

Phase shifts in $l = 2 \pi\pi$ heavy-pion-scattering from two approaches

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Introduction

- Goal:
 - explain nuclear physics by dynamics of QCD (*binding energies, scattering phases, dense matter EoS, etc.*)
- Well-known approaches (from ab-initio QCD):
 - Lüscher's finite volume method ([Commun.Math.Phys. 105 \(1986\)](#))
 - non-zero chemical potential
- new approach:
 - potential method ([arXiv:1203.3642 \(HAL-QCD 2012\)](#))
 - untested, potential systematic drawbacks
- ☞ test and compare potential method to the Lüscher method
([Commun.Math.Phys. 105 \(1986\)](#)) in case of less complicated $J=2$
 $\pi\pi$ -scattering problem

Ingredients

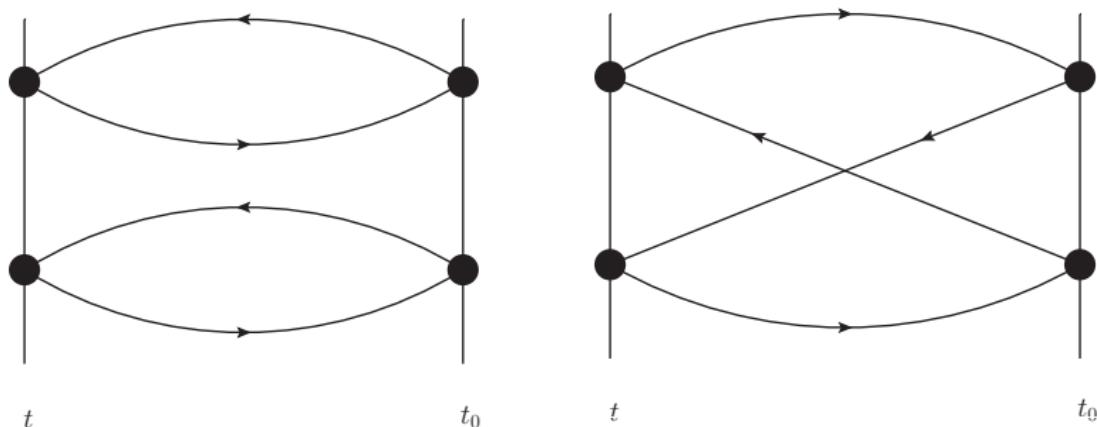
- compute two- and single-pion correlators:

$$C_{\pi\pi}(t, \mathbf{r}, \mathbf{P}) \equiv \sum_{\mathbf{R}} e^{-i\mathbf{RP}} \langle \pi^+(t, (\mathbf{R} + \mathbf{r})/2) \pi^+(t, (\mathbf{R} - \mathbf{r})/2) J_{\pi^-}(0, \mathbf{P}) J_{\pi^-}(0, \mathbf{0}) \rangle,$$

$$C_\pi(t, \mathbf{P}) \equiv \sum_{\mathbf{R}} e^{-i\mathbf{RP}} \langle \pi^+(t, \mathbf{R}) J_{\pi^-}(0, \mathbf{P}) \rangle$$

- J_{π^-} : (random) wall or Gaussian sources (Coulomb gauge)
- AP-BC or Dirichlet-BC ($\bar{\psi}(N_T/2, \mathbf{x}) = \psi(N_T/2, \mathbf{x}) = 0$)

Contractions



- contractions for $C_{\pi\pi}(t, \mathbf{r}, \mathbf{P})$
- ✓ no disconnected contributions in $I=2$ channel

Time dependent potential method I

- Nambu-Bethe-Salpeter (NBS) wave-function with asymptotic momentum \mathbf{k} given by [\(arXiv:1203.3642 \(HAL-QCD 2012\)\)](https://arxiv.org/abs/1203.3642)

$$\psi_{\mathbf{k}}(\mathbf{r}) \equiv \sum_{\mathbf{x}} \langle 0 | \pi^+(\mathbf{x}) \pi^+(\mathbf{x} + \mathbf{r}) | \pi^-(\mathbf{k}) \pi^-(-\mathbf{k}) \rangle$$

- ☞ NBS-WF satisfies Schrödinger-equation for $E < E_{\text{thres}}$

$$\left(\mathbf{k}^2 + \nabla^2 \right) \psi_{\mathbf{k}}(\mathbf{r}) = m_\pi \int_{\mathbb{R}^3} d^3 r' U(\mathbf{r}, \mathbf{r}') \psi_{\mathbf{k}}(\mathbf{r}')$$

- ✓ asymptotic behavior

$$\psi_{\mathbf{k}}(\mathbf{r}) \xrightarrow{|\mathbf{r}|=r \rightarrow \infty} e^{i\delta(k)} \frac{\sin(kr + \delta(k))}{kr} + \dots$$

Time dependent potential method II

- define

$$R(t, \mathbf{r}) \equiv C_{\pi\pi}(t, \mathbf{r}, 0) / C_\pi^2(t, 0)$$

- ☞ decompose to

$$R(t, \mathbf{r}) = \sum_{\mathbf{k}} \psi_{\mathbf{k}}(\mathbf{r}) a_{\mathbf{k}} e^{-t\Delta E(\mathbf{k})}$$

where $\Delta E(\mathbf{k}) = 2\sqrt{\mathbf{k}^2 + m_\pi^2} - 2m_\pi$.

- it is

$$\left(-\frac{\partial}{\partial t} + \frac{1}{4m_\pi} \frac{\partial^2}{\partial t^2} \right) R(t, \mathbf{r}) = \sum_{\mathbf{k}} \frac{\mathbf{k}^2}{m_\pi} \psi_{\mathbf{k}}(\mathbf{r}) a_{\mathbf{k}} e^{-t\Delta E(\mathbf{k})}$$

Time dependent potential method III

- this yields

$$\left(-\frac{\partial}{\partial t} + \frac{1}{4m_\pi} \frac{\partial^2}{\partial t^2} + \frac{\nabla^2}{m_\pi} \right) R(t, \mathbf{r}) = \int_{\mathbb{R}^3} d^3 r' \textcolor{blue}{U}(\mathbf{r}, \mathbf{r}') R(t, \mathbf{r}')$$

- expand non-local potential for $I=2$ case:

$$\textcolor{blue}{U}(\mathbf{r}, \mathbf{r}') \rightarrow V_C(\mathbf{r}) \delta(\mathbf{r} - \mathbf{r}') + \mathcal{O}(\nabla^2)$$

- this allows us to compute LO potential V_C :

$$V_C(\mathbf{r}) = \frac{\nabla^2 R(t, \mathbf{r})}{m_\pi R(t, \mathbf{r})} - \frac{(\partial/\partial t) R(t, \mathbf{r})}{R(t, \mathbf{r})} + \frac{1}{4m_\pi} \frac{(\partial/\partial t)^2 R(t, \mathbf{r})}{R(t, \mathbf{r})}$$

Cooking recipe

- ✓ compute $R(t, r)$ on the lattice
- ✗ extract LO potential $V_C(r)$
- ✗ model potential, $V \rightarrow \infty$ limit
- ✓ solve SE for arbitrary k (ext. parameter) and obtain $\psi_k(r)$
- ✓ compute scattering phases using $\beta \equiv [r d \ln \psi_k / dr]_{r=R}$

$$\tan \delta(k) = \frac{kR j_0'(kR) - \beta j_0(kR)}{kR n_0'(kR) - \beta n_0(kR)}$$

where $j_0(\rho) = \sin \rho / \rho$ and $n_0(\rho) = -\cos \rho / \rho$

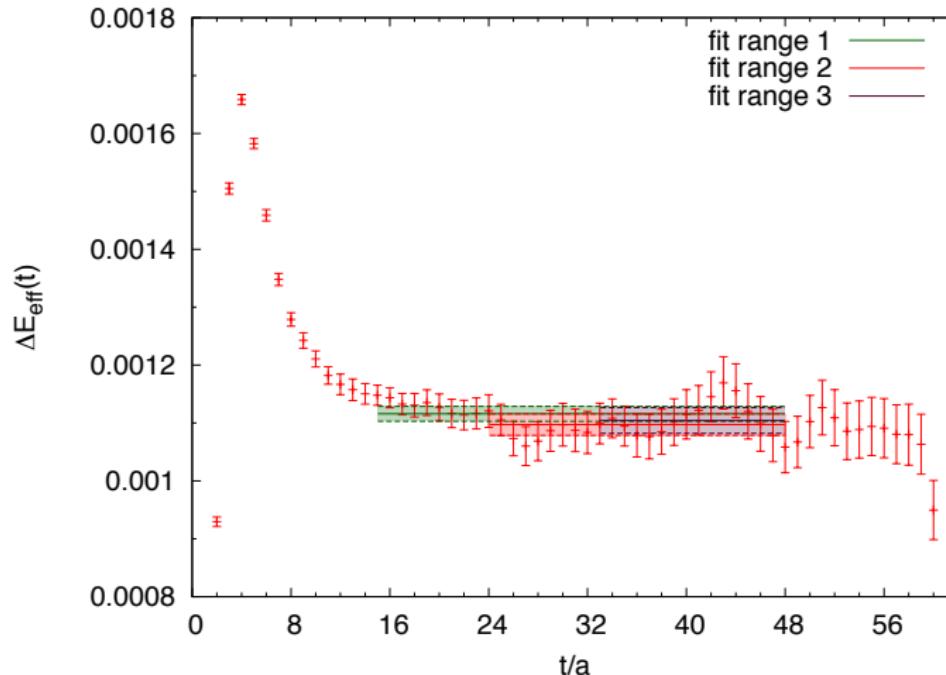
- ☞ compute low energy observables by fitting $\delta(k)$ to ERE:

$$\frac{k \cot \delta(k)}{m_\pi} = \frac{1}{m_\pi} \textcolor{magenta}{a_{\pi\pi}^{l=2}} + \frac{1}{2} m_\pi \textcolor{blue}{r} \left(\frac{k^2}{m_\pi^2} \right) + \textcolor{brown}{P}(m_\pi \textcolor{blue}{r})^3 \left(\frac{k^2}{m_\pi^2} \right)^2 + \dots$$

Setup

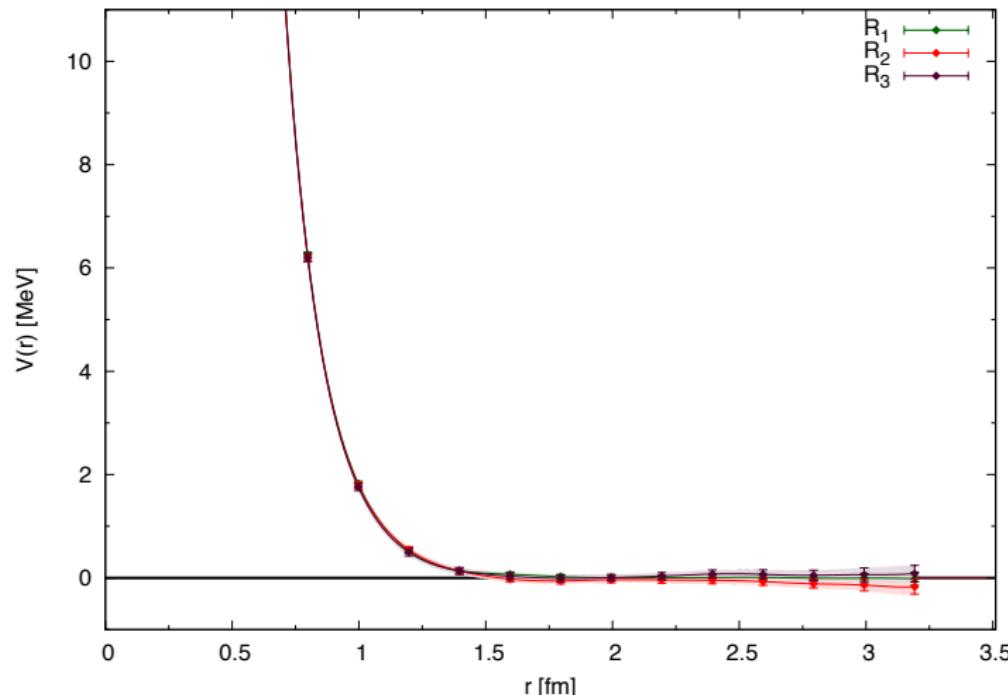
- ☞ test method on quenched setup using
 $M_\pi \sim (330 - 940) \text{ MeV}$ with 2 HEX smeared tree-level improved clover-Wilson quarks
- $a \approx 0.115 \text{ fm}$ and $L \approx 3.7 \text{ fm}$ (wall source: $1.84 < L < 4.9 \text{ fm}$)
- statistical error \Rightarrow 2000 bootstrap samples
- systematic uncertainties \Rightarrow histogram method
 - ✓ rotational invariance breaking \Rightarrow extract potential along axis, surface-diagonals and cubic diagonal
 - ✓ source dependence \Rightarrow use wall and gauss sources (radius $r \approx 0.3 \text{ fm}$)
 - ✓ ground state saturation (energy dependence) \Rightarrow use different time-slices
 - ✓ potential modeling \Rightarrow use different potential models (empirical) ☞ (barycentric) rational interpolation
 - ✓ asymptotic regime of ψ_k \Rightarrow different distances R

Mass Plateau



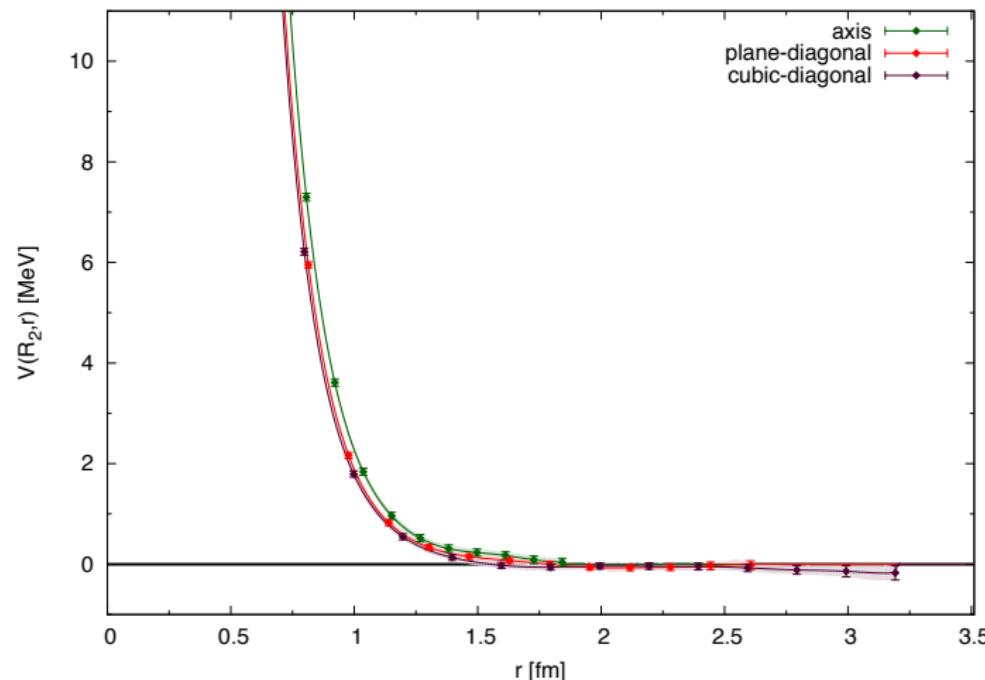
✓ clean plateaus

Potentials and excited state effects



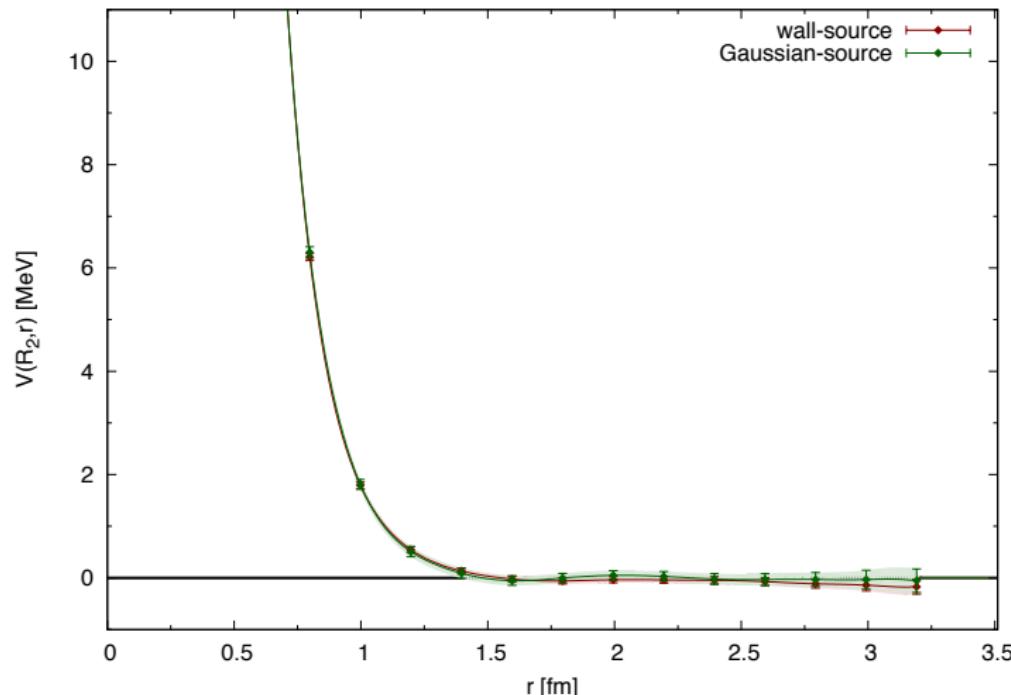
✓ no time-dependence (for $t \gtrsim 1.73$ fm)

Direction dependence



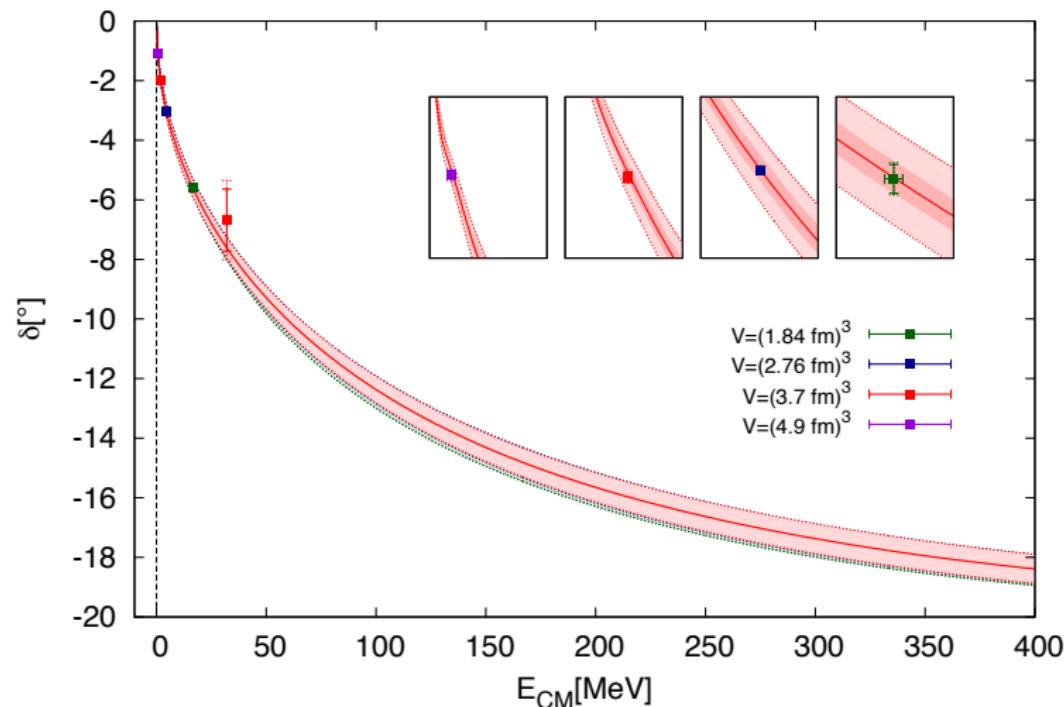
👉 breaking of rotational invariance dominant

Source-dependence



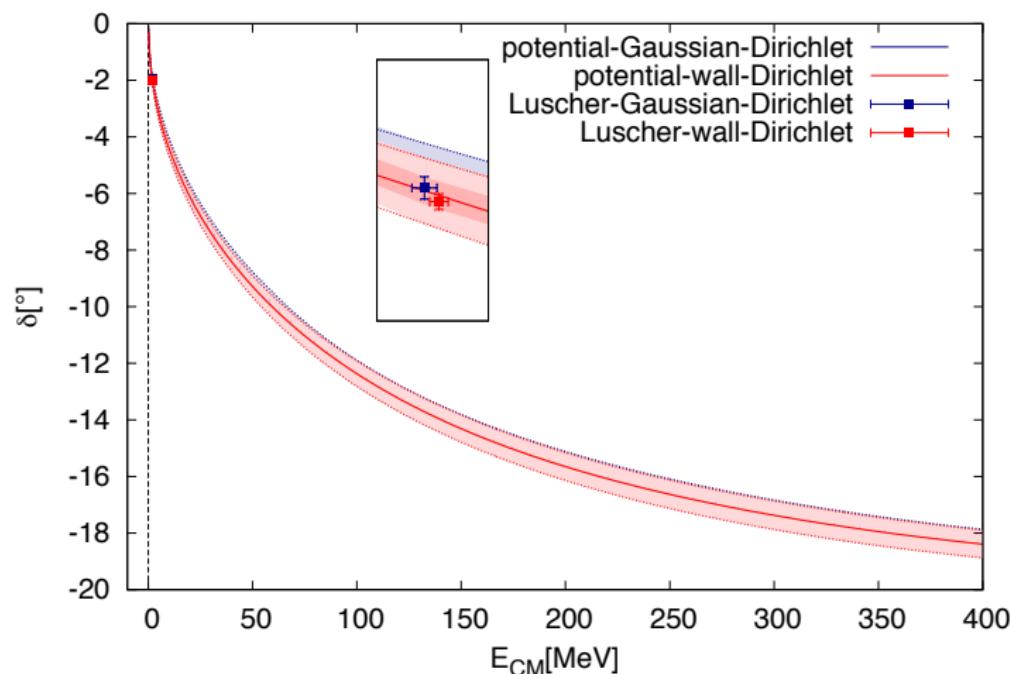
✓ no source-dependence

Scattering phase shifts for different volumes



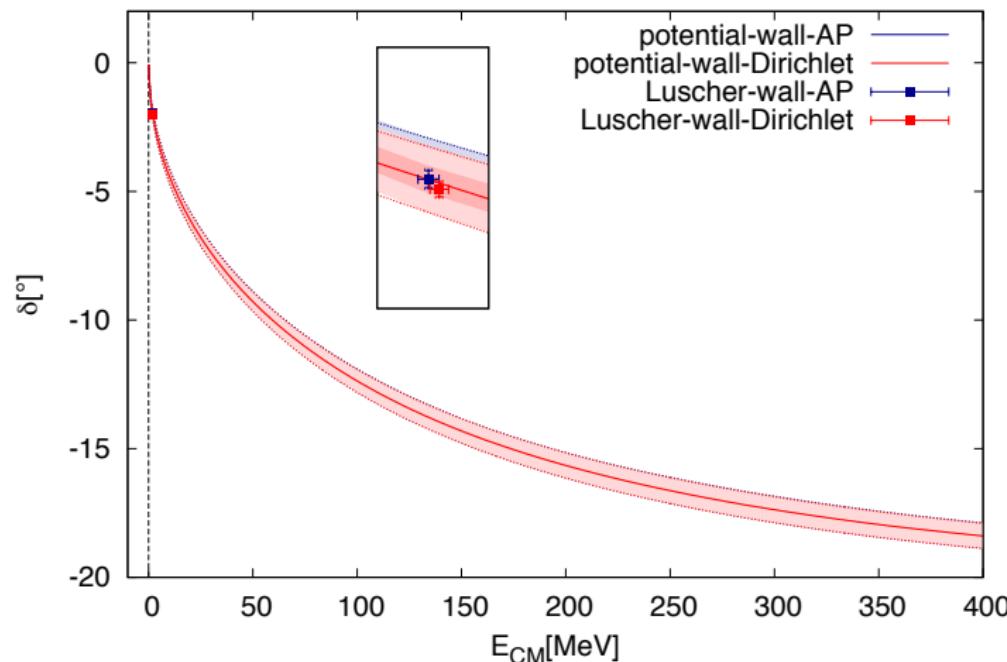
✓ results from Lüscher's and potential method agree

Different sources



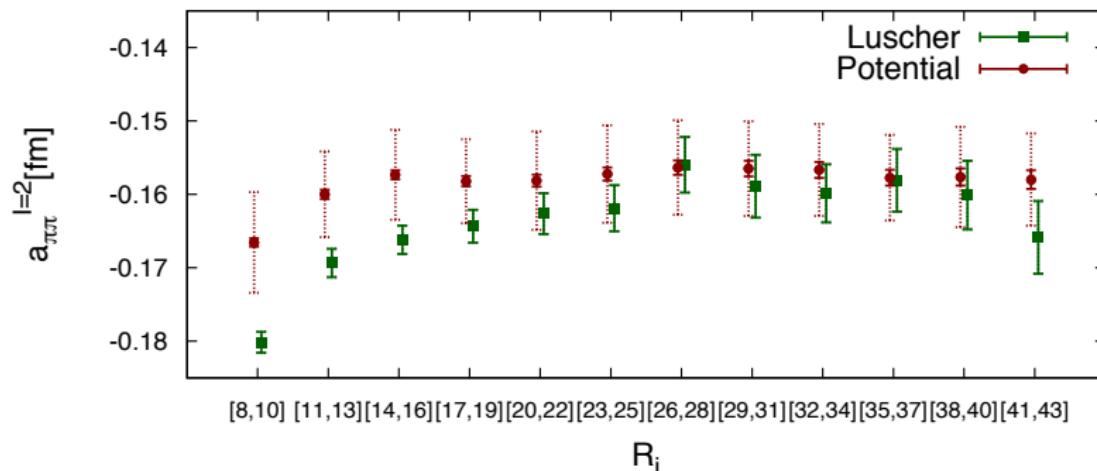
✓ phases independent of source-type

Different BC



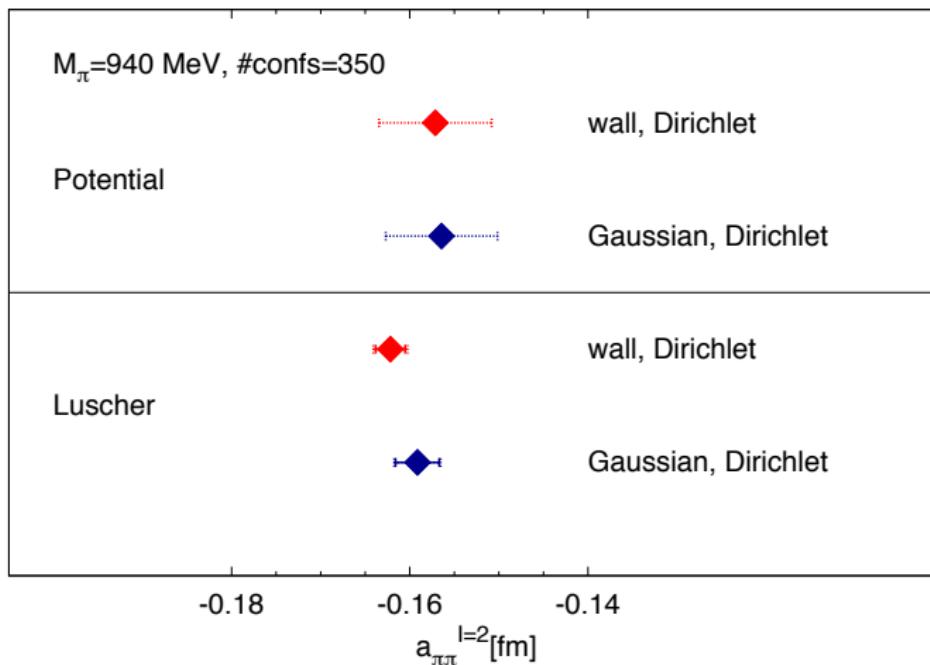
✓ phases independent of BC

Phases V



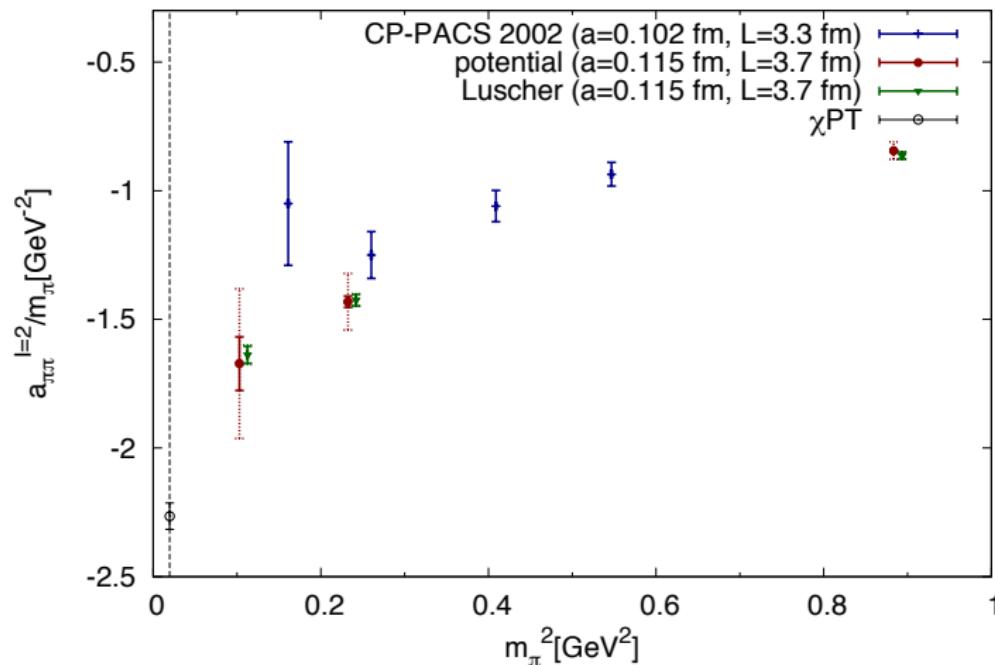
- ✓ potential method less sensitive to excited states contaminations

Scattering lengths



✓ scattering lengths agree for $M_\pi \approx 940 \text{ MeV}$

Mass dependence of scattering lengths



✓ agreement with previous quenched calculations

Summary

- ☛ compared potential method to Lüscher's approach to scattering problems
- ✓ both allow for extracting scattering phases and lengths
- ✓ compatible results
- ✗ Lüscher method: mapping out $\delta(k)$ requires use of different volumes, excited state extractions or boosted frames
- ✓ potential method: k^2 is free parameter
- ✓ potential method can use earlier time-slices \Rightarrow advantageous for multi-baryon scattering