

Controlling Excited-State Contributions to Nucleon Isovector Charges using Distillation

Colin P. Egerer

In collaboration with Dr. David Richards & Dr. Frank Winter

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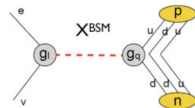
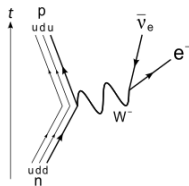


WILLIAM
& MARY



Why Calculate Nucleon Isovector Charges?

- $n \rightarrow pe^{-}\bar{\nu}_e$ coupling & proton-proton fusion
- $g_A^{u-d} = \langle 1 | \Delta u - \Delta d$
- $g_T^{u-d} = \langle 1 | \delta u - \delta d$
- g_S^{u-d} as related to DM direct detection
- novel BSM interactions at TeV scale



[LANL: T-2 Nuclear Physics]

Ab-initio QCD

- Despite actual QCD on Lattice ... inherent systematics

- ① finite lattice spacing a effects

[J. Liang, et al., arXiv:hep-lat/1612.04388
Y.-B. Yang, et al., Phys. Rev. **D93**, 034503 (2016), arXiv:hep-lat/1509.

- ② volume dependence V

[G. S. Bali, et al., Phys. Rev. **D91**, 054501 (2015).]
[R. Horsley, et al., Phys. Lett. **B732** 41 (2014).]

- ③ pion mass m_π

[S. Ohta (LHP, RBC, UKQCD collabs.), arXiv:/hep-lat/1511.05126]

- ④ number of quark flavors

[H.-W. Lin, et al., arXiv:hep-lat/1806.10604].

- ⑤ chiral symmetry

[R. Gupta, et al., arXiv:hep-lat/1806.09006].

[R. Gupta, et al., arXiv:hep-lat/1801.03130]

- ⑥ signal-to-noise

[C. C. Chang, et al., Nature **558**, 91-94 (2018).]

- ⑦ excited states

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Excited-State Contamination

$$C_{2\text{pt}}(t) = \sum_{\vec{x}} \langle \mathcal{O}(\vec{x}, t) \overline{\mathcal{O}}(\vec{0}, 0) \rangle = \sum_n \frac{1}{2M_n} |\langle 0 | \mathcal{O} | n \rangle|^2 e^{-M_n t}$$

- Interpolating fields of point-like quark/gluon fields
 - hadronic states at all energy scales
- For ground-state quantities
 - large-Euclidean time
 - judicial choice of $\mathcal{O} \rightarrow \langle 0 | \mathcal{O} | n \rangle|_{n>0} \ll \langle 0 | \mathcal{O} | n = 0 \rangle$

Spatial Smearing

$$\tilde{\psi}(\vec{x}, t) = \sum_{\vec{y}} S[U](\vec{x}, \vec{y}) \psi(\vec{y}, t)$$

- Jacobi smearing

[C. Allton et al. (UKQCD Collaboration), Phys. Rev. D47, 5128 (1993).]

$$S[U](\vec{x}, \vec{y}) = J_{\sigma, n_{\sigma}}(t) \equiv \left(1 + \frac{\sigma \nabla^2(t)}{n_{\sigma}} \right)^{n_{\sigma}}$$
$$\longrightarrow \text{Exp}(\sigma \nabla^2(t))$$

- exp. suppression of high eigenmodes of Laplacian
 - lowest modes contribute appreciably

$$\tilde{\psi}(\vec{x}, t) = J_{\sigma, n_{\sigma}}(t) \psi(\vec{x}, t)$$

Distillation

- solutions to $-\nabla^2 \xi^{(k)} = \lambda^{(k)} \xi^{(k)}$ - ordered by $\lambda^{(k)}$

$$-\nabla_{ab}^2 (\vec{x}, \vec{y}; t) = 6\delta_{xy}\delta_{ab} - \sum_{j=1}^3 \left[\tilde{U}_j(\vec{x}, t)_{ab} \delta_{x+\hat{j},y} + \tilde{U}_j^\dagger(\vec{x} - \hat{j}, t)_{ab} \delta_{x-\hat{j},y} \right]$$

- define *Distillation* op. of rank $N \ll M = N_c \times N_X \times N_Y \times N_Z$

$$\square (\vec{x}, \vec{y}; t)_{ab} = \sum_{k=1}^N \xi_a^{(k)}(\vec{x}, t) \xi_b^{(k)\dagger}(\vec{y}, t)$$

[M. Peardon, et. al., (2009) arXiv: 0905.2160v1]

Distillation - Properties & Advantages

- interpolator construction separable from quark propagation
- can construct interpolators to probe angular structure of states w/o recalculating $M^{-1}(t, t')$
- momentum projection at source & sink

$$C(t', t) = \sum_{\vec{x}, \vec{y}} e^{-i\vec{p}\cdot\vec{x}} e^{i\vec{p}\cdot\vec{y}} \langle \mathcal{O}(\vec{x}, t') \overline{\mathcal{O}}(\vec{y}, t) \rangle$$

- number of eigenvectors scales with V_3
 - baryon contractions $\mathcal{O}(V_3^4)$

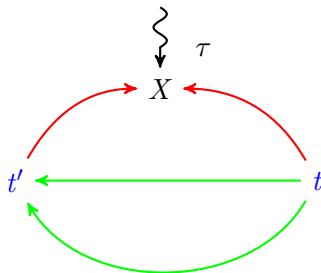
Components of a Calculation using Distillation

- **Solution Vectors**

$$S^k(\vec{x}, t'; t) = M_f^{-1}(t', t) \xi^{(k)}(t)$$

- **Perambulators**

$$\tau_{\alpha\beta}^{kl}(t', t) = \xi^{(k)\dagger}(t') M_{\alpha\beta}^{-1}(t', t) \xi^{(l)}(t)$$



- **Elementals** - operator insertions in Distillation space

$$\Phi_{\alpha_1, \alpha_2, \alpha_3}^{(i, j, k)}(t) = \epsilon^{abc} (\mathcal{D}_1 \xi^{(i)})^a (\mathcal{D}_2 \xi^{(j)})^b (\mathcal{D}_3 \xi^{(k)})^c(t) S_{\alpha_1, \alpha_2, \alpha_3}$$

- Correlators comprised of perambulators and elementals

Variational Analysis

- $\mathcal{B} = \{\mathcal{O}_1, \dots, \mathcal{O}_N\}$
- optimal linear combination of interpolators to project onto $|n\rangle$, solve

$$C(t) u_n(t, t_0) = \lambda_n(t, t_0) C(t_0) u_n(t, t_0)$$

- solved for fixed reference time t_0 and all later times t
- solutions organized by magnitude of eigenvalues $\lambda_n(t, t_0)$ - principal correlators $\lambda_n(t, t_0) \underset{t \gg t_0}{\sim} e^{-m_n(t-t_0)}$
- components of eigenvectors $u_n(t, t_0)$ yield weight of each $\mathcal{O}_i \in \mathcal{B}$ to interpolate $|n\rangle$ from vacuum
- projected operator $\equiv \mathcal{O}_{\text{proj}}^{n\dagger} = \sum_i u_n^i \mathcal{O}_i^\dagger$

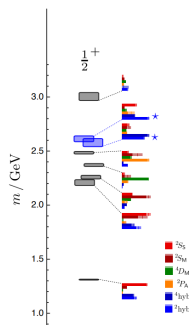
[B. J. Owen, et al., Phys. Lett. **B723**, 217 (2013).] $\sim 8\%$ improvement

Operator Construction

$$\mathcal{O}_i(t) \propto \epsilon^{abc} \mathcal{S}_i^{\alpha\beta\gamma} (\mathcal{D}_1 \square d)_a^\alpha (\mathcal{D}_2 \square u)_b^\beta (\mathcal{D}_3 \square d)_c^\gamma(t)$$

- $\left(N_M \otimes \left(\frac{1}{2}^+ \right)_M^1 \otimes D_{L=0,S}^{[0]} \right)^{J^P = \frac{1}{2}^+}$
- $\left(N_M \otimes \left(\frac{1}{2}^+ \right)_M^1 \otimes D_{L=0,M}^{[2]} \right)^{J^P = \frac{1}{2}^+}$
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- $\left(N_M \otimes \left(\frac{1}{2}^+ \right)_M^1 \otimes D_{L=1,A}^{[2]} \right)^{J^P = \frac{1}{2}^+}$
- $\left(N_M \otimes \left(\frac{1}{2}^+ \right)_M^1 \otimes D_{L=1,M}^{[2]} \right)^{J^P = \frac{1}{2}^+}$
- $\left(N_M \otimes \left(\frac{3}{2}^+ \right)_S^1 \otimes D_{L=1,M}^{[2]} \right)^{J^P = \frac{1}{2}^+}$
- $\left(N_M \otimes \left(\frac{3}{2}^+ \right)_S^1 \otimes D_{L=2,M}^{[2]} \right)^{J^P = \frac{1}{2}^+}$

$$N^{2S+1} L_P J^P$$



N spectrum ordered by J^P for $m_\pi \sim 520$ MeV
 [J. J. Dudek and R. G. Edwards (2012) arXiv:hep-ph/1201.2349]
 [R. Edwards, et. al., (2011) arXiv: 1104.5152v2]

Lattice & Operator Specifics

- $32^3 \times 64$ isoclover lattices ($\beta = 6.3$) - 350cfigs
- $u/d + s$ flavor QCD $m_s \rightsquigarrow (2M_{K^+}^2 - M_{\pi^0}^2) / M_{\Omega^-}^2 = 0.1678$
- tree-level tadpole-improved Symanzik gauge action
- $m_\pi = 356$ MeV
- $a = 0.098$ fm
- Distilled operators
 - 64 eigenvectors
- **Our Aim:** rather than precise determination of g_A, g_S, g_T , can Distillation demonstrate controlled excited-states?

First application of Distillation to structure calculations

2pt Decomposition

- Direct comparisons between 'local' ($N^2 S_S \frac{1}{2}^+$) and 'proj' ($\mathcal{O}_{\text{proj}}^{\text{n}\dagger} = \sum_i u_n^i \mathcal{O}_i^\dagger$)

$$C_{2\text{pt}}(\vec{p}, t) = \sum_{\vec{x}} e^{-i\vec{p}\cdot\vec{x}} \sum_{\vec{y}} e^{i\vec{p}\cdot\vec{y}} \langle \mathcal{O}(\vec{x}, t) \mathcal{O}^\dagger(\vec{y}, 0) \rangle$$

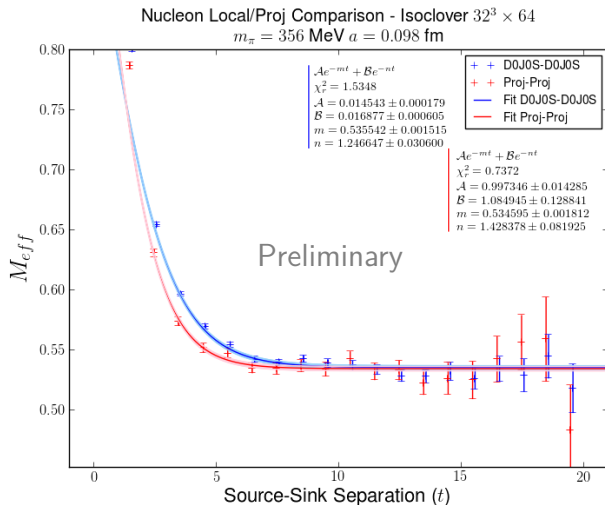
- For zero-momentum states, keeping ground and first-excited states:

$$C_{2\text{pt}}(t) = V_3 \left(\frac{|Z_0|^2}{2m_0} e^{-m_0 t} + \frac{|Z_1|^2}{2m_1} e^{-m_1 t} \right)$$

- Simple 2-state fit for masses/overlap factors

$$C_{2\text{pt}}(t) = \mathcal{A} e^{-m_0 t} + \mathcal{B} e^{-m_1 t}$$

Effective Mass Comparison - Two State



3pt Decomposition

$$C_{3\text{pt}}(\vec{p}', t; \vec{q}, \tau; t_0) = V_3 \sum_{\vec{x}} e^{-i\vec{p}' \cdot \vec{x}} \sum_{\vec{z}} e^{i\vec{q} \cdot \vec{z}} \langle 0 | \mathcal{O}(\vec{x}, t) \mathcal{J}(\vec{z}, \tau) \overline{\mathcal{O}}(t_0) | 0 \rangle$$

- For zero-momentum states at fixed t_{sep}

$$C_{3\text{pt}}(\tau) = \left(\frac{V_3 |Z_0|^2}{4m_0^2} \mathcal{J}_{00} e^{-m_0 t_{\text{sep}}} + \frac{V_3 |Z_1|^2}{4m_1^2} \mathcal{J}_{11} e^{-m_1 t_{\text{sep}}} \right) \begin{array}{c} \text{---} \tau \\ \downarrow \\ X \\ \leftarrow t \quad \rightarrow t_0 \end{array}$$

$$+ \frac{V_3 Z_0 Z_1^\dagger}{4m_0 m_1} \mathcal{J}_{01} e^{-m_0 t_{\text{sep}}} e^{-(m_1 - m_0)\tau} + \frac{V_3 Z_1 Z_0^\dagger}{4m_0 m_1} \mathcal{J}_{10} e^{-m_1 t_{\text{sep}}} e^{(m_1 - m_0)\tau}$$

- Fit form: $C_{3\text{pt}}(\tau) = \mathcal{C} + \mathcal{D} \cosh[\Delta m (\tau - t_{\text{sep}}/2)]$

Extraction of Ground-state Matrix Element

$$\mathcal{C} = \frac{V_3 |Z_0|^2}{4m_0^2} \mathcal{J}_{00} e^{-m_0 t_{\text{sep}}} + \frac{V_3 |Z_1|^2}{4m_1^2} \mathcal{J}_{11} e^{-m_1 t_{\text{sep}}}$$

- Fit t_{sep} -dependence of \mathcal{C} to extract \mathcal{J}_{00}

$$\begin{aligned}\mathcal{C}(t_{\text{sep}}) &= X e^{-m_0 t_{\text{sep}}} + Y e^{-m_1 t_{\text{sep}}} \\ &\Rightarrow \mathcal{J}_{00} \propto X\end{aligned}$$

- “effective charges”

$$g_{\Gamma}^{\text{eff}}(t_{\text{sep}}, \tau) = \frac{C_{\Gamma}^{\text{3pt}}(t_{\text{sep}}, \tau)}{C_{\text{fit}}^{\text{2pt}}(t_{\text{sep}})}$$

Extraction of Ground-state Matrix Element

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- Fit t_{sep} -dependence of \mathcal{C} to extract \mathcal{J}_{00}

$$\mathcal{C}(t_{\text{sep}}) = X e^{-m_0 t_{\text{sep}}} + Y e^{-m_1 t_{\text{sep}}}$$

$$\Rightarrow \mathcal{J}_{00} \propto X$$

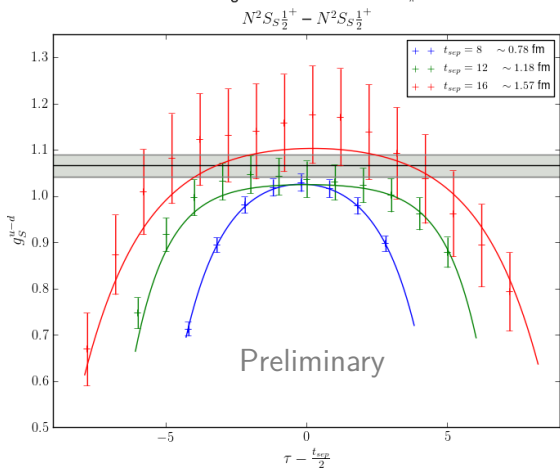
little effect in extracting \mathcal{J}_{00} ...

- “effective charges”

$$g_{\Gamma}^{\text{eff}}(t_{\text{sep}}, \tau) = \frac{C_{\Gamma}^{\text{3pt}}(t_{\text{sep}}, \tau)}{C_{\text{fit}}^{\text{2pt}}(t_{\text{sep}})}$$

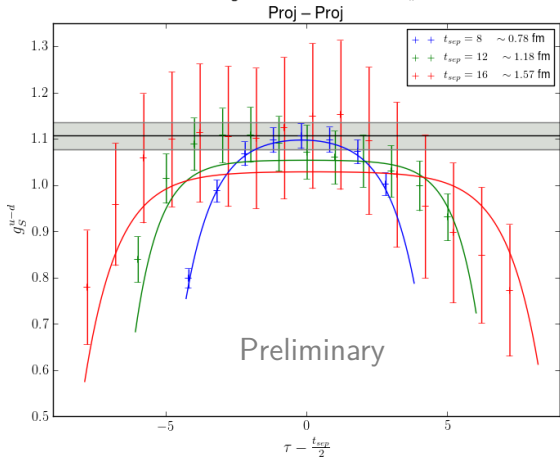
Determination of g_S - Local Operator

Nucleon effective scalar Charge - Isoclover $32^3 \times 64 m_\pi = 356 \text{ MeV } a = 0.098 \text{ fm}$

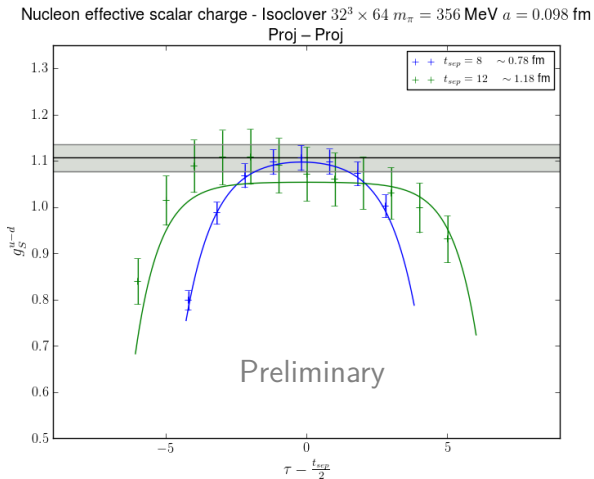


Determination of g_S - Projected Operator

Nucleon effective scalar Charge - Isoclover $32^3 \times 64 m_\pi = 356 \text{ MeV } a = 0.098 \text{ fm}$



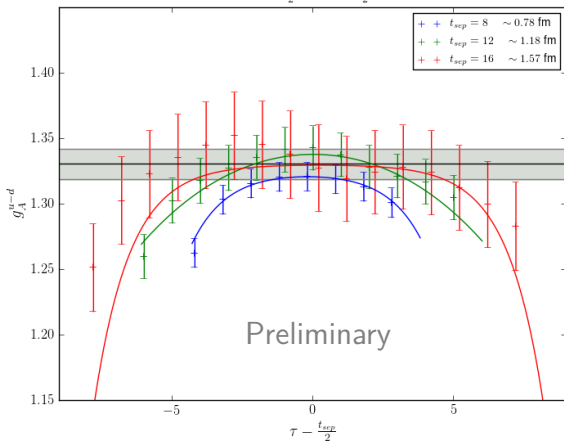
Determination of g_S - Projected Operator (w/o $t_{\text{sep}} = 16$)



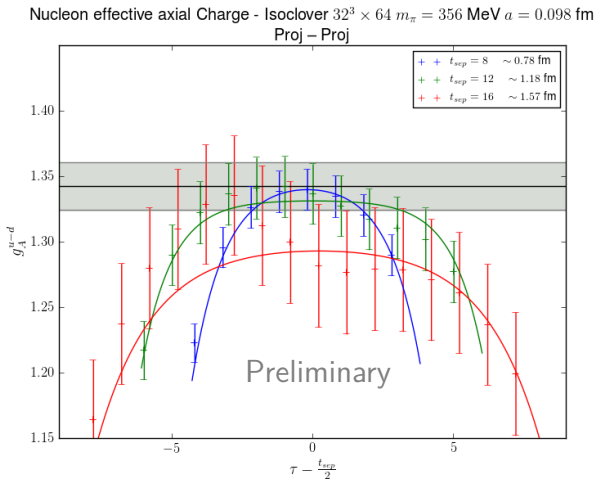
Determination of g_A - Local Operator

Nucleon effective axial Charge - Isoclover $32^3 \times 64$ $m_\pi = 356$ MeV $a = 0.098$ fm

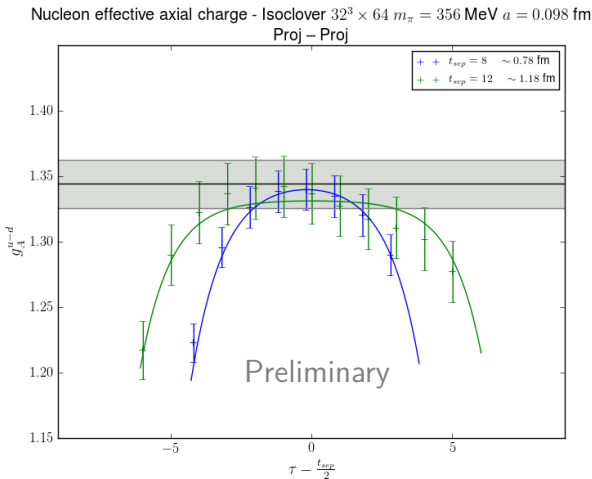
$$N^2S_{S\frac{1}{2}}^{1+} - N^2S_{S\frac{1}{2}}^{1+}$$



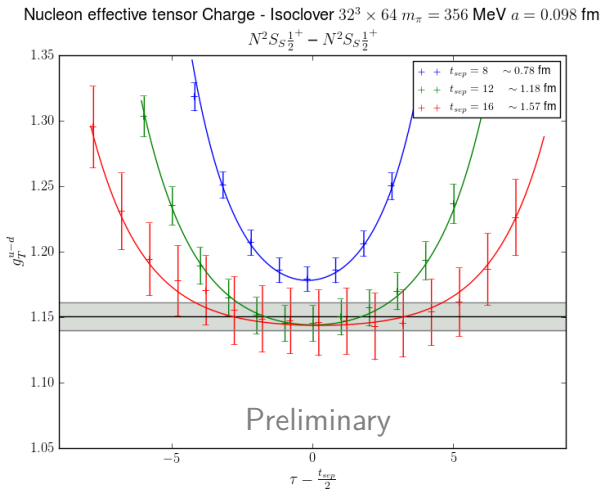
Determination of g_A - Projected Operator



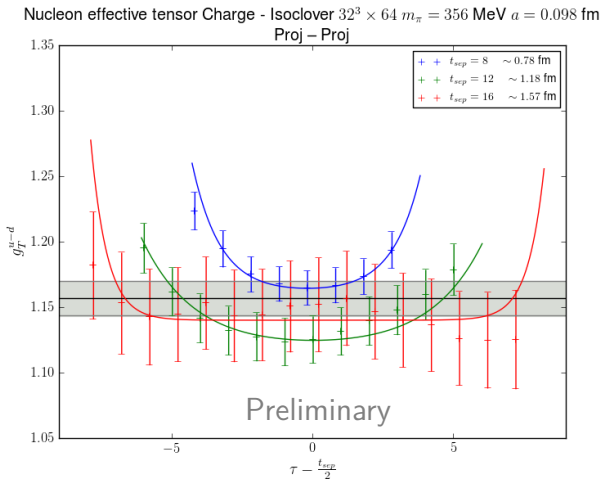
Determination of g_A - Projected Operator (w/o $t_{\text{sep}} = 16$)



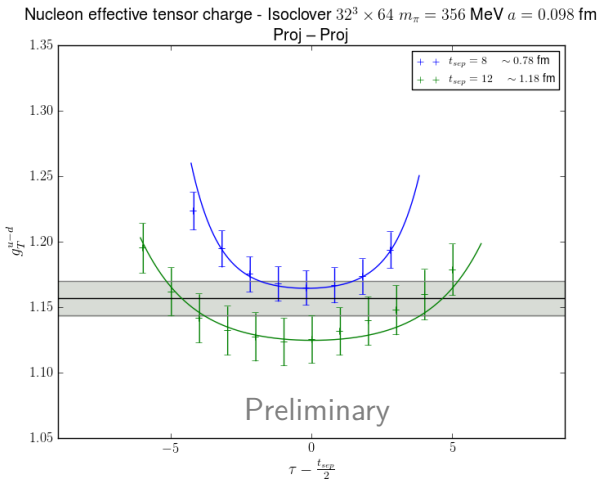
Determination of g_T - Local Operator



Determination of g_T - Projected Operator



Determination of g_T - Projected Operator (w/o $t_{\text{sep}} = 16$)



Closing Thoughts

- Success of Distillation in spectroscopy begs study when considering structure
- Variational method applied to an extended basis of Distilled interpolators
 - effective mass plateaus for much smaller source-sink separations $\sim 0.6 \text{ fm}$ \rightarrow better statistics
 - clear separation of ground & excited states - c.f. projected 2-state fit
- $C_{3\text{pt}}(t_{\text{sep}}, \tau) / C_{2\text{pt}}(t_{\text{sep}})$ plateaus
 - little src-snk dependence when using variationally optimized nucleon interpolators ****noisy $t_{\text{sep}}=16$**
- Considered forward-scattering of at rest nucleons
 - $\vec{p} \neq \vec{0}$ states - address, say, g_{A_i} and g_{A_4} [J. Liang, et al., arXiv/1612.04388.]
- Q^2 axial/vector form factors? Other structure calculations?

BACKUPS

Principal Correlators - $t_0 = 2, 3$

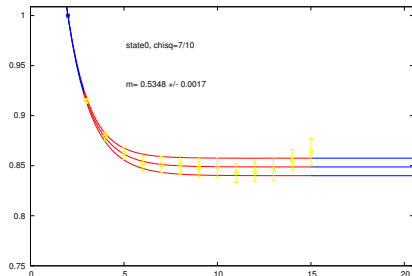


Figure 2: Ground-state principal correlator for $t_0 = 2$ and $t_Z = 5$.

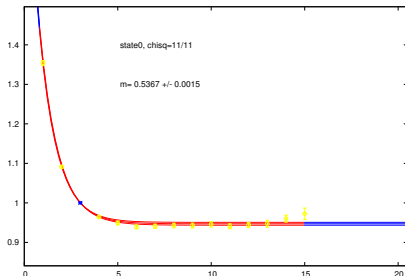


Figure 3: Ground-state principal correlator for $t_0 = 3$ and $t_Z = 5$.

Method for renormalized g_Γ Extraction

- Relation between continuum and lattice charges

$$g_\Gamma = Z_\Gamma g_\Gamma^{\text{lat}}$$

- Z_V set based on knowledge of continuum vector charge
 - i.e. $1 + \mathcal{O}(a^2) = Z_V g_V^{\text{lat}} \implies Z_V = 1/g_V^{\text{lat}} + \mathcal{O}(a^2)$
- Utilized definition of arbitrary renormalized charge

$$g_\Gamma = \frac{Z_\Gamma}{Z_V} \frac{g_\Gamma^{\text{lat}}}{g_V^{\text{lat}}}$$

- Z_Γ/Z_V on isoclover lattices with $a = 0.094$ fm and $m_\pi = 280$ MeV
 - n.b. Z_Γ appears to be weakly dependent on pion mass, but depends on lattice spacing

[B. Yoon, et. al., (2017) arXiv: 1611.07452v3]