

Baryon-baryon potentials in the flavor SU(3) limit from lattice QCD

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We investigate baryon-baryon (BB) interactions in the flavor SU(3) symmetric world, by using the 3-flavor full QCD simulations. In the SU(3) limit, six independent two-baryon states exist for a given orbital angular momentum. We forcus on S-wave states and extract BB potentials using the method recently developed. We discuss the flavor dependence of the potentials, in particular, the strength of the repulsion at short distance. We find that the SU(3) flavor structure of the repulsive core is compatible with that expected from the Pauli principle for the quarks.

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1. Introduction

The generalized nuclear force, which includes not only the nucleon-nucleon (NN) interaction but also hyperon-nucleon (YN) and hyperon-hyperon (YY) interactions, has been one of the central topics in hadron and nuclear physics. Experimental studies on the ordinary and hyper nuclei as well as the observational studies of the neutron stars and supernova explosions are deeply related to the physics of the generalized nuclear force [1].

As far as the NN interaction is concerned, significant number of scattering data have been accumulated and are parametrized in terms of the phase shifts or of the phenomenological potentials [2]. In particular, the moderate attraction at low energies and the strong repulsion at high energies seen in the S-wave NN scattering phase shift is one of the characteristic features of the NN force.

On the other hand, the properties of the YN and YY interactions are hardly known even today due to the lack of high precision YN and YY scattering data. The approximate flavor SU(3) symmetry does not necessarily create a bridge between the NN force and the hyperon forces because there are six independent flavor-channels for the scatterings between octet baryons. Although there have been previous theoretical attempts to fill the gap on the basis of the phenomenological quark models [3] and of the one-boson-exchange models [4], it is much more desirable to study the generalized nuclear force from first principle QCD.

This report is our exploratory attempt to unravel the similarities and differences among different flavor representations of the baryon-baryon (BB) potentials on the basis of the lattice QCD simulations. To capture the essential part of physics, we take the finite but degenerated u,d,s quark masses, so that the system is in the flavor SU(3) limit. In the next section, we briefly review our method to study the baryon-baryon interaction in lattice QCD. In Section 3, we introduce the classification of baryon-baryon systems in the flavor SU(3) limit. In Section 4, we show a setup of our numerical simulations. In Section 5, we show the baryon-baryon potentials in the SU(3) limit and discuss their implications. The last section is devoted to summary and outlook.

2. Baryon-baryon potentials from lattice QCD

The potential is a useful concept to describe elastic BB interactions at low-energies. Recently a method to extract the potential from lattice QCD simulations has been proposed and applied to the NN system [5, 6], and ΞN and ΛN systems [7, 8, 9] as well. In this method, the potential is defined from the Bethe-Salpeter (BS) amplitude $\phi_E(\vec{r})$ through the Schrödinger equation as

$$\left[\frac{\nabla^2}{2\mu} + T\right]\phi_E(\vec{r}) = \int d^3\vec{r}' U(\vec{r},\vec{r}')\phi_E(\vec{r}'), \quad \text{with} \quad T = \frac{k^2}{2\mu}, \quad k^2 = \frac{E^2}{4} - \frac{M^2}{4}$$
(2.1)

where μ and M are reduced and total masses of the system. Here $U(\vec{r}, \vec{r}')$ is an energy independent but non-local potential. At low energies, it can be expanded by the local velocity $\vec{v} = \vec{\nabla}/\mu$ as $U(\vec{r}, \vec{r}') = V(\vec{r}, \vec{v})\delta(\vec{r} - \vec{r}') = (V_{LO} + V_{NLO} + V_{NNLO} + \cdots)\delta(\vec{r} - \vec{r}')$, where N^nLO term is of $O(v^n)$. Thus, in the leading-order (LO) approximation, we have

$$V(\vec{r}) = \frac{1}{2\mu} \frac{\nabla^2 \phi_E(\vec{r})}{\phi_E(\vec{r})} + T .$$
 (2.2)



HadronMass [1/a]Mass [MeV] π, K 0.5112(6)834.4(1.0) N, Λ, Σ, Ξ 1.068(2)1744.0(3.8)

Table 1: Hadron masses in the present lattice QCD simulation $(16^3 \times 32 \text{ lattice and spatial volume of about (2.0 fm)}^3)$. The lattice spacing a = 0.1209 fm is used to convert them in physical unit.

Figure 1: A part of all 64 octet-baryon pairs. Eight pairs composed of the proton and octet baryons are expanded in several multiplets.

It is shown in Ref. [10] that the LO potential $V(\vec{r})$ defined above is a good approximation of $U(\vec{r}, \vec{r}')$ at low energies between $T \sim 0$ MeV and $T \sim 45$ MeV for the *NN* system in quenched QCD. The BS wave function ϕ_E is extracted from the four point function, for example,

$$W(t - t_0, \vec{r}) = \sum_{\vec{x}} \langle 0 | p(t, \vec{x} + \vec{r}) n(t, \vec{x}) \bar{n}(t_0) \bar{p}(t_0) | 0 \rangle , \qquad (2.3)$$

which is dominated by the lowest energy state with the total energy E_0 at large time separation, so that it is proportional to $\phi_{E_0}(\vec{r})$ at large $t - t_0$.

3. Baryon-baryon potentials in the flavor SU(3) limit

In the flavor SU(3) limit, the ground states of a single baryon forms octet and decuplet. Systems of two octet baryons are classified into six irreducible multiplets in flavor SU(3);

$$8 \otimes 8 = \underbrace{1 \oplus 8 \oplus 27}_{\text{symmetric}} \oplus \underbrace{8 \oplus 10^* \oplus 10}_{\text{anti-symmetric}} . \tag{3.1}$$

Since baryons obey the Fermi statistics, there exist six independent states for a given orbital angular momentum. In particular for the S-wave, we have singlet, octet and 27-plet for the spin-singlet state $({}^{1}S_{0})$, while octet, anti-decuplet and decuplet for the spin-triplet state $({}^{3}S_{1})$. The six independent potentials for these S-wave states are the targets of our study. For example, we have computed $\sum_{\vec{x}} \langle 0|B_{i}B_{j}\overline{BB}^{(27)}|0\rangle$ in order to extract $V^{(27)}$ through eq.(2.2), where $\overline{BB}^{(27)}$ is a source operator of two octet baryons in the 27-plet representation, and $B_{i}B_{j}$ is a sink operator with two baryons which has an overlap with the 27-plet component.

We have to employ the source operator which belongs to the definite flavor representation in order to extract the corresponding BS wave functions at large time separation, since all two octet baryon states have the same energy in the infinite volume and possible energy differences due to the interactions in the finite volume are generally small. Fig 1 illustrates a part of all octet-baryon pairs. As shown in the figure, some of baryon pairs belong exclusively to 27-plet, anti-decuplet or decuplet. Numbers of wick contractions are relatively small for such operators, while more complicated source operators with the SU(3) Clebsh-Gordan coefficients are needed for other multiplets. Due to the limitation of computational resources, we consider only the potentials $V^{(27)}$, $V^{(10^*)}$ and $V^{(10)}$ for the moment, which can be extracted by these simpler source operators.



Figure 2: Effective mass plot of the 4-point func- Figure 3: Two baryon wave functions in 27-plet tion for the two-baryon operator in the 27-plet. The (red), anti-decuplet (green) decuplet (blue) represenhorizontal lines stand for the threshold of two free tations, normalized to one at the maximum distance. baryons and its error band.

4. Numerical setup

In order to calculate the potentials in 3-flavor full QCD, we have utilized a gauge configuration set of Japan Lattice Data Grid(JLDG)/International Lattice Data Grid(ILDG) generated by CP-PACS/JLQCD Collaborations on $16^3 \times 32$ lattice with the RG improved Iwasaki gauge action at $\beta = 1.83$ and the non-perturbatively O(a) improved Wilson quark action at lattice spacing a = 0.129fm and spatial volume of about (2.0 fm)³ [11, 12]. The hopping parameters of the configuration set correspond to the flavor SU(3) symmetric point, $\kappa_u = \kappa_d = \kappa_s = 0.13760$. Quark propagators are calculated from the spatial wall source at t_0 with the Dirichlet boundary condition in temporal direction at $t = t_0 \pm 16 \mod 32$. The wall source is placed at 4 different time slices on each of 800 gauge configurations, in order to enhance the signals, together with the average over forward and backward propagations in time. The average over discrete rotations of the cubic group is taken for the sink operator, in order to obtain A_1^+ state, whose lowest energy state is assumed to be dominated by the S-wave. The statistical data are divided into bins of the size eight and the jackknife prescription is adopted to estimate the statistical error. All of our numerical computations are carried out at KEK supercomputer system, Blue Gene/L and SR11000.

5. Results

Table 1 lists masses of the octet pseudo-scalar meson and baryon in our simulation, obtained by fitting corresponding 2-point functions. The mass of the octet pseudo-scalar meson is 834 MeV, which is much heavier than the physical π or K meson. For our present purpose, however, the flavor SU(3) symmetry is more important than the lighter *u*, *d* quark masses.

Fig 2 shows the effective mass plot of the 4-point function for two baryon operator in the 27plet representation. A plateau is observed at large time separation, $t - t_0 \ge 9$, so that the 4-point function is well saturated by the two-baryon ground state. A similar plateaus are obtained for antidecuplet and decuplet representation. We take $t - t_0 = 10$ to extract baryon-baryon potentials in this report.





Figure 4: The 27-plet potential, $V^{(27)}(r)$, without energy shift T_0 .

Figure 5: Scattering phase shift computed from the Schrödinger equation with $V^{(27)}$, as a function of the laboratory energy T_{lab} .

Fig 3 shows the BS wave function obtained from lattice QCD simulations. It is clearly seen that the wave functions depend on their flavor representation(27, 10 or 10^{*}). An agreement between *pn* ${}^{1}S_{0}$ state and $p\Sigma^{+} {}^{1}S_{0}$ state, both of which are the members of the 27-plet, gives a good consistency check of our calculations. These wave functions suggest presences of repulsive interaction at short distance (around r < 0.4 fm) for all channels. The shape of these wave functions indicates that no bound state seems to exist for all channels, though more detailed analysis is needed in the finite volume for the definite conclusion.

In Fig 4, the potential extracted from the BS wave function is plotted as a function of r in the case of the 27-plet (spin singlet) state. Note that here only the first term of eq.(2.2) is presented without the constant term T_0 , which is expected to be small from Fig 2. The potential $V^{(27)}$ in Fig 4 corresponds to the NN ${}^{1}S_0$ potential in the SU(3) limit. We see a repulsive core at short distance and an attractive pocket around 0.6 fm. These qualitative features are similar to what was found in the NN system [5]. Assuming that the leading order term in the velocity expansion works well as demonstrated for the NN potentials [10], we can calculate the "physical" observables in the SU(3) symmetric world. For example, Fig 5 shows the scattering phase shift as a function of energy in the laboratory system, computed from the Schrödinger equation with the baryon mass of 1744 MeV in the present simulation and with the potential $V^{(27)}$ where $T_0 = -10$ MeV is included.

The potential $V^{(10^*)}$ without T_0 in Fig 6 corresponds to the NN 3S_1 effective central potential in the SU(3) limit. We call it the "effective" central potential, since the mixing with the 3D_1 state is not separated from the true central potential in this calculation. We observe that $V^{(10^*)}$ has a weaker repulsive core and a broader attractive pocket than $V^{(27)}$ which corresponds to the NN 1S_0 potential. Accordingly, it is more attractive as a whole.

The potential $V^{(10)}$ without T_0 in Fig 7 looks different from others: its repulsive core is higher and broader, and no clear attractive pocket is observed. Accordingly, it is much more repulsive than others. This feature can be also seen from the wave function in Fig 3. The stronger repulsion in the decuplet channel has been predicted from the phenomenological quark model where the decuplet states are semi-forbidden state due to the Fermi statistics of quarks [3]. Our lattice QCD result is compatible with this Pauli principle picture. This is more clearly seen in Fig 8, where the







 $W_{01,01}$

W01.10

 $p\Sigma^+$ W_{01,10}

4

1.0

0.8

0.6

0.4

0.2 0.0

-0.2

-0.4

0

2

Components of BS amplitudes





Figure 8: Component BS amplitudes with Dirac in- Figure 9: Potential $V^{(10)}(r)$ without T_0 , extracted dexes of baryon operator. Wave function of decuplet from data at three time slice. is given by sum of $p\Sigma^+ W_{0101}$ and $p\Sigma^+ W_{0110}$.

6

8

10

4-point functions $W_{\alpha\beta,\alpha'\beta'}(r)$ with spinor indices $\alpha\beta,\alpha'\beta'$ are plotted. The spin triplet (flavor antisymmetric) component is given by the sum $W_{01,01} + W_{01,10}$. One can see that two $p\Sigma^+$ (decuplet) 4-point functions tend to cancel each other, so that the decuplet wave function tends to vanish as if it were the totally forbidden state, while no such cancellation is observed for the anti-decuplet. The potential $V^{(10)}$ likely has a shallow attractive pocket around 0.8 fm, as seen more clearly in Fig 9 at time slice $t - t_0 = 8$.

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6. Summary and outlook

We have performed the 3-flavor full QCD simulation to study the general features of the baryon-baryon interaction in the flavor SU(3) limit. From the BS wave function measured on the lattice, we extracted the leading-order potentials in the S-wave, $V^{(27)}$, $V^{(10^*)}$ and $V^{(10)}$, labeled by irreducible representations of flavor SU(3). All three potentials have a repulsive core at short distance in common, but its strength depends on the representation. In particular, $V^{(10)}$ is most repulsive among three, which may indicate the importance of the Pauli principle for quarks in baryon-baryon interactions at short distance. Also, both $V^{(27)}$ and $V^{(10^*)}$ show attractive pockets at intermediate distance. We are currently working on the the potentials in the flavor singlet and octet channels to obtain a complete picture of the BB interaction in the SU(3) limit.

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