} \Gamma V[t]$, $\Gamma = \gamma_5, \gamma_i, \nabla_i, \dots$

Example two-point meson correlation:

$$\begin{split} C(t) &= -\left\langle \mathrm{Tr}\left(\Phi[t]\tau[t,0]\bar{\Phi}[0]\tau[0,t]\right)\right\rangle_{\mathrm{gauge}} \\ &+ \left\langle \mathrm{Tr}\left(\Phi[t]\tau[t,t]\right)\mathrm{Tr}\left(\bar{\Phi}[0]\tau[0,0]\right)\right\rangle_{\mathrm{gauge}} \end{split}$$

High inversion cost but matrices have manageable sizes and can be reused for different choices of correlations, e.g one-particle, two-particle, etc...

Improved Distillation J. A. Urrea-Niño, F. Knechtli, T. Korzec & M. Peardon. Phys. Rev. D 106. 034501 (2022)

- ► Exploit further freedom: $V[t]V[t]^{\dagger} \rightarrow V[t]J[t]V[t]^{\dagger}$ with quark distillation profile $J[t]_{ij} = \delta_{ij}g(\lambda_i[t])$.
- Build a GEVP using different quark profiles:

$$\begin{split} \psi_k(t) &= V[t] J_k[t] V[t]^{\dagger} \psi(t) \rightarrow \text{ Different quark profiles...} \\ \mathcal{O}_k(t) &= \bar{\psi}_k(t) \Gamma \psi_k(t) \rightarrow \dots \text{define different meson operators...} \\ \Phi^k[t] &= J_k[t]^{\dagger} \Phi[t] J_k[t] \rightarrow \dots \text{ with different elementals...} \\ C_{ab}(t) &= \left\langle \mathcal{O}_a(t) \bar{\mathcal{O}}_b(0) \right\rangle \rightarrow \dots \text{ for a correlation matrix.} \end{split}$$

E.g optimal meson profile

$$f^{(\Gamma,n)}\left(\lambda_{i}[t],\lambda_{j}[t]\right) = \sum_{k} a_{k}^{(\Gamma,n)} g_{k}\left(\lambda_{i}[t]\right)^{*} g_{k}\left(\lambda_{j}[t]\right)$$

- $\checkmark~$ No additional inversion cost for improvement.
- \checkmark One optimal profile for each Γ and energy level n.
- ✓ Profiles introduced at contraction level, which can be made efficient via BLAS, Numpy, etc...



Standard vs **Improved** distillation in connected-only charmonium correlations.





Gauge ensembles

Wilson fermion action with non-perturbatively determined clover improvement + Lüscher-Weisz gauge action. R. Höllwieser et al. Eur. Phys. J. C 80, 349, P. Fritzsch et al. J. High Energ. Phys. 2018, 25 (2018)



Our setup is very convenient:

! Quenched lattice QCD predicts a 0^{++} glueball at \approx 1800 MeV. c.

Morningstar and M. Peardon, Phys. Rev. D 60, 034509

We can restrict pion decays in each ensemble:

- A1: Glueball $\rightarrow \pi\pi$, $\pi\pi\pi\pi$
- A1h: Glueball $\rightarrow \pi\pi$



Hadron creation operators

We have to saturate the spectrum as best as possible.

One-meson operators: $\bar{c} \mathbb{I} c$, $\frac{1}{\sqrt{3}} \sum_{q=u,d,s} \bar{q} \mathbb{I} q$

- Build correlation matrix with 7 Gaussian profiles.
- Prune via SVD to 3(5) charmonium(light meson) operators. J. Balog et al., Phys. Rev. D 60, 094508, F. Niedermayer et al., Nuclear Physics B 597, 413–450
- $\checkmark\,$ Pruned operators are almost "aligned" with energy eigenstates.

Gluonic operators: Sum of Laplacian eigenvalues. C. Morningstar et al. Phys. Rev. D 88, 014511

 \checkmark Consistent with 3D Wilson loops but slightly cleaner signal.

Two-pion operators: π ($\vec{p} = 0$) π ($-\vec{p} = 0$) with standard distillation

We account for most expected states but improvement is possible...

What will I show?

- Finite volume energy spectrum of the scalar channel taking into account mixing between mesonic and gluonic operators.
- Effect of including different types of operators on the spectrum.
- Overlaps between energy eigenstates and the different types of operators.

The overlaps are calculated as

 $\langle n | \hat{\mathcal{O}}_{i}^{\dagger} | \Omega \rangle \propto [C(t_{0})\vec{\omega}_{n}(t,t_{0})]_{i}$

to see which type of operator contributes most to each state. J.J. Dudek et al. Phys. Rev. D 77, 034501 (2008)

We obtain a finite volume spectrum and a scattering study requires the Lüscher formalism. M. Lüscher, Nucl. Phys. B 354, 531 (1991)

0^{++} flavor-singlet correlation matrix at t = a

$C_{ij}(t) \to C_{ij}(t)/\sqrt{C_{ii}(a)C_{jj}(a)}.$







Figure: A1h ensemble

- Charmonium, light meson and gluonic operators have non-zero correlations: mixing!
- 2-pion operator mixes very little: creates a mostly multi-particle state.

Single-particle operator mixing in A1 $(m_{\pi} \approx 420 \text{ MeV})$



- Charmonium operators alone see a light state: charm annihilation (disconnected) effects!
- ► Including *O*_g does not change the low-lying spectrum. Similar to R. Brett et al. 1909.07306.

Full operator mixing in A1 $(m_{\pi} \approx 420 \text{ MeV})$



▶ 10 operators: 3 × O_c, 5 × O_l, O_{2π} and O_g.
 ▶ 2-pion operator introduces an additional state.



Single-particle operator mixing in A1h ($m_{\pi} \approx 800 \text{ MeV}$)



Charm annihilation effects are again important!

• Including \mathcal{O}_g does not change the low-lying spectrum.

Full operator mixing in A1h ($m_{\pi} \approx 800 \text{ MeV}$)



▶ 10 operators: 3 × O_c, 5 × O_l, O_{2π} and O_g.
 ▶ 2-pion operator introduces an additional state.



What have we learned from the spectrum?

- ✓ Including flavor-mixing of operators is fundamental to map out the energy spectrum.
- ✓ Including multi-particle operators is necessary to avoid **missing** existing states.
- ✓ Including purely gluonic operators does **not** introduce a new state.

What have we learned from the overlaps?

- \checkmark All flavor-singlet operators "talk" to each other to some degree.
- ✓ Charmonium operators "see" light states and viceversa: charm annihilation effects are important!
- ✓ 2-pion operators overlap strongly with a state somewhat separate from the 1-particle ones.



Where can we improve?

Tackle the signal-to-noise problem for *disconnected*-like correlations, e.g gluonic operators or meson loops.



Multi-level sampling was recently systematically studied for 3D Wilson loops in SU(3) pure gauge theory. L. Barca *et. al.* Phys. Rev. D 110, 054515 (2024)

Application to case with dynamical quarks + distillation is currently under study. Talk by L. Barca at Lattice 2024

This *algorithmic* improvement happens at the level of generating the gauge configurations.

Improve our operator basis.

One-meson operators:

- Include multiple Γ with profiles.
- ✓ Derivative-based operators can better sample spatial structure or additional gluonic content.
- ✓ E.g $\mathbb{B}_i = \epsilon_{ijk} \nabla_j \nabla_k$ is very sensitive to gluonic background.

Two-pion operators:

- ▶ Include non-zero back-to-back momentum in $\pi(\vec{p}) \pi(-\vec{p})$.
- Include distillation profiles to each pion.
- ✓ Profiles have been shown to help also with one-meson operators at non-zero momentum.

These *operator* improvements happen at contraction time.

FOR5269: Future methods for studying confined gluons in QCD

Spokesperson: Prof. Dr. Francesco Knechtli

Collaboration between **physics** and **applied mathematics** at BUW, DESY Zeuthen and Trinity College Dublin.

Some main pillars:

- ► Contributions from charm-annihilation effects.
- Glueballs and their mixing in dynamical QCD.
- Improved numerical methods, e.g for distillation.
- Multi-level sampling with/without dynamical quarks.
- String breaking in hybrid potentials.
- New schemes for molecular dynamics.

We are tackling *physical, mathematical and numerical* problems particularly relevant for hadron spectroscopy.



Thank you for your attention!



There are computationally intensive steps worth optimizing.

Eigenvector calculation:

- ► N_v lowest eigenvectors of $\nabla^2[t]$ are calculated for every t.
- ✓ Thick-Restart Lanczos with periodic reorthogonalization and Chebyshev acceleration.
- ✓ C + MPI + LAPACK code written in the *qcdlib* library of T. Korzec.

Perambulator calculation:

- ► $\tau[t_1, t_2]$ requires solving $4 \times N_v \times N_t$ linear systems Dx = b.
- $\checkmark\,$ C + MPI code in *qcdlib* which calls the openQCD linear solver.

Matrix contractions:

- $\tau[t_1, t_2]$ and $\Phi[t]$ have size $4N_v \times 4N_v$.
- \checkmark Numpy is flexible and efficient for such "small" matrices.
- ✓ Sequences Tr $(\Phi[t_1]\tau[t_1, t_2]\Phi[t_2]\tau[t_2, t_3]...\Phi[t_M]\tau[t_M, t_1])$ can be efficiently calculated, e.g via Numpy's *einsum* + mpi4py.

Example profiles for connected-only charmonium



- Suppression of higher modes for ground states.
- Non-constant profiles.

- Nodes in excited states.
- Non-trivial structure.

