Strangeness S=-2 baryon-baryon interactions from Lattice QCD

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HAL (Hadrons to Atomic nuclei from Lattice) QCD Collaboration

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Introduction

Introduction

BB interactions are crucial to investigate the nuclear phenomena

Once we obtain "proper" nuclear potentials, we apply them to the structure of (hyper-) nucleus.



How do we obtain the nuclear force?

Phenomenological descriptions

Traditional process to research the BB interaction / potential



The models would be highly ambiguous if experimental data are scarce!

Derivation of hadronic interaction from QCD

Start with the fundamental theory,QCD



HAL QCD method

Nambu-Bethe-Salpeter wave function



It satisfies the Helmholtz eq. in asymptotic region : $(p^2 + \nabla^2) \Psi(E, \vec{r}) = 0$

 $\Psi(E, \vec{r}) \simeq A \frac{\sin(pr + \delta(E))}{pr}$

Using the reduction formula.

C.-J.D.Lin et al., NPB619 (2001) 467.

$$\Psi^{\alpha}(E,\vec{r}) = \sqrt{Z_{H_1}} \sqrt{Z_{H_2}} \left(e^{i\vec{p}\cdot\vec{r}} + \int \frac{d^3q}{2E_q} \frac{T(q,p)}{4E_p(E_q - E_p - i\epsilon)} e^{i\vec{q}\cdot\vec{r}} \right)$$

Phase shift is defined as

$$S \equiv e^{i\delta}$$

NBS wave function has a same asymptotic form with quantum mechanics. (NBS wave function is characterized from phase shift)

Potential in HAL QCD method

We define potentials which satisfy Schrödinger equation

$$\left(p^{2}+\nabla^{2}\right)\Psi^{\alpha}(E,\vec{x})\equiv\int d^{3}y \,\underline{U}^{\alpha}_{\alpha}(\vec{x},\vec{y})\,\Psi^{\alpha}(E,\vec{y})$$

Energy independent potential

$$\begin{pmatrix} p^2 + \nabla^2 \end{pmatrix} \Psi^{\alpha}(E, \vec{x}) = K^{\alpha}(E, \vec{x}) K^{\alpha}(E, \vec{x}) \equiv \int dE' K^{\alpha}(E', \vec{x}) \int d^3 y \, \widetilde{\Psi}^{\alpha}(E', \vec{y}) \Psi^{\alpha}(E, \vec{y}) = \int d^3 y \Big[\int dE' K^{\alpha}(E', \vec{x}) \widetilde{\Psi}^{\alpha}(E', \vec{y}) \Big] \Psi^{\alpha}(E, \vec{y}) = \int d^3 y U^{\alpha}_{\alpha}(\vec{x}, \vec{y}) \Psi^{\alpha}(E, \vec{y})$$

We can define an energy independent potential but it is fully non-local.

This potential automatically reproduce the scattering phase shift

0

0.5

1

1.5

r [fm]

2.5

2

Time-dependent method

Start with the normalized four-point correlator.

$$R_{I}^{B_{1}B_{2}}(t,\vec{r}) = F_{B_{1}B_{2}}(t,\vec{r})e^{(m_{1}+m_{2})t}$$
Each wave functions satisfies
Schrödinger eq. with proper energy

$$= A_{0}\Psi(\vec{r},E_{0})e^{-(E_{0}-m_{1}-m_{2})t} + A_{1}\Psi(\vec{r},E_{1})e^{-(E_{1}-m_{1}-m_{2})t} + \cdots$$

$$\left(\frac{p_{0}^{2}}{2\mu} + \frac{\nabla^{2}}{2\mu}\right)\Psi(\vec{r},E_{0}) = \int U(\vec{r},\vec{r}\,')\Psi(\vec{r}\,',E_{0})d^{3}r'$$

$$\left(\frac{p_{1}^{2}}{2\mu} + \frac{\nabla^{2}}{2\mu}\right)\Psi(\vec{r},E_{1}) = \int U(\vec{r},\vec{r}\,')\Psi(\vec{r}\,',E_{1})d^{3}r'$$

$$\left(-\frac{\partial}{\partial t} + \frac{\nabla^{2}}{2\mu}\right)R_{I}^{B_{1}B_{2}}(t,\vec{r}) = \int U(\vec{r},\vec{r}\,')R_{I}^{B_{1}B_{2}}(t,\vec{r})d^{3}r'$$

A single state saturation is not required!!

Treatment of non-local potential

$$\left(-\frac{\partial}{\partial t}+\frac{\nabla^2}{2\mu}\right)R_I^{B_1B_2}(t,\vec{r})=\int U(\vec{r},\vec{r}')R_I^{B_1B_2}(t,\vec{r})d^3r'$$

Derivative (velocity) expansion of U is performed to deal with its nonlocality.

For the case of oct-oct system,

$$U(\vec{r},\vec{r}') = \begin{bmatrix} V_C(r) + S_{12}V_T(r) \end{bmatrix} + \begin{bmatrix} \vec{L} \cdot \vec{S}_s V_{LS}(r) + \vec{L} \cdot \vec{S}_a V_{ALS}(r) \end{bmatrix} + O(\nabla^2)$$

Leading order part

For the case of dec-oct and dec-dec system,

$$U(\vec{r},\vec{r}') = \begin{bmatrix} V_C(r) + S_{12}V_{T_1}(r) + S_{ii}V_{T_2}(r) + O(Spin op^3) \end{bmatrix} + O(\nabla^2)$$

Leading order part

$$\equiv \begin{bmatrix} V_C^{eff}(r) \end{bmatrix} + O(\nabla^2) \qquad ((\vec{r}\cdot\vec{S_1})^2 - \frac{\vec{r}^2}{3}\vec{S_1}^2 + (\vec{r}\cdot\vec{S_2})^2 - \frac{\vec{r}^2}{3}\vec{S_2}^2)V_{T^2}(r)$$

We consider the effective central potential which contains not only the genuine central potential but also tensor parts.

HAL QCD method

NBS wave function

$$\begin{split} \Psi(E,\vec{r})e^{-E(t-t_0)} &= \sum_{\vec{x}} \langle 0|B_i(t,\vec{x}+\vec{r})B_j(t,\vec{x})|E,t_0 \rangle \\ \text{E: Total energy of system} \\ &= \text{In asymptotic region }: (p^2+\nabla^2)\Psi(E,\vec{r})=0 \\ &= \text{In interaction region }: (p^2+\nabla^2)\Psi(E,\vec{r})=K(E,\vec{r}) \\ &= \text{Modified Schrödinger equation} \\ &= \left(-\frac{\partial}{\partial t} + \frac{\nabla^2}{2\mu}\right)R_I^{B_1B_2}(t,\vec{r}) = \int U(\vec{r},\vec{r}')R_I^{B_1B_2}(t,\vec{r})d^3r' \\ &= \left(-\frac{\partial}{\partial t} + \frac{\nabla^2}{2\mu}\right)R_I^{B_1B_2}(t,\vec{r}) = \int U(\vec{r},\vec{r}')R_I^{B_1B_2}(t,\vec{r})d^3r' \\ &= V_C(r) + S_{12}V_T(r) + \vec{L}\cdot\vec{S}_s V_{LS}(r) + \vec{L}\cdot\vec{S}_a V_{ALS}(r) + O(\nabla^2) \\ &= \left(-\frac{\partial}{\partial t} + \frac{\nabla^2}{2\mu}\right)R_I^{B_1B_2}(t,\vec{r})/R_I^{B_1B_2}(t,\vec{r}) \end{split}$$

HAL QCD method (coupled-channel)

NBS wave function

 $\Psi^{\alpha}(E_i,\vec{r})e^{-E_it} = \langle 0|(B_1B_2)^{\alpha}(\vec{r})|E_i\rangle$ $\Psi^{\beta}(E_{i},\vec{r})e^{-E_{i}t} = \langle 0|(B_{1}B_{2})^{\beta}(\vec{r})|E_{i}\rangle \qquad R_{E}^{B_{1}B_{2}}(t,\vec{r}) = \Psi_{B_{1}B_{2}}(\vec{r},E)e^{(-E_{1}+E_{2})}$

$$\int dr \tilde{\Psi}_{\beta}(E', \vec{r}) \Psi^{\gamma}(E, \vec{r}) = \delta(E'-E) \delta_{\beta}^{\gamma}$$

Leading order of velocity expansion and time-derivative method.

Modified coupled-channel Schrödinger equation

$$\begin{pmatrix} \left(-\frac{\partial}{\partial t}+\frac{\nabla^{2}}{2\mu_{\alpha}}\right)R_{E_{0}}^{\alpha}(t,\vec{r})\\ \left(-\frac{\partial}{\partial t}+\frac{\nabla^{2}}{2\mu_{\beta}}\right)R_{E_{0}}^{\beta}(t,\vec{r}) \end{pmatrix} = \begin{pmatrix} V_{\alpha}^{\alpha}(\vec{r}) & V_{\beta}^{\alpha}(\vec{r})\Delta_{\beta}^{\alpha}(t)\\ V_{\alpha}^{\beta}(\vec{r})\Delta_{\alpha}^{\beta}(t) & V_{\beta}^{\beta}(\vec{r}) \end{pmatrix} \begin{pmatrix} R_{E_{0}}^{\alpha}(t,\vec{r})\\ R_{E_{0}}^{\beta}(t,\vec{r}) \end{pmatrix} \\ \begin{pmatrix} \left(-\frac{\partial}{\partial t}+\frac{\mathbf{v}}{2\mu_{\beta}}\right)R_{E_{1}}^{\beta}(t,\vec{r}) & \Delta_{\beta}^{\alpha}=\frac{\exp\left(-\left(m_{\alpha_{1}}+m_{\alpha_{2}}\right)t\right)}{\exp\left(-\left(m_{\beta_{1}}+m_{\beta_{2}}\right)t\right)} \end{pmatrix} \begin{pmatrix} \vec{r})\Delta_{\beta}^{\alpha}(t)\\ V_{\beta}^{\beta}(\vec{r}) \end{pmatrix} \begin{pmatrix} R_{E_{1}}^{\alpha}(t,\vec{r})\\ R_{E_{1}}^{\beta}(t,\vec{r}) \end{pmatrix}$$

S.Aoki et al [HAL QCD Collab.] Proc. Jpn. Acad., Ser. B, 87 509 K.Sasaki et al [HAL QCD Collab.] PTEP no 11 (2015) 113B01

Considering two different energy eigen states

$$\begin{array}{l} \textbf{Potential} \\ \begin{pmatrix} V^{\alpha}_{\ \alpha}(\vec{r}) & V^{\alpha}_{\ \beta}(\vec{r})\Delta^{\alpha}_{\beta} \\ V^{\beta}_{\ \alpha}(\vec{r})\Delta^{\beta}_{\alpha} & V^{\beta}_{\ \beta}(\vec{r}) \end{pmatrix} = \begin{pmatrix} (\frac{\nabla^{2}}{2\mu_{\alpha}} - \frac{\partial}{\partial t})R^{\alpha}_{E0}(t,\vec{r}) & (\frac{\nabla^{2}}{2\mu_{\beta}} - \frac{\partial}{\partial t})R^{\alpha}_{E1}(t,\vec{r}) \\ (\frac{\nabla^{2}}{2\mu_{\alpha}} - \frac{\partial}{\partial t})R^{\beta}_{E0}(t,\vec{r}) & (\frac{\nabla^{2}}{2\mu_{\beta}} - \frac{\partial}{\partial t})R^{\beta}_{E1}(t,\vec{r}) \end{pmatrix} \begin{pmatrix} R^{\alpha}_{E0}(t,\vec{r}) & R^{\alpha}_{E1}(t,\vec{r}) \\ R^{\beta}_{E0}(t,\vec{r}) & R^{\beta}_{E1}(t,\vec{r}) \end{pmatrix}^{-1} \end{array}$$

S=-2 BB interaction

Interests of S=-2 multi-baryon system

H-dibaryon

The flavor singlet state with J=0 predicted by R.L. Jaffe.

- Strongly attractive color magnetic interaction.
- No quark Pauli principle for flavor singlet state.

Double- Λ hypernucleus

Conclusions of the "NAGARA Event"

K.Nakazawa and KEK-E176 & E373 Collaborators

 Λ −N attraction Λ − Λ weak attraction $m_{H} \ge 2m_{\Lambda} - 6.9$ MeV



Ξ hypernucleus

Conclusions of the "KISO Event"
 K.Nakazawa and KEK-E373 Collaborators

 Ξ –N attraction



SU(3) feature of BB interaction





Short range repulsion in BB interaction could be a result of Pauli principle and color-magnetic interaction for the quarks.

$$V_{OGE}^{CMI} \propto \frac{1}{m_{q1}m_{q2}} \langle \lambda_1 \cdot \lambda_2 \sigma_1 \cdot \sigma_2 \rangle f(r_{ij})$$

For the s-wave BB system, no repulsive core is predicted in flavor singlet state which is known as H-dibaryon channel.

	Flavor symmetric states			Flavor anti-symmetric states		
	27	8	1	<u>10</u>	10	8
Pauli	mixed	forbidden	allowed	mixed	forbidden	mixed
СМІ	repulsive	repulsive	attractive	repulsive	repulsive	repulsive
					Oka, Shimizu a	and Yazaki NPA464

B-B potentials in SU(3) limit

*m*π**= 469MeV**



Two-flavors

Three-flavors

- Quark Pauli principle can be seen at around short distances
 - No repulsive core in flavor singlet state
 - Strongest repulsion in flavor 8s state
- Possibility of bound H-dibaryon in flavor singlet channel.

Baryon-baryon system with S=-2



Relations between BB channels and SU(3) irreducible representations



Features of flavor singlet interaction is integrated into the S=-2 J^p=0⁺, I=0 system.

Numerical setup

2+1 flavor gauge configurations.

Iwasaki gauge action & O(a) improved Wilson quark action

- *a* = 0.086 [*fm*], a⁻¹ = 2.300 GeV.
- $96^3 \times 96$ lattice, L = 8.24 [fm].
- 414 confs x 28 sources x 4 rotations.

Flat wall source is considered to produce S-wave B-B state.



I=0 channel

Lists of channels



$\Sigma\Sigma$ (I=2) ¹S_o channel



• $\Sigma\Sigma$ (I=2) potential belongs to the 27plet in flavor SU(3) limit.

The potential has an attractive pocket and repulsive core.

In this time range, potentials are qualitatively similar.

Potential saturation is achieved at after t=11??

$$N\Xi$$
 (I=0) ${}^{3}S_{1}-{}^{3}D_{1}$ channel



NE (I=0,spin triplet) potential belongs to the 8plet in flavor SU(3) limit.

The potential has an attractive pocket and repulsive core.

Tensor potential is weaker than the phenomenological NN tensor potential.

$N\Xi$ (I=0) ${}^{3}S_{1}$ channel



Effective NE (I=0,spin triplet) central potential is plotted.

From this figure, we find that the tensor potential effects are small.

Phase shifts are same within the error bars.

NΞ, $\Lambda \Sigma$ (I=1) ¹S₀ channel







Diagonal elements are repulsive in whole range.

- Diagonal NE potential is strongly repulsive.
 - It means that the NE potential is strongly depend on the channel.
- Potentials are not saturated in this time range.

NΞ, $\Lambda \Sigma$, $\Sigma \Sigma$ (I=1) ³S₁ channel



ΛΛ, ΝΞ, ΣΣ (I=0) 1 S_o channel



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Comparison of potential matrices

Transformation of potentials

from the particle basis to the SU(3) irreducible representation (irrep) basis.

 $\begin{aligned} & \left(\begin{vmatrix} 1 \\ 8 \\ 27 \end{matrix} \right) = U \begin{pmatrix} |\Lambda\Lambda\rangle \\ N\Xi \\ |\Sigma\Sigma \rangle \end{pmatrix}, \quad U \begin{pmatrix} V^{\Lambda\Lambda} & V^{\Lambda\Lambda}_{N\Xi} & V^{\Lambda\Lambda}_{\Sigma\Sigma} \\ V^{N\Xi}_{\Lambda\Lambda} & V^{N\Xi} & V^{N\Xi}_{\Sigma\Sigma} \\ V^{\Sigma\Sigma}_{\Lambda\Lambda} & V^{\Sigma\Sigma}_{N\Xi} & V^{\Sigma\Sigma} \end{pmatrix} U^{t} \rightarrow \begin{pmatrix} V_{1} & V_{1} & V_{1} & V_{1} \\ V_{27} & V_{27} & V^{27} & V^{27} \end{pmatrix} \\ & \left(\begin{pmatrix} 1\Sigma\Sigma \\ N\Xi \\ N\Xi \end{pmatrix} \\ (\Lambda\Lambda) \end{pmatrix} = \frac{1}{\sqrt{40}} \begin{pmatrix} -1 & -\sqrt{24} & \sqrt{15} \\ \sqrt{12} & \sqrt{8} & \sqrt{20} \\ \sqrt{27} & -\sqrt{8} & -\sqrt{5} \end{pmatrix} \begin{pmatrix} 127 \\ 18_{s} \\ 1 \end{pmatrix} \end{aligned}$

In the SU(3) irreducible representation basis, the potential matrix should be diagonal in the SU(3) symmetric configuration.

Off-diagonal part of the potential matrix in the SU(3) irrep basis would be an effective measure of the SU(3) breaking effect.

1, 8, 27plet (I=0) ¹S_o channel





H-dibaryon (flavor SU(3) situations)



NPL : PRL106(2011)162001

The bound H-dibaryon state in heavy pion region.
 Potential in flavor singlet channel is getting more attractive as decreasing quark masses





Works on H-dibaryon state



Effective two channel potential

Original coupled channel equation

$$\begin{pmatrix} \left(E^{\Lambda\Lambda} - H_{0}^{\Lambda\Lambda}\right)R^{\Lambda\Lambda}(\vec{r},t) \\ \left(E^{\Xi N} - H_{0}^{\Xi N}\right)R^{\Xi N}(\vec{r},t) \\ \left(E^{\Sigma\Sigma} - H_{0}^{\Sigma\Sigma}\right)R^{\Sigma\Sigma}(\vec{r},t) \end{pmatrix} = \begin{pmatrix} V^{\Lambda\Lambda}_{\Lambda\Lambda}(\vec{r}) & V^{\Lambda\Lambda}_{\Xi N}(\vec{r}) & V^{\Lambda\Lambda}_{\Sigma\Sigma}(\vec{r}) \\ V^{\Xi N}_{\Lambda\Lambda}(\vec{r}) & V^{\Xi N}_{\Xi N}(\vec{r}) & V^{\Xi N}_{\Sigma\Sigma}(\vec{r}) \\ V^{\Sigma\Sigma}_{\Lambda\Lambda}(\vec{r}) & V^{\Sigma\Sigma}_{\Xi N}(\vec{r}) & V^{\Sigma\Sigma}_{\Sigma\Sigma}(\vec{r}) \end{pmatrix} \begin{pmatrix} R^{\Lambda\Lambda}(\vec{r},t) \\ R^{\Xi N}(\vec{r},t) \\ R^{\Sigma\Sigma}(\vec{r},t) \end{pmatrix}$$

Truncation of $\Sigma\Sigma$ channel

Reduced coupled channel equation

$$\begin{pmatrix} (E^{\Lambda\Lambda} - H_0^{\Lambda\Lambda}) R^{\Lambda\Lambda}(\vec{r},t) \\ (E^{\Xi N} - H_0^{\Xi N}) R^{\Xi N}(\vec{r},t) \end{pmatrix} = \begin{pmatrix} \overline{V_{\Lambda\Lambda}^{\Lambda\Lambda}}(\vec{r}) & \overline{V_{\Xi N}^{\Lambda\Lambda}}(\vec{r}) \\ \overline{V_{\Lambda\Lambda}^{\Xi N}}(\vec{r}) & \overline{V_{\Xi N}^{\Xi N}}(\vec{r}) \end{pmatrix} \begin{pmatrix} R^{\Lambda\Lambda}(\vec{r},t) \\ R^{\Xi N}(\vec{r},t) \end{pmatrix}$$

Effective $\Lambda\Lambda$ -N Ξ potential

The same scattering phase shift would be expected in a low energy region.

Non-locality (energy dependence, higher derivative contribution)

of potential matrix could be enhanced.

$\Lambda\Lambda$, $N\Xi$ (I=0) ¹S_o potential (Effective 2ch calc.)

N_f = 2+1 full QCD with L = 8fm, $m\pi = 146 \text{ MeV}$



Potential calculated by only using ΛΛ and ΝΞ channels.

Long range part of potential is almost stable against the time slice.

- Short range part of NE potential changes as time t goes.
- •ΛΛ–NΞ transition potential is quite small in r > 0.7fm region



Preliminary!

$\Lambda\Lambda$ and $N\Xi$ phase shift and inelasticity \square

N_f = 2+1 full QCD with L = 8fm, $m\pi = 146 \text{ MeV}$

Preliminary!



AA and NE phase shift is calculated by using 2ch effective potential.
A sharp resonance is found just below the NE threshold.
Inelasticity is small.



Breit-Wigner mass and width

N_f = 2+1 full QCD with L = 8fm, $m\pi = 146 \text{ MeV}$

Preliminary!

50



Summary and outlook

We have investigated coupled channel baryonic interactions from lattice QCD.

We find that

• Potential in $\Lambda\Lambda$ ¹S₀ channel is weakly attractive.

NE potential is largely depends on its channel.

Potential in flavor singlet ¹S₀ channel is strongly attractive.

We have studied dibaryon candidate states

- H-dibaryon channel
 - We perform $\Lambda\Lambda$ -NE coupled channel calculation.
 - Sharp resonance is found just below the NE threshold (Time slice saturation is not achieved yet.)

We continue to study it by using higher statistical data.