

Decay of ρ and a_1 mesons on the lattice using distillation

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We extract the P-wave $\pi\pi$ phase shift for five values of pion relative momenta, which gives information on the ρ resonance. The Breit-Wigner formula describes the $\pi\pi$ phase shift dependence nicely and we extract $m_{\rho} = 792(7)(8)$ MeV and the coupling $g_{\rho\pi\pi} = 5.13(20)$ at our $m_{\pi} = 266$ MeV. We extract the P-wave scattering length $a_{l=1}^{\pi\pi} = 0.082(10)(3)$ fm³ from the state with the lowest pion relative momenta.

We also determine the S-wave $\rho\pi$ phase shift for two values of relative momenta, which provides parameters of the lowest axial resonance $a_1(1260)$. Using the Breit-Wigner fit we extract $m_{a1} = 1.44(4)$ GeV and the coupling $g_{a_1\rho\pi} = 1.1(3)$ GeV. From the lowest state we also extract the $\rho\pi$ scattering length $a_{l=0}^{\rho\pi} = 0.23(12)$ fm for our m_{π} .

The simulation is performed using one $N_f = 2$ ensemble of gauge configurations with cloverimproved Wilson quarks. The phase shifts are determined from the lowest two energy-levels, which are obtained by the variational analysis with a number of quark-antiquark and mesonmeson interpolators. The correlation functions are calculated using the distillation method with the Laplacian Heaviside (LapH) smearing of quarks.

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1. Introduction

Extracting the width of a hadronic resonance *R* from lattice QCD is challenging. The only proper method used up to now applies to resonances *R* that appear in the elastic scattering of two hadrons $H_1H_2 \rightarrow R \rightarrow H_1H_2$. First the elastic phase shift $\delta(s)$ for H_1H_2 scattering has to be determined from the lattice for several values of $s = E_{CM}^2 = E^2 - \mathbf{P}^2$, where *E* and **P** are the energy and the total momentum of the H_1H_2 system. Lüscher has shown that the energy *E* of two hadrons in a box of size $L \simeq$ few fm provides the value of the infinite-volume elastic phase shift $\delta(s)$ at $s = E^2 - \mathbf{P}^2$ [1]. His relation between δ and *E* for $\mathbf{P} = 0$ was generalized to $\mathbf{P} \neq 0$ in [2, 3, 4]. In practice, one or two lowest energy levels *E* are extracted and a few choices of **P** are used in order to extract $\delta(s)$ at different values of $s = E^2 - \mathbf{P}^2$.

The resulting $\delta(s)$ can be fit with a Breit-Wigner (or any other desired) form, where both are related via the scattering amplitude a_l for the *l*-th partial wave

$$a_{l} = \frac{-\sqrt{s}\Gamma_{R}(s)}{s - m_{R}^{2} + i\sqrt{s}\Gamma_{R}(s)} = \frac{e^{2i\delta(s)} - 1}{2i} \quad \text{or} \quad \sqrt{s}\Gamma_{R}(s)\cot\delta(s) = m_{R}^{2} - s \,, \quad \Gamma_{R}(s) \propto g_{RH_{1}H_{2}}^{2} \frac{p^{*2l+1}}{s}$$
(1.1)

and p^* is the momentum of H_1 and H_2 in their center-of-momentum (CMF) frame. This relation can be used to extract the mass m_R and the width $\Gamma_R = \Gamma_R(m_R^2)$ of the resonance from lattice data on $\delta(s)$. The width depends significantly on the phase space and therefore on m_{π} , so it is common to extract the coupling $g_{RH_1H_2}$, which is expected to depend only mildly on m_{π} .

Among all the meson resonances only the ρ meson width has been determined properly using this method. The first lattice determination was done by PACS-CS in 2007 [5]. Since then, several studies of the ρ have been carried out [6, 7], with the most recent ones [8, 9, 10]. In this talk we present our recent study of the ρ [9], which achieves the smallest statistical errors (on one ensemble only, however) on the resulting $\delta(s)$, m_{ρ} and Γ_{ρ} due to several improvements listed below.

We also extract the S-wave $\rho\pi$ elastic phase shift, which enables us to extract the mass m_{a1} and the width Γ_{a1} of the lowest lying axial resonance $a_1(1260)$. The lattice study of this resonance is especially welcome as the experimental knowledge on it is very poor: the width has a wide range $\Gamma_{a1}^{exp} = 250 - 600$ MeV [11], and none of its branching ratios have been reliably determined ¹ [11]. To our knowledge, this is the first lattice study aimed at the $\rho\pi$ scattering and Γ_{a1} .

2. Lattice simulation

We use 280 $N_f = 2$ configurations with tree-level clover-improved Wilson dynamical and valence quarks, corresponding to $m_{\pi}a = 0.1673(16)$ or $m_{\pi} = 266(3)(3)$ MeV [12]. The lattice spacing a = 0.1239(13) fm was determined using the Sommer parameter r_0 [9] and our $N_L^3 \times N_T = 16^3 \times 32$ is rather small, allowing us to use the powerful but costly full distillation method [13]. We combine periodic and anti-periodic propagators in time to reduce the finite N_T effects [9].

¹All final states are quoted just as "seen" in [11].



Figure 1: Contractions for our ρ and a_1 correlators with $\bar{q}q$ and meson – meson interpolators (I = 1).

3. ρ resonance and $\pi\pi$ phase shift

The details of our lattice simulation aimed at $\pi\pi$ phase shifts and the ρ resonance have been published in [9]. In this talk, we emphasize the most important steps and results.

The $\pi^+\pi^- \to \rho^0 \to \pi^+\pi^-$ scattering is elastic below the 4π threshold $\sqrt{s} < 4m_{\pi}$ and we can apply Lüscher's method. We determine the lowest two energy-levels of the $\rho^0 \leftrightarrow \pi^+\pi^-$ coupled system with $J^{PC} = 1^{--}$ and $|I, I_3\rangle = |1, 0\rangle$ for the following cases of total momentum *P*

Р	group	irrep	decay
0	O_h	T_1^-	$ ho_{3}(0) ightarrow \pi(\mathbf{e_3})\pi(-\mathbf{e_3})$
$\frac{2\pi}{L}$ e ₃	D_{4h}	A_2^-	$ ho_3(\mathbf{e_3}) ightarrow \pi(\mathbf{e_3})\pi(0)$
$\frac{2\pi}{L}(\mathbf{e_1}+\mathbf{e_2})$	D_{2h}	B_1^-	$ ho_{1,2}(e_1+e_2) o \pi(e_1+e_2)\pi(0)$

and all permutations in direction **P** and ρ -polarization. We display the symmetry group, the irreducible representation and the decay mode, which applies to three cases of **P** [2, 4, 8, 9].

Other simulations aimed at Γ_{ρ} used at most one quark-antiquark interpolator and one $\pi\pi$ interpolator for each **P**. We use 15 quark-antiquark interpolators $\mathcal{O}_{i=1-5}^{s=n,m,w}$ and one $\pi\pi$ interpolator for each **P**, where each pion is projected to a definite momentum:

$$\mathscr{O}_{i=1,\dots,5}^{s}(t) = \sum_{\mathbf{x}} \frac{1}{\sqrt{2}} \left[\bar{u}_{s}(x) \,\mathscr{F}_{i} \, \mathrm{e}^{\mathrm{i}\mathbf{P}\mathbf{x}} \, u_{s}(x) - \bar{d}_{s}(x) \,\mathscr{F}_{i} \, \mathrm{e}^{\mathrm{i}\mathbf{P}\mathbf{x}} \, d_{s}(x) \right] \qquad (s=n,m,w) \,, \tag{3.1}$$

$$\mathcal{O}_{6}^{n}(t) = \frac{1}{\sqrt{2}} [\pi^{+}(\mathbf{p_{1}})\pi^{-}(\mathbf{p_{2}}) - \pi^{-}(\mathbf{p_{1}})\pi^{+}(\mathbf{p_{2}})], \qquad \pi^{\pm}(\mathbf{p_{i}}) = \sum_{\mathbf{x}} \bar{q}_{n}(x)\gamma_{5}\tau^{\pm}e^{i\mathbf{p_{i}x}}q_{n}(x).$$

Quark-antiquark interpolators have five different color-spin-space structures \mathscr{F}_i . The quarks are smeared using the Laplacian Heaviside (LapH) smearing proposed in [13], i.e.,

$$q_{s} \equiv \Theta(\sigma_{s}^{2} + \nabla^{2}) q = \sum_{k=1}^{N_{v}} \Theta(\sigma_{s}^{2} + \lambda^{(k)}) v^{(k)} v^{(k)\dagger}, \quad s = n \text{ (narrow)}, m \text{ (middle)}, w \text{ (wide)}, (3.2)$$

where different truncations $N_v = 96$, 64, 32 correspond to three different widths s = n, m, w [9].

The 16×16 correlation matrix $C_{ij}(t_f, t_i) = \langle 0 | \mathcal{O}_i(t_f) \mathcal{O}_j^{\dagger}(t_i) | 0 \rangle$ necessitates the inclusion of the contractions in Fig. 1. The contractions were computed using the full distillation method, which is based on the LapH smeared quarks (3.2) [13] and leads to relatively precise results for all types of contractions in Fig. 1. We apply this method for the first time to extract a meson width. We also propose how to apply it for interpolators with different smearing widths in the same variational basis [9]. All correlators are expressed in terms of the so-called perambulators in Appendix A of [9]. The resulting correlators are averaged over all source time-slices t_i , over all directions of **P** and ρ polarization.

The lowest two energies of the system are determined using the Generalized Eigenvalue Method (GEVP) [14] and the dependence on the choice of the interpolators in the variational basis is explored in [9]. The lowest energy level is robust to this choice. We find that the first excited energy



Figure 2: The phase shift δ (in degrees) for $\pi\pi$ scattering in P-wave and $((ap^*)^3/\sqrt{sa^2})\cot\delta$ as a function of *s*, together with a Breit-Wigner fit.

level cannot be reliably obtained without the $\pi\pi$ interpolator in the basis, and that more than two interpolators are required at least in the case $P = \frac{2\pi}{L}(1,1,0)$. The extracted six energy levels for our preferred interpolator choice [9] are given in Table III of [9].

Each of the six energy levels gives the value of the phase shift $\delta(s)$ at $s = E^2 - \mathbf{P}^2$ (²) via the Lüscher formula for $\mathbf{P} = 0$ or its generalization to $\mathbf{P} \neq 0$ [2, 4]. We independently confirmed the needed relations and compiled them in [9]. One of these levels, $E_2(\mathbf{P} = 0)$, is above the inelastic threshold $\sqrt{s} > 4m_{\pi}$ and we omit it from further analysis.

The resulting phase shifts for five different values of *s* are plotted in Fig. 2. The phase shift has relatively small errors and exhibits a resonating behavior, which allows us to extract m_{ρ} and Γ_{ρ} or rather the coupling $g_{\rho\pi\pi}$. We use the Breit-Wigner relation (1.1) together with $\Gamma(s) \equiv g_{\rho\pi\pi}^2 p^{*3}/(6\pi s)$, which leads to

$$\frac{p^{*3}}{\sqrt{s}}\cot\delta(s) = \frac{6\pi}{g_{\rho\pi\pi}^2}(m_{\rho}^2 - s) .$$
(3.3)

This allows a linear fit in *s* (Fig. 2) to extract m_{ρ} and $g_{\rho\pi\pi}$ given in Table 1. The resulting $m_{\rho}a = 0.4972(42)$ is slightly lower than the naive value $m_{\rho}^{naive}a = 0.5107(40)$, which is extracted from the ground state with $\mathbf{P} = 0$. We also extract the P-wave scattering length $a_{l=1}^{\pi\pi} = 0.082(10)(3)$ fm³ (defined as $a_l \equiv lim_{\delta\to 0}\delta(p^*)/p^{*2l+1}$ [15]) from the state with the lowest³ $p^*a = 0.1076(36)$ and $\delta = 3.03(6)^{\circ}$. This quantity is not directly measured, so we compare it to the typical value $a_{l=1}^{\pi\pi} \simeq 0.038(2) (m_{\pi}^{phy})^{-3}$ obtained by combining experiment and ChPT or Roy equations [15].

A comparison of the resulting m_{ρ} and $g_{\rho\pi\pi}$ to two recent lattice simulations [8, 10] is compiled in [10]. The $N_f = 2$ simulation with twisted mass quarks [8] and the $N_f = 2 + 1$ simulation with Wilson quarks were done at four/two values of m_{π} and explicitly demonstrate the mild dependence of $g_{\rho\pi\pi}$ on m_{π} . All three results on $g_{\rho\pi\pi}$ are relatively close to each other and close to the $g_{\rho\pi\pi}^{exp} = 5.97$ extracted from Γ_{ρ}^{exp} . The resonance mass m_{ρ} of [10] is $\simeq 11\%$ higher than ours, while m_{ρ} of [8] is $\simeq 21\%$ higher than ours, at comparable m_{π} . Note that all three simulations get the resonance m_{ρ} within 3% from the value of m_{ρ}^{naive} , which implies that the simulations differ already in m_{ρ}^{naive} . Possible causes for different m_{ρ}^{naive} could be discretization effects or scale fixing of all

²We use the discrete dispersion relation $cosh(\sqrt{s}a) = cosh(Ea) - 2\sum_{k=1}^{3} sin^{2}(P_{k}a/2)$ instead of the continuum one $s = E^{2} - \mathbf{P}^{2}$ to analyze the ρ [9, 5]. We analyze the a_{1} using the continuum dispersion relation.

³The next state leads to $a_{l=1}^{\pi\pi}$ consistent with the value obtained from the lowest state.

_	$m_{ ho}$ [MeV]	8ρππ	$a_{l=1}^{\pi\pi}$	m_{a1} [GeV]	$g_{a_1\rho\pi}$ [GeV]	$a_{l=0}^{\rho\pi}$ [fm]	
latt	792(7)(8)	5.13(20)	0.082(10)(3)	1.44(4)	1.1(3)	0.23(12)	using $m_{ ho}$
				1.43(5)	1.7(4)	0.56(23)	using $m_{ ho}^{naive}$
exp	775.5	5.97	0.108(5) *	1.23(4)	< 1.35(30)	not meas.	

Table 1: Our lattice results for the resonance properties [9], compared to the experimental values. The results related to a_1 depend on the choice of the input ρ mass: m_{ρ} or m_{ρ}^{naive} . The experimental value of $a_{l=1}^{\pi\pi}$ is obtained combining experiment with ChPT or Roy equations.

three simulations, flavor breaking of twisted mass quarks [8] or partial quenching of the strange quark [8, 9]. Additional causes for the different m_{ρ} values could be the small interpolator basis in [8, 10] or the small box $L \simeq 2$ fm in [9]. The exponentially suppressed terms, which are neglected in Lüscher formulae, may not be completely negligible for our $L \simeq 2$ fm, which is a systematic uncertainty of our simulation. We are planing a simulation at larger *L* to explore possible finite size effects. We believe, however, that our small *L* does not influence our m_{ρ}^{naive} , as the first excited state $\pi(2\pi/L)\pi(-2\pi/L)$ at $\mathbf{P} = 0$ hardly affects the m_{ρ}^{naive} ground state.

Our $\delta(s)$ agrees reasonably well with the prediction of the lowest⁴ order of Unitarized Chiral Perturbation Theory [16], which has been recalculated for our $m_{\pi} = 266$ MeV.

4. The $\rho\pi$ phase shift and a_1 resonance

We study the S-wave scattering of $\rho \pi$, where the resonance $a_1(1260)$ appears, for the total momentum $\mathbf{P} = 0$. The scattering is elastic at least until $a_1(1260)$ on our lattice since \bar{K}^*K cannot be created on our $N_f = 2$ ensemble. The ground scattering state is $\rho(\mathbf{0})\pi(\mathbf{0})$ in the non-interacting limit. The scattering particle $\rho(\mathbf{0})$ is almost stable on our lattice, since its lowest decay channel $\pi(2\pi/L)\pi(-2\pi/L)$ is significantly higher in energy.

We use 9 quark-antiquark interpolators $\mathscr{O}_{i=1-3}^{s=n,m,w}$ and one $\rho(\mathbf{0})\pi(\mathbf{0})$ interpolator, all with $J^{PC} = 1^{++}$, $|I, I_3\rangle = |1, 0\rangle$ and $\mathbf{P} = 0$:

$$\mathcal{O}_{1}^{s}(t) = \sum_{\mathbf{x},i} \frac{1}{\sqrt{2}} \bar{u}_{s}(x) A_{i} \gamma_{i} \gamma_{5} e^{i\mathbf{P}\mathbf{x}} u_{s}(x) - \{u_{s} \leftrightarrow d_{s}\} \qquad (s = n, m, w) , \qquad (4.1)$$

$$\mathcal{O}_{2}^{s}(t) = \sum_{\mathbf{x},i,j,k} \frac{1}{\sqrt{2}} \bar{u}_{s}(x) \overleftarrow{\nabla}_{j} A_{i} \gamma_{i} \gamma_{5} e^{i\mathbf{P}\mathbf{x}} \overrightarrow{\nabla}_{j} u_{s}(x) - \{u_{s} \leftrightarrow d_{s}\} \qquad (s = n, m, w) , \qquad (4.1)$$

$$\mathcal{O}_{3}^{s}(t) = \sum_{\mathbf{x},i,j,k} \frac{1}{\sqrt{2}} \varepsilon_{ijl} \bar{u}_{s}(x) A_{i} \gamma_{j} \frac{1}{2} [e^{i\mathbf{P}\mathbf{x}} \overrightarrow{\nabla}_{l} - \overleftarrow{\nabla}_{l} e^{i\mathbf{P}\mathbf{x}}] u_{s}(x) - \{u_{s} \leftrightarrow d_{s}\} \qquad (s = n, m, w) , \qquad (4.1)$$

$$\mathcal{O}_{4}^{n}(t) = \frac{1}{\sqrt{2}} [\pi^{+}(\mathbf{0})\rho^{-}(\mathbf{0}) - \pi^{-}(\mathbf{0})\rho^{+}(\mathbf{0})] , \qquad \pi^{\pm}(\mathbf{0}) = \sum_{\mathbf{x}} \bar{q}_{n} \gamma_{5} \tau^{\pm} q_{n} , \qquad \rho^{\pm}(\mathbf{0}) = \sum_{\mathbf{x}} \bar{q}_{n} A_{i} \gamma_{i} \tau^{\pm} q_{n} ,$$

where ∇ is the covariant derivative. The contractions in Fig. 1 are calculated using the full distillation method and averaged over all source time slices t_i and all a_1 polarizations **A**.

The effective mass for the lowest two eigenvalues are shown in Fig. 3 and the resulting *E* and p^* are given in Table 2. The CMF momentum p^* is extracted using $E = \sqrt{p^{*2} + m_{\pi}^2} + \sqrt{p^{*2} + m_{\rho}^2}$:

⁴One cannot make a fair comparison between our lattice result and the NLO prediction, since it depends on a number of LECs, and some of them have been fixed using m_{ρ} from another lattice study, which gets a significantly higher m_{ρ} .



Figure 3: The effective mass for lowest two eigenvalues in a_1 channel (left). The combination $p^* \cot \delta / \sqrt{s}$ as a function of $s = E^2 - \mathbf{P}^2$, where δ is $\rho \pi$ phase shift in S-wave (right).

it is imaginary for the ground state below $m_{\pi} + m_{\rho}$ threshold, and real for the first excited state. We take two choices for the input ρ mass: our main results are based on the resonance mass m_{ρ} (green lines in Fig. 3), while m_{ρ}^{naive} is taken for comparison (blue lines in Fig. 3).

The S-wave phase shift δ for **P** = 0 is extracted using the well known Lüscher relation [1]

$$p^{*} \cot \delta = \frac{2}{\sqrt{\pi} L} Z_{00} \left(1, \left(\frac{p^{*}L}{2\pi} \right)^{2} \right) \xrightarrow{p^{*} \to 0} \frac{1}{a_{l=0}^{\rho \pi}} , \qquad (4.2)$$

which applies above and below threshold. The results are compiled in Table 2 for both choices of ρ mass. The first excited level gives $\delta \approx 90^\circ$, so it is sitting close to the top of the a_1 resonance and $m_{a1} \approx E_2$ holds. The ground state with imaginary p^* gives imaginary δ , but the product $p^* \cot \delta$ is real since $Z_{00}(1, (\frac{p^*L}{2\pi})^2)$ is real.

We parametrize $\Gamma_{a1}(s) \equiv g_{a_1\rho\pi}^2 p^*/s$ and apply the Breit-Wigner relation (1.1) to get

$$\frac{p^*}{\sqrt{s}}\cot\delta(s) = \frac{1}{g_{a_1\rho\pi}^2}(m_{a_1}^2 - s) , \qquad (4.3)$$

which applies in the vicinity of the resonance above or below threshold. Given the values of $p^* \cot \delta$ at two different values of *s*, we apply a linear fit (4.3) in *s* (shown in Fig. 3) to extract m_{a1} and $g_{a_1\rho\pi}$. The results are compiled in Table 1. Our m_{a1} at $m_{\pi} = 266$ MeV is about 14% higher than the experimental resonance $a_1(1260)$. The first lattice result for $g_{a_1\rho\pi}$ is valuable, since this coupling is not known experimentally. None of the a_1 branching ratios have been measured, so we provide only the upper limit for $g_{a_1\rho\pi}^{exp}$ resulting from the total width $\Gamma_{a1}^{exp} = 250 - 600$ MeV. Our lattice result $g_{a_1\rho\pi} = 1.1(3)$ GeV is in agreement with the value $g_{a_1\rho\pi}^{phen} \approx 0.9$ GeV obtained using Unitarized Effective Field Theory approach [17] and converted to our convention. We extract also $a_{l=0}^{\rho\pi}$ from the ground state, which is sufficiently close to the threshold. The scattering experiment cannot be carried out since ρ is a quickly decaying particle, so we compare our $a_{l=0}^{\rho\pi}(m_{\pi} = 266 \text{ MeV}) = 0.23(12)$ fm to $a_{l=0}^{\rho\pi}(m_{\pi}^{phy}) \approx 0.37$ fm obtained from Unitarized Effective Field Theory [18].

5. Conclusions

The lattice extraction of the phase shifts for elastic scattering has recently become possible also for the attractive resonant channels. We simulated the scattering in the ρ and a_1 channels and

level	fit	$Ea = \sqrt{sa}$	p^*a	δ	$p^* cos(\delta)/\sqrt{s}$	
1	7-10	0.6468(73)	i 0.065(13)	i 7.1(54)°	0.82(44)	(using m_{ρ})
			i 0.086(9)	i 23(14)°	0.34(14)	(using m_{ρ}^{naive})
2	6-9	0.897(13)	0.280(10)	83.7(59)°	0.034(33)	(using m_{ρ})
			0.272(10)	$88.9(59)^{\circ}$	0.005(31)	(using m_{ρ}^{naive})

Table 2: The results for the $a_1 \leftrightarrow \pi \rho$ coupled channel with interpolators $\mathscr{O}_{1,2,4}^n$ and GEVP reference time $t_0 = 5$. The ground state is below $\rho \pi$ threshold, so p^* and δ are imaginary, while the $p^* \cot \delta$ is real.

extracted the mass and the width of these two resonances as well as the scattering lengths in the corresponding meson-meson channels.

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