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Omega-Omega interaction on the Lattice

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We investigate interactions between Omega baryons in lattice QCD. Employing the HAL QCD method, we extract the Omega-Omega potential from the Nambu-Bethe-Salpeter (NBS) wave function, which are calculated on 2+1 flavor full QCD gauge configurations generated by the CP-PACS/JLQCD Collaboration at $m_{\pi} = 875$ MeV and $m_{\Omega} = 2108$ MeV. Both a shape of the potential and the phase shift of the Omega-Omega scattering calculated with it indicate that the interaction is strongly attractive. We finally discuss a possibility for an existence of a weak bound state in this system.

31st International Symposium on Lattice Field Theory LATTICE 2013 July 29, 2013 - August 3, 2013 Mainz, Germany

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1. Introduction

While hyperon interactions become important in hight density matters such as the core of the neutron star [1], investigations so far have been mainly focused on the octet sector. A main reason is, of course, that all decuplet baryons except Omega are unstable due to strong decays. Furthermore, even for the Omega-Omega case, it is still difficult to investigate their interaction experimentally due to their short-life time via weak decays. Theoretically, there exists two model calculations for the omega-omega inyteraction in the J = 0 channel [2, 3] leading to controversial results: while one reported a weakly repulsive interaction, the other indicated a strong attraction. The recent lattice QCD calculation using the standard finite volume method[Lueshcer] concluded that the Omega-Omega interaction is weakly repulsive in this channel: it's scattering length a =- 0.16±0.22 fm [4].

Recently, a new but first-principle method was proposed to investigate nucleon-nucleon interactions in QCD on the lattice [5–7], where the potential can be extracted from the Nambu-Bethe-Salpeter (NBS) wave function. This method has been generalized to derive potentials including hyperons (YN and YY) [8–13] and the three-nucleon force [14–16]. In this report, we therefore employ the HAL QCD method to study the Omega-Omega interaction, by calculating the corresponding potential in lattice QCD. Our results suggest an attractive Omega-Omega interaction in the J = 0 channel, which seems rather strong.

2. Extraction of potentials

The potential method was originally introduced by the HAL QCD collaboration [5], where a non-local potential which is defined from the equal time Nambu-Bethe-Salpeter (NBS) wave functions. For the two Omega system, the NBS wave function is defined as

$$\psi_n(\vec{r}) \equiv \langle 0 | \Omega(\vec{r}, 0) \Omega(\vec{0}, 0) | \Omega(k_n) \Omega(-k_n); in \rangle$$
(2.1)

where $|\Omega(k_n)\Omega(-k_n);in\rangle$ is an eigenstate of two-omega in QCD with the energy $2\sqrt{m_{\Omega}^2 + k_n^2}$, $\Omega(x)$ and $\overline{\Omega}(x)$ are local operators for Omega, whose explicit definition will be given in section 3.1. One of the most import ant properties here is that the NBS wave function at large $r = |\vec{r}|$ in QCD has the same asymptotic form to that of the scattering wave in quantum mechanics. From this fact, one can define a non-local but energy independent potential from the NBS wave function as

$$(E_n - H_0)\psi_n(\vec{r}) = \int d^3r' U(\vec{r}, \vec{r'})\psi_n(\vec{r'})$$
(2.2)

where $H_0 \equiv -\frac{1}{2\mu_{\Omega}} \nabla^2$ is the non-interaction part of the Hamiltonian, $\mu_{\Omega} \equiv \frac{m_{\Omega}}{2}$ is reduced mass, $E_n = k_n^2/(2\mu_{\Omega})$ is the kinetic energy in the center-of-mass frame, in two Omega system. The nonlocal potential $U(\vec{r}, \vec{r'})$ can be made to be energy-independent. Although we use the non-relativistic Schrödinger equation to define the potential, no non-relativistic approximation is made here [8]. Of course, potentials are not physical observables, but the potential defined above can reproduce physical observables such as phases of the S-matirx correctly by construction. In lattice QCD simulations, NBS wave unctions can be extracted from the two-Omega correlation function as

$$C_{\Omega\Omega}(\vec{r},t,t_0) = \frac{1}{V} \sum_{\vec{x}} \left\langle \Omega(t,\vec{x}+\vec{r})\Omega(t,\vec{x})\overline{\Omega}(t_0)\overline{\Omega}(t_0) \right\rangle = \sum_n \psi_n(\vec{r})a_n e^{-E_n(t-t_0)} + \cdots$$
(2.3)

$$\simeq a_0 \psi_0(\vec{r}) a_n e^{-E_0(t-t_0)}, \quad t-t_0 \to \infty$$
 (2.4)

with $a_n = \langle \Omega(k_n)\Omega(-k_n); in |\overline{\Omega}(t_0)\overline{\Omega}(t_0)| 0 \rangle$, where E_0 is the smallest energy of the system, and ellipses represent inelastic contributions.

In this report, we have employed the improved extraction, called the time dependence method [7], which is given by

$$\left(\frac{1}{4m}\frac{\partial^2}{\partial t^2} + \frac{1}{m}\nabla^2 - \frac{\partial}{\partial t}\right)R(\vec{r}, t, t_0) = \int d^3r' U(\vec{r}, \vec{r'})R(\vec{r'}, t, t_0)$$
(2.5)

where

$$R(\vec{r},t,t_0) \equiv \frac{C_{\Omega\Omega}(\vec{r},t,t_0)}{e^{-2m(t-t_0)}} = \sum_n a_n \psi_n(r) e^{-\Delta W_n(t-t_0)} + \cdots$$
(2.6)

with $\Delta W_n \equiv E_n - 2m_{\Omega}$. Since all two-Omega elastic scattering states with different *n* satisfy the same Schrödinger equation with the same energy-independent non-local potential, the large $t - t_0$ limit necessary for the ground state saturation is no longer required. A condition necessary for this method to work is that $t - t_0$ should be large enough to suppress both inelastic contributions in the two-Omega system and excited states in the single-Omega correlation function.

To extract potentials in practice, we employ the derivative expansion of the non-local potential [5]. We consider the leading-order term at low energies as

$$U(\vec{r},\vec{r}') = V_{\alpha'\beta'\alpha\beta}(\vec{r})\delta(\vec{r}-\vec{r}') + \mathcal{O}(\vec{\nabla})$$
(2.7)

where $\alpha', \beta', \alpha, \beta = 0 \sim 3$ are spin indices. In this report, we consider a case that the total spin of two Omega baryons is zero, whose potential depend on $r = |\vec{r}|$ only.

3. Symmetry of the Omega-Omega system

In this section we discuss a symmetry of the Omega-Omega system.

3.1 Symmetry

A single Omega baryon can not decay to a pair of an octet baryon and a pseudo-scalar meson in QCD, since the lowest energy pair, Ξ and K, has larger energy than the Omega baryon mass. On the lattice, we simply define local operators Omega and anti-Omega baryons as

$$\Omega_{\alpha,k}(x) \equiv \varepsilon^{abc} s_a^T(x) C \gamma_k s_b(x) s_{c\alpha}(x), \quad \overline{\Omega}_{\alpha,k}(x) \equiv \Omega_{\alpha,k}^{\dagger} \gamma^0 = \varepsilon^{abc} \overline{s}_{a\alpha}(x) \overline{s}_b^T(x) \gamma_{k_1} C \overline{s}_c(x)$$
(3.1)

where *a*, *b*, *c* are color indices, ε^{abc} is the totally anti-symmetric tensor, γ_k is the gamma matrix, α is the spinor index, and $C \equiv \gamma_4 \gamma_2$ is the charge conjugation matrix.

If a distance between two Omega baryons becomes large, we can neglect interactions between them. Such asymptotic Omega-Omega states can be classified by the orbital angular momentum

	P = +	P = -
J = 0	S = 0, L = 0; S = 2, L = 2	S = 1, L = 1; S = 3, L = 3
J = 1	S = 2, L = 2	S = 1, L = 1; S = 3, L = 3
J = 2	S = 2, L = 0; $S = 0, L = 2$; $S = 2, L = 2$;	S = 1, L = 1; S = 3, L = 1; S = 1, L = 3;
	S = 2, L = 4	S = 3, L = 3; S = 3, L = 5
J = 3	S = 2, L = 2; S = 2, L = 4	S = 3, L = 1; S = 1, L = 3; S = 3, L = 3;
		S = 3, L = 5
J = 4	S = 2, L = 2; S = 0, L = 4; S = 2, L = 4;	S = 3, L = 1; S = 1, L = 3; S = 3, L = 3;
	S = 2, L = 6	S = 1, L = 5; S = 3, L = 5; S = 3, L = 7

Table 1: Decomposition of the Omega-Omega system with conserved quantum numbers J and P in terms of states with a given (L, S).

(*L*), the total spin (*S*), the total angular momentum (*J*) and parity(*P*), even though conserved quantum numbers in QCD are *J* and *P* only. The two fermion state must change a sign under an exchange of them, while the asymptotic Omega-Omega state with given *L* and *S* has a factor $(-1)^{S+L+1}$ by the exchange, so that we should have S + L = even. In table1 we decompose states with conserved quantum numbers *J* and *P* in terms of asymptotic Omega-Omega states with given *L* and *S*.

3.2 Spin Projection

Since we employ wall sources, which has L = 0 and P = +, in our lattice calculations, the *J* is determined by the total spin *S* generated by the sources.

To construct the Omega-Omega operator which creates the state with the definite total spin S, let us first define a single spin 3/2 Omega baryon operator with a given S_z as

$$\Omega_{\frac{3}{2},\frac{3}{2}} \equiv -(\psi\Gamma_{+}\psi)\psi_{\frac{1}{2}}$$
(3.2)

$$\Omega_{\frac{3}{2},\frac{1}{2}} \equiv \frac{1}{\sqrt{3}} \left[\sqrt{2} (\psi \Gamma_Z \psi) \psi_{\frac{1}{2}} + (\psi \Gamma_+ \psi) \psi_{-\frac{1}{2}} \right]$$
(3.3)

$$\Omega_{\frac{3}{2},-\frac{1}{2}} \equiv \frac{1}{\sqrt{3}} [(\psi \sqrt{2} \Gamma_Z \psi) \psi_{-\frac{1}{2}} + (\psi \Gamma_- \psi) \psi_{\frac{1}{2}}]$$
(3.4)

$$\Omega_{\frac{3}{2},-\frac{3}{2}} \equiv (\psi \Gamma_{-} \psi) \psi_{-\frac{1}{2}}$$
(3.5)

where $\Gamma_{\pm} \equiv \frac{1}{2}(C\gamma^2 \pm iC\gamma^1)$, $\Gamma_Z \equiv \frac{-i}{\sqrt{2}}C\gamma^3$, so that spin 1 di-quark operators, $\psi\Gamma_+\psi$, $\psi C\Gamma_Z\psi$ and $\psi\Gamma_-\psi$, have $S_z = 1, 0, -1$, respectively, in non-relativistic limit.

Combining these operators, we can construct spin 3 ,spin 2 , spin 1,spin 0 states of Omega-Omega. The spin 0 state, used in this research, is given by

$$(\Omega\Omega)_{0,0} \equiv \frac{1}{2} \left(\Omega_{\frac{3}{2},\frac{3}{2}} \Omega_{\frac{3}{2},-\frac{3}{2}} - \Omega_{\frac{3}{2},\frac{1}{2}} \Omega_{\frac{3}{2},-\frac{1}{2}} + \Omega_{\frac{3}{2},-\frac{1}{2}} \Omega_{\frac{3}{2},\frac{1}{2}} - \Omega_{\frac{3}{2},-\frac{3}{2}} \Omega_{\frac{3}{2},\frac{3}{2}} \right).$$
(3.6)

4. Results

4.1 Lattice set up

In our calculation we employ 700 gauge configurations generated by CP-PACS and JLQCD Collaborations [17] with the renormalization group improved gauge action and the non-perturbatively

 $\mathcal{O}(a)$ improved Wilson quark action [17] at $\beta = 1.83$ ($a \simeq 0.12$ fm) on the $16^3 \times 32$ lattice, whose physical extension becomes L = 1.92 fm. The hopping parameter is (κ_{ud} , κ_s) = (0.13760, 0.13710), which gives $m_{\pi} = 875(1)$ MeV and $m_K = 916(1)$ MeV. The periodic boundary condition is imposed on the quark fields along the spatial direction, while the Dirichlet boundary condition is employed along the temporal direction on the time-slice t = $\frac{T}{2}$. To improve statistics, we calculate nine sources on different time slices per config, where the Dirichlet boundary is always separated from the source by $\frac{T}{2}$. Statistical errors are estimated by the Jackknife method where the bin-size is taken to be 1 configuration.

4.2 Omega-Omega Potential

We show the central potential between Omega-Omega in the ${}^{1}S_{0}$ channel at $t - t_{0} = 7, 8, 9$ in Fig1, where we use the notation ${}^{2S+1}L_{J}$ to specify quantum numbers of the channel. Overall structures of potentials are similar to those of NN potentials previously obtained in the lattice QCD [5]. This potential have repulsive core and deep attractive pocket, whose depth is 70~80 MeV. We observe that time dependence is small except for a long distance part at t = 9, which seems to be affected by finite volume effect. In future studies, we should investigate a volume dependence of the potential.



Figure 1: The central potential $V_c(r)$ between Omega-Omega in the ¹ S_0 channel at $t - t_0 = 7$ (blue) ,8(green), 9(red).

4.3 Phase shift and Scattering length

To calculate phase shift and scattering length, we first fit the potential with the form that

$$V(r) = a_1 e^{-a_2 r^2} + a_3 e^{-a_4 r^2} + a_5 e^{-a_6 r^2},$$
(4.1)

which gives, for example, $a_1 = -1.5(0.5) \times 10^2 \text{MeV}$, $a_2 = 2.0(0.6) \text{fm}^{-2}$, $a_3 = 2.0(0.9) \times 10^2 \text{MeV}$, $a_4 = 1.2(1.1) \times 10 \text{fm}^{-2}$, $a_5 = 1.0(0.1) \times 10^3 \text{MeV}$, $a_6 = 6.9(1.1) \times 10 \text{fm}^{-2}$ at $t - t_0 = 7$.

Using this fit result, we solve the Schrödinger equation in the infinite volume, and we obtain the phase shift $\delta(k)$ in the ¹S₀ channel. The scattering phase shift $\delta(k)$ at a given momentum k can be obtained as ratio

$$\tan \delta(k) = \lim_{x_1, x_2 \to \infty} \frac{\psi_k(x_2) \sin(kx_1) - \psi_k(x_1) \sin(kx_2)}{\psi_k(x_1) \cos(kx_2) - \psi_k(x_2) \cos(kx_1)}.$$
(4.2)

We show a center of mass energy dependence of the scattering phase shift in Fig. 2, where $E_{\rm CM} = k^2/(2\mu_{\Omega})$. As shown in the figure, while phase shift calculated from the potential at t=7 suggests an existence for a bound state, phase shifts from data at t=8 and 9 indicate that the Omega-Omega interaction is strongly attractive but it may not be strong enough to form a bound state at this quark mass.

The scattering length a and effective range r_e are defined by

$$k\cot\delta(k) = \frac{1}{a} + \frac{1}{2}r_ek^2 + O(k^4).$$
(4.3)

The fit of $k \cot \delta(k)$ near k = 0 by the above formula gives $1/a = -2.4(2.5) \times 10^{-1}$, $9.0(6.9) \times 10^{-2}$, $2.9(1.2) \times 10^{-1}$ fm⁻¹, and $r_e = 5.0(5.1) \times 10^{-3}$, $5.1(1.0) \times 10^{-3}$, $5.2(1.8) \times 10^{-3}$ fm at $t - t_0 = 7, 8, 9$, respectively. Unfortunately, errors of both 1/a and r_e are quite large, as expected from Fig. 2.



Figure 2: Scattering phase shift at $t - t_0 = 7$ (blue) ,8(green) , 9(red), as a function of the center of mass energy E_{CM} .

5. Summary

In this paper, we have investigated the Omega-Omega interaction in the channel with the orbital angular momentum L=0 and the total spin S=0 in 2+1 flavor lattice QCD, using the method developed by the HAL QCD collaboration. We find that the central potential obtained from the

NBS wave function for the ${}^{1}S_{0}$ channel shows attractions at long distance and the repulsive core at short distance. The phase shift derived from this potential shows that the Omega-Omega interaction is strongly attractive, while an existence for an Omega-Omega bound sate is unfortunately inconclusive at the pion mass in this study. In future we will examine the Omega-Omega interaction at more lighter pion masses in larger volumes.

Acknowledgements

We thank CP-PACS/JLQCD collaborations and ILDG/JLDG for providing us the 2 + 1 flavor gauge configurations [18]. We also thank authors and maintainer of CPS++ [19], whose modified version is used in this report. Numerical computations of this work have been carried out at KEK supercomputer system (BG/Q) and T2K at University of Tsukuba. This research is supported in part by JICFuS [20].

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