

η and η' mesons from $N_f=2+1$ lattice QCD at physical point using topological charge operator

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Quark model: nandn'

For exact SU(3) flavor symmetry one expects:

Flavor octet state:
$$|\eta_8\rangle = \frac{1}{\sqrt{6}}(|\bar{u}u\rangle + |\bar{d}d\rangle - 2|\bar{s}s\rangle)$$

Flavor singlet state:
$$|\eta_1\rangle = \frac{1}{\sqrt{3}}(|\bar{u}u\rangle + |\bar{d}d\rangle + |\bar{s}s\rangle)$$

However, SU(3) flavor symmetry is broken by large $m_s \gg m_u \approx m_d \equiv m_l$.

Physical η , η' states are not flavor eigenstates but mixtures,

$$\begin{pmatrix} |\eta\rangle \\ |\eta'\rangle \end{pmatrix} = M \begin{pmatrix} |\eta_8\rangle \\ |\eta_1\rangle \end{pmatrix}$$

M is the mixing matrix of η and η' .

η and η' on the lattice

 \clubsuit Generally, η and η' are calculated using fermion fields on the lattice.

$$\eta(x) = \sum_{i}^{N_f} \bar{\psi}_i(x) \gamma_5 \psi_i(x) \qquad C_{\vec{p}}^{\eta}(t, t_0) = \langle \tilde{\eta}(t, \vec{p}) (\tilde{\eta}(t_0, \vec{p}))^{\dagger} \rangle$$

$$C_{\vec{p}}^{\eta}(t, t_0) = \frac{1}{V_3} \sum_{\vec{x}, \vec{y}} e^{i\vec{p}(\vec{y} - \vec{x})} \langle \mathcal{Q}_{\beta, b, t_0, \vec{y}; \alpha, a, t, \vec{x}}^{-1} \mathcal{Q}_{\alpha, a, t, \vec{x}; \beta, b, t_0, \vec{y}}^{-1} - \mathcal{Q}_{\beta, b, t_0, \vec{y}; \alpha, b, t_0, \vec{y}}^{-1} \mathcal{Q}_{\alpha, a, t, \vec{x}; \alpha, a, t, \vec{x}}^{-1} \rangle$$

$$t, \vec{x} \qquad t_0, \vec{y} \qquad t, \vec{x} \qquad t_0, \vec{y}$$

Disconnected & Require high statistics

Using the topological charge operator avoids the difficulties of handling disconnected insertions.

$$\eta I^{G}(J^{PC}) = 0^{+}(0^{-+})$$

$$F_{\mu\nu}\tilde{F}^{\mu\nu} \propto \vec{E} \cdot \vec{B}$$

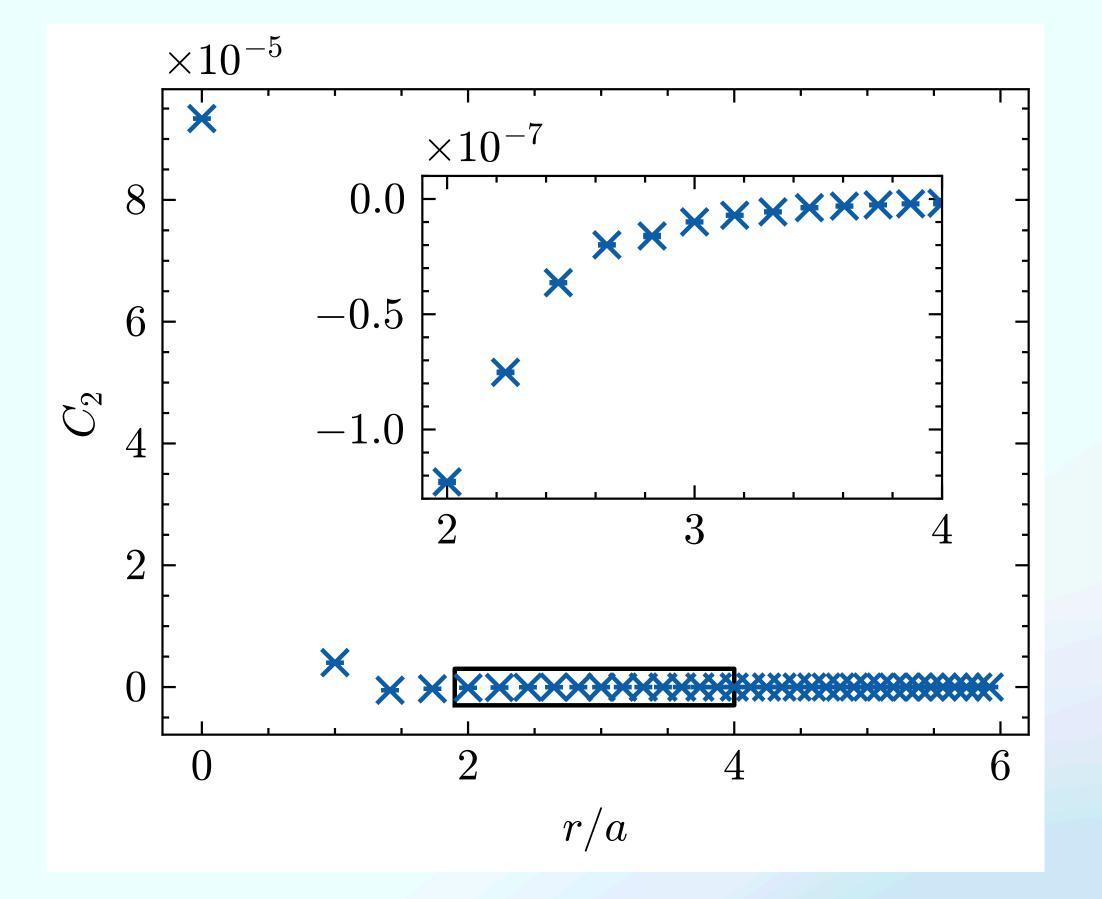
$$q(x) = \frac{1}{32\pi^{2}} \epsilon_{\mu\nu\rho\sigma} \text{Tr} F^{\mu\nu}(x) F^{\rho\sigma}(x)$$

$$\eta' I^{G}(J^{PC}) = 0^{+}(0^{-+})$$

η and η' on the lattice

Constructing two-point correlation functions using topological charge density operators

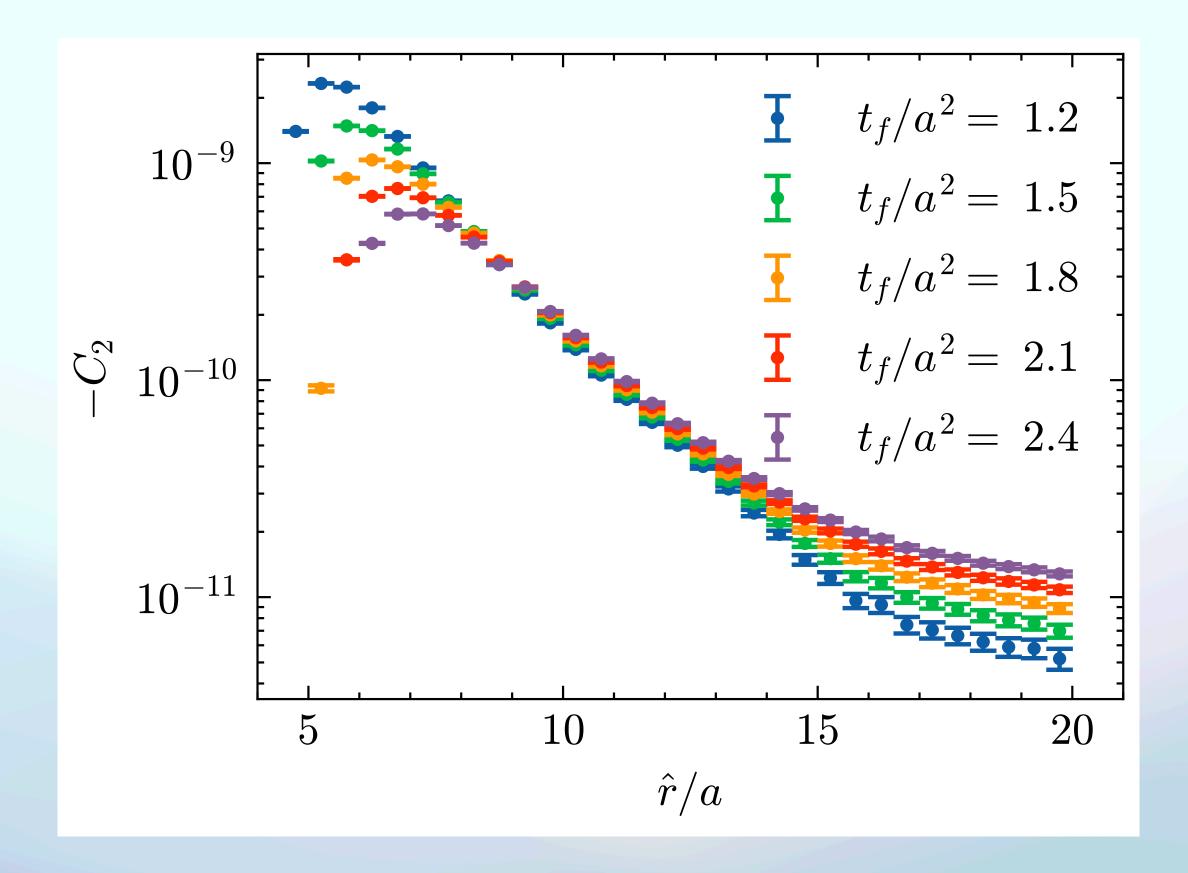
$$C_2(r) = \frac{1}{VN_r} \sum_{y \in \{|y-x|=r|\}} \sum_{x} \langle q(y) | q(x) \rangle$$



- The reflection positivity and the fact that the topological charge is reflection-odd ensure that the correlation functions are negative at non-zero distance.
- The positive contact contribution is a delta function at r=0. However, the topological charges are not locally defined on the lattice, the positive contact contribution is smeared to cover several lattice spacings.
- The data points in the negative tail contain important physical information.

η and η' on the lattice

- \bullet Bin data in $\delta r = 0.5a$ intervals to reduce strongly correlated between data points
- Use Wilson flow to smear gauge fields for improved signal-to-noise ratio



Ensemble details

Symbol	L/T	Mpi (MeV)	Mk(MeV)	a (fm)	Ncfg
481	48/96	139	499	0.1141(2)	356
641	64/128	139	508	0.0837(2)	330

R. Arthur et al., PRD87, 094514 (2013)

T. Blum et al., PRD93, 074505 (2016)

P. Boyle et al., PRD 93, 054502 (2016)

- We employ two 2+1-flavor gauge ensembles of domain wall fermions generated by the RBC/UKQCD collaboration.
- Both of them are at the physical pion mass point.
- The discretization effects of our results are carefully estimated.

Effective mass

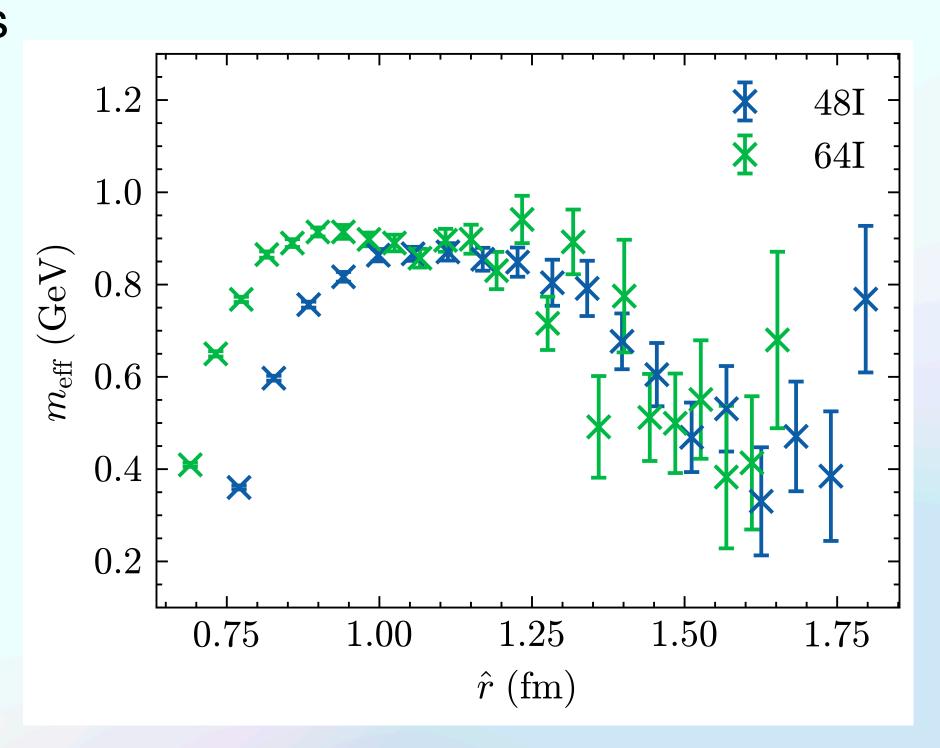
The correlation functions of topological charge density operators can be expressed as

$$C_2(\hat{r}) = \sum_n A_n \frac{m_n}{4\pi^2 \hat{r}} K_1(m_n \hat{r})$$

 K_1 is the modified Bessel function of the second kind.

The effective mass at each \hat{r} is obtained by solving the following equation numerically.

$$\frac{K_1(m_{\text{eff}}\hat{r})}{K_1\left((m_{\text{eff}}(\hat{r}-\delta_r)\right)}\frac{\hat{r}-\delta_r}{\hat{r}} = \frac{C_2(\hat{r})}{C_2(\hat{r}-\delta_r)}$$

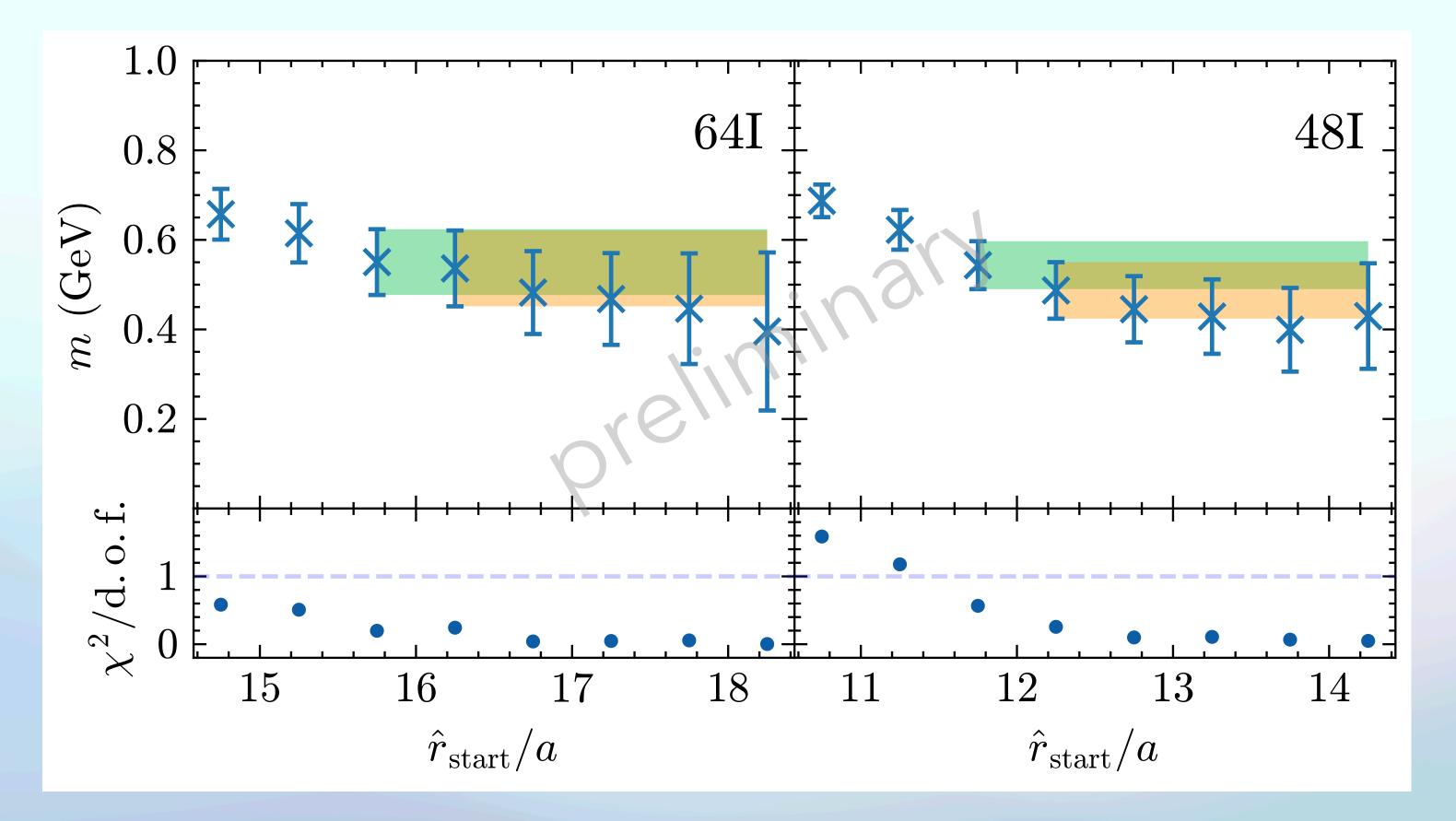


When \hat{r} is larger than ~1.4 fm, the effective masses are close to the η mass. And in the middle \hat{r} region, the effective masses are close to the η' mass.

Fittings and Results: mass of η

From two-point correlation functions to extracting the mass of η and η' :

Single state fitting:
$$C_2(\hat{r}) = A_{\eta} \frac{m_{\eta}}{4\pi^2 \hat{r}} K_1(m_{\eta} \hat{r})$$



$$m_{\eta}^{64{
m I}} = 0.550(73)\,{
m GeV}$$
 $m_{\eta}^{48{
m I}} = 0.543(53)\,{
m GeV}$ Consistent within errors

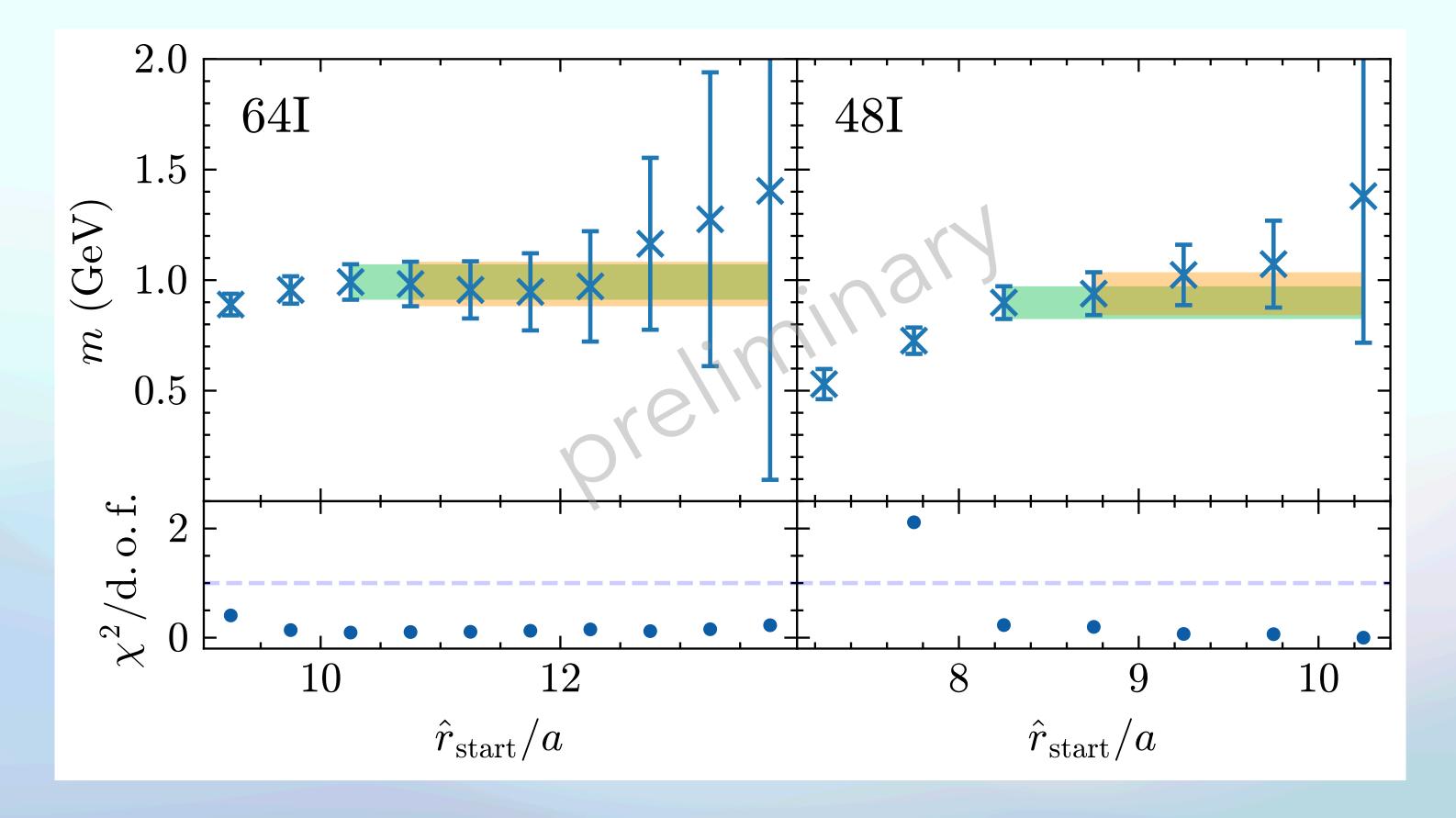
Our final results represent an average of the results from the two lattices.

$$m_{\eta} = 0.546(43)(5) \text{ GeV}$$

Fittings and Results: mass of η'

Double states fitting:
$$C_2(\hat{r}) = A_{\eta'} \frac{m_{\eta'}}{4\pi^2 \hat{r}} K_1 \left(m_{\eta'} \hat{r} \right) + A_{\eta} \frac{m_{\eta}}{4\pi^2 \hat{r}} K_1 \left(m_{\eta} \hat{r} \right)$$

Use previously fitted η mass as priors in these fits



$$m_{\eta'}^{64\text{I}} = 0.992(80) \text{ GeV}$$

 $m_{\eta'}^{48\text{I}} = 0.898(74) \text{ GeV}$

Following the same argument and strategy for the η case.

$$m_{\eta'} = 0.941(54)(50) \text{ GeV}$$

Fittings and Results: mixing angle

Double states fitting:
$$C_2(\hat{r}) = A_{\eta'} \frac{m_{\eta'}}{4\pi^2 \hat{r}} K_1\left(m_{\eta'}\hat{r}\right) + A_{\eta} \frac{m_{\eta}}{4\pi^2 \hat{r}} K_1\left(m_{\eta}\hat{r}\right)$$

In the leading-order, the octet–singlet eigenstates (η_8, η_1) are related to the physical mass eigenstates (η, η') by:

$$\begin{pmatrix} |\eta\rangle \\ |\eta'\rangle \end{pmatrix} = \begin{pmatrix} \cos\theta & -\sin\theta \\ \sin\theta & \cos\theta \end{pmatrix} \begin{pmatrix} |\eta_8\rangle \\ |\eta_1\rangle \end{pmatrix}$$

 θ is the $\eta - \eta'$ mixing angle in the octetsinglet basis.

Modern treatment of the mixing is parametrized by the matrix elements using two mixing angles as:

$$\begin{cases} F_{\eta'}^1 = F^1 \cos \theta_1, F_{\eta'}^8 = F^8 \sin \theta_8 \\ F_{\eta}^1 = -F^1 \sin \theta_1, F_{\eta}^8 = F^8 \cos \theta_8 \\ \text{L. Gan, et al. Phys.Rept. 945 (2022)} \\ \begin{cases} F_{\eta/\eta'}^1 = \langle 0 \, | \, O_1 \, | \, \eta/\eta' \rangle \\ F_{\eta/\eta'}^8 = \langle 0 \, | \, O_8 \, | \, \eta/\eta' \rangle \end{cases}$$

 O_8 and O_1 being the flavor octet and singlet operators.

Fittings and Results: mixing angle

Double states fitting:
$$C_2(\hat{r}) = A_{\eta'} \frac{m_{\eta'}}{4\pi^2 \hat{r}} K_1 \left(m_{\eta'} \hat{r} \right) + A_{\eta} \frac{m_{\eta}}{4\pi^2 \hat{r}} K_1 \left(m_{\eta} \hat{r} \right)$$

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In our case, the topological charge operator is a flavor single operator, we use $O_1 = q$, and thus we arrive at:

$$\tan \theta_1 = -\frac{\langle 0 | q | \eta \rangle}{\langle 0 | q | \eta' \rangle}$$

In the two-state fits:

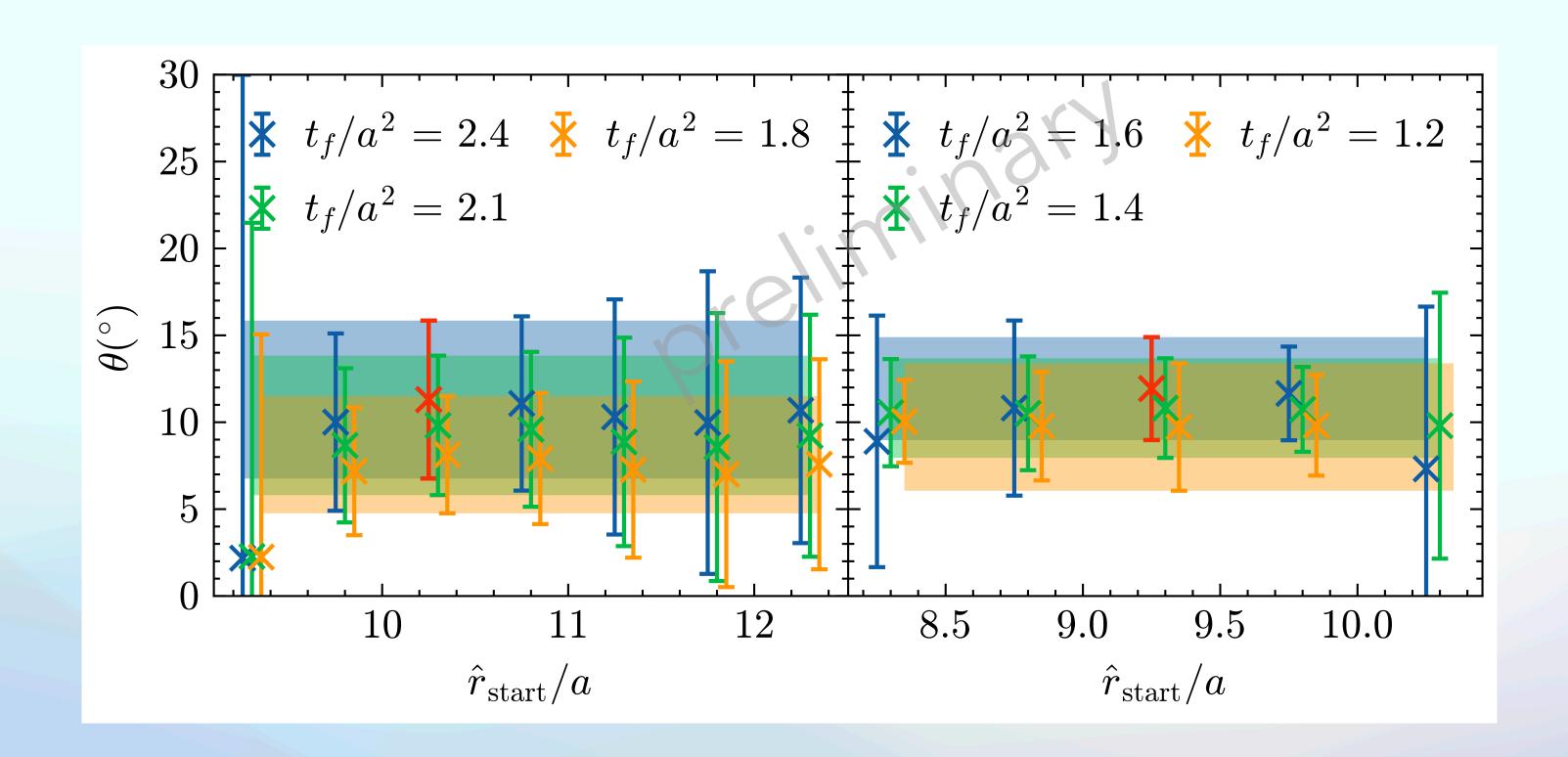
$$A_{\eta/\eta'} = \left| \left\langle 0 \mid q \mid \eta/\eta' \right\rangle \right|^2$$

Mixing angle:
$$\theta_1 = -\arctan \sqrt{A_{\eta}/A_{\eta'}}$$

 O_8 and O_1 being the flavor octet and singlet operators.

Fittings and Results: mixing angle

Double states fitting:
$$C_2(\hat{r}) = A_{\eta'} \frac{m_{\eta'}}{4\pi^2 \hat{r}} K_1 \left(m_{\eta'} \hat{r} \right) + A_{\eta} \frac{m_{\eta}}{4\pi^2 \hat{r}} K_1 \left(m_{\eta} \hat{r} \right)$$



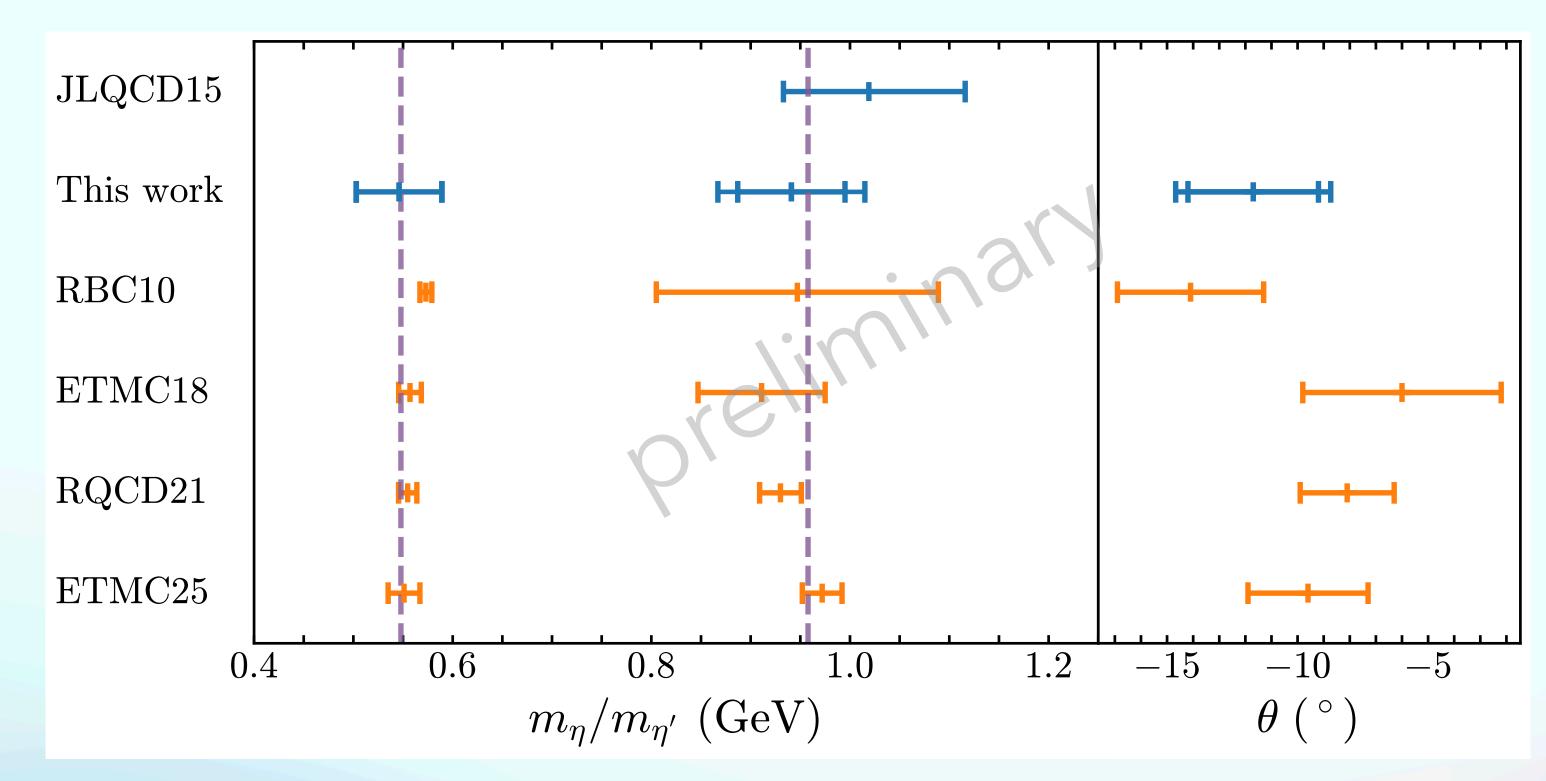
$$\theta_1^{64I} = -11.3(4.5)^\circ$$

$$\theta_1^{48I} = -11.9(3.0)^{\circ}$$

Consideration of continuum limit, and estimation of system uncertainty

$$\theta_1 = -11.7(2.5)(1.6)^{\circ}$$

Results



$$m_{\eta} = 0.546(43)(5) \text{ GeV}$$
 $m_{\eta'} = 0.941(54)(50) \text{ GeV}$
 $\theta_1 = -11.7(2.5)(1.6)^{\circ}$

- H. Fukaya, et al., Phys. Rev. D 92, 111501 (2015)
- N. H. Christ, et al., Phys. Rev. Lett. 105, 241601 (2010)
- K. Ottnad, et al., Phys. Rev. D 97, 054508 (2018)
- G. S. Bali, et al., JHEP 08, 137(2021)
- K. Ottnad, et al., arXiv:2503.09895 [hep-lat]
- *We first extract both the η and η' masses and also the mixing angle θ_1 by using topological charge operators.
- *The topological charge operator is a flavor-singlet operator which couples primarily to the η' state $(|A_{\eta}/A_{\eta'}| = \tan^2\theta_1 \sim 0.036)$

Summary and outlook

- For the first time, based on gauge configurations with the physical pion mass, we extract both the η and η' masses and also the mixing angle θ_1 by using topological charge operators.
- The systematic uncertainties are carefully considered.
- Compared with quark bilinear operators, the topological charge operator achieves high precision while avoiding complex computations and reducing computational resource consumption.

Summary and outlook

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- The systematic uncertainties are carefully considered.
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Thank you!