# Baryon interactions from lattice QCD with physical masses -- S=-2 sector --

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#### for HAL QCD Collaboration



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### Introduction

#### **BB** interactions are crucial to investigate (hyper-)nuclear structures



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### 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<sup>p</sup>=0<sup>+</sup>, I=0 system.

### Keys to understand H-dibaryon

#### A strongly bound state predicted by Jaffe in 1977 using MIT bag model.

H-dibaryon state is

- SU(3) flavor singlet [uuddss], strangeness S=-2.
- spin and isospin equals to zero, and J<sup>P</sup>= 0<sup>+</sup>

Strongly attractive interaction is expected in flavor singlet channel.

Short range one-gluon exchange contributions

**Strongly attractive Color Magnetic Interaction** 

Symmetry of two-baryon system (Pauli principle)

Flavor singlet channel is free from Pauli blocking effect

	27	8	1	<u>10</u>	10	8
Pauli	mixed	forbidden	allowed	mixed	forbidden	mixed
CMI	repulsive	repulsive	attractive	repulsive	repulsive	repulsive

#### SU(3) breaking effects

- Threshold separation
- Changes of interactions



Oka, Shimizu and Yazaki NPA464 (1987)

Non-trivial contributions

### Hunting for H-dibaryon in SU(3) limit

Strongly attractive interaction is expected in flavor singlet channel.



 Strongly attractive potential was found in the flavor singlet channel.
 Bound state was found in this mass range with SU(3) symmetry.

What happens at the physical point?



### Works on H-dibaryon state



### 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

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

### Numerical setup

2+1 flavor gauge configurations.

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

- *a* = 0.086 [*fm*], a<sup>-1</sup> = 2.300 GeV.
- 96<sup>3</sup>x96 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.



## $\Lambda\Lambda$ , $N\Xi$ (I=0) <sup>1</sup>S<sub>o</sub> potential (2ch calc.)

#### N<sub>f</sub> = 2+1 full QCD with L = 8fm, $m\pi = 146 \text{ MeV}$



#### Potential calculated by only using ΛΛ and NΞ 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 NE phase shift and inelasticity [199]

**Preliminary!** 

#### N<sub>f</sub> = 2+1 full QCD with L = 8fm, $m\pi = 146 \text{ MeV}$

 $\Lambda\Lambda$  phase shift Inelasticity 180 0.9 180 150 150 0.95 0.8 120 120 0.9 δ [deg] F 90 0.85 90 0.8 0.7 60 0.75 60 30 0.7 0.6 0.65 41.7 41.8 41.9 42 42.1 41.5 41.6 30 0.6 42.2 42.4 42.8 42 42.6 43 0.5 0 30 0 10 20 40 50 30 10 20 40 50 0 E [MeV] E [MeV] 200 C 1 V<sub>NΞ-NΞ</sub>(t=09) ⊢ 1  $V_{N\equiv-N\equiv}(fit)$  $V_{\Lambda\Lambda-N\Xi}(t=09)$ 150  $V_{\Lambda\Lambda-N=}(fit)$ -30 11  $V_{\Lambda\Lambda-\Lambda\Lambda}(t=09)$ 1 -30  $V_{\Lambda\Lambda-\Lambda\Lambda}(fit)$ 100 V(r)[MeV] -60 -60 δ [deg] 50 -90 -90 42.4 42.6 42.8 42 42.2 0 N<del>E</del> phase shift -120 -50 0 0.5 1.5 2 2.5 3 1 0 10 20 30 40 50 r[fm]

### $\Lambda\Lambda$ and $N\Xi$ phase shift and inelasticity [t=10]

#### N<sub>f</sub> = 2+1 full QCD with L = 8fm, $m\pi = 146 \text{ MeV}$

**Preliminary!** 



### $\Lambda\Lambda$ and $N\Xi$ phase shift and inelasticity [t=1]

#### N<sub>f</sub> = 2+1 full QCD with L = 8fm, $m\pi = 146 \text{ MeV}$

**Preliminary!** 



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

#### N<sub>f</sub> = 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<sub>f</sub> = 2+1 full QCD with L = 8fm, $m\pi = 146 \text{ MeV}$

#### **Preliminary!**



$$\delta(E) = \delta_B - \arctan\left(\frac{\Gamma/2}{E - E_r}\right)$$
thus

$$\frac{d \,\delta(E)}{d E} = \frac{\Gamma/2}{(E - E_r)^2 + (\Gamma/2)^2}$$

 Fitting the time delay of ΛΛ scattering by the Breit-Wigner type finction,

Resonance enargy and width

#### t=09

$$E_{R} - E_{\Lambda\Lambda} = 41.894 \pm 0.039 [MeV]$$
  

$$\Gamma = 0.099 \pm 0.059 [MeV]$$

$$E_{R} - E_{\Lambda\Lambda} = 41.917 \pm 0.056 [MeV]$$
  

$$\Gamma = 0.077 \pm 0.021 [MeV]$$

#### t=11

$$E_{R} - E_{\Lambda\Lambda} = 41.927 \pm 0.105 [MeV]$$
  

$$\Gamma = 0.050 \pm 0.053 [MeV]$$

### Invariant mass spectrum of $\Lambda\Lambda$ channel





#### Sharp peak below NE threshold

Direct comparison with our simulation results and experimental data will be performed in near future?

### Summary

H-dibaryon state is investigated using 414confs x 28src x 4rot.

•We perform  $\Lambda\Lambda$ -NE coupled channel calculation.

Sharp resonance is found just below the NE threshold.

Resonance position and width from Breit-Wigner type fit



We continue to study it by using higher statistical data.

# Backup slides

# HAL QCD method

### Derivation of hadronic interaction from QCD



### Nambu-Bethe-Salpeter wave function



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

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

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)

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### Potential in HAL QCD method

We define potentials which satisfy Schrödinger equation

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

Energy independent potential

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

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

This potential automatically reproduce the scattering phase shift

2.5

r [fm]

### 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 satisfy  
Schrödinger eq. with proper energy  

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

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

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

$$E_{n}-m_{1}-m_{2}\approx\frac{p_{n}^{2}}{2\mu}$$

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

#### A single state saturation is not required!!

### **BB** interaction from NBS wave function

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

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.

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# ΛΛ, ΝΞ, ΣΣ (I=0) $^{1}S_{0}$ channel (t=11)



Allediagonaliteleaventelaavelyalangeutsive compating to the tentialiptestrangly repulsive.

#### •Diagonal NE potential is more attractive than the $\Lambda\Lambda$ potential.

We need more statistics to discuss physical observables through this potential.



#### Potential of flavor singlet channel does not have a repulsive core

Potential of flavor octet channel is strongly repulsive which reflects Pauli effect.
 Off-diagonal potentials are visible only in r<1fm region.</li>

