### Omega-Omega interaction from 2+1 flavor QCD

### Masanori Yamada (University of Tsukuba) for HAL QCD Collaboration

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# Outline

- Introduction
- Formulation
  - Construction of the potential [HAL QCD method]
  - Symmetry of Omega-Omega system
- Lattice QCD Simulation results
  - Potential
  - phase shift & binding energy
- Conclusion

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Introduction

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#### **Decuplet Baryon**

#### Motivation



- $\cdot$  Omega baryon is stable in QCD
- $\cdot$  It's a first target of Decuplet-Decuplet

interactions at HAL QCD method.

• There have been different model calculations in the J=0 channel

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interaction energy

 $\Delta M_{\Omega\Omega} = E_{\Omega\Omega} - 2M_\Omega = -166 {
m MeV}$ 

(SU(3) Chiral Quark Model)

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$$\Delta M_{\Omega\Omega} = E_{\Omega\Omega} - 2M_{\Omega} = 43 \pm 18 {
m MeV}$$

(Quark Disloc./Color-screen Model)

[Z.Y.Zhang et al. Phys.Rev.C .61, 065204] [F.Wang et al. Phys Rev C. 51, 3411]

#### There have been different model calculations in the J=0 channel



### Report from another group (Lattice QCD simulation)

Lüscher's method [Lüscher CMP105(86)153, NPB354(91)531]

Buchoff et al. : L=3fm Ω=1628[MeV]

 $J=0: weak repulsion a= -0.16 \pm 0.22 \text{ fm}$  [arXiv:1201.3596]

J=2 : strong repulsion

J.Wasem @Lattice2012

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no definite conclusion, attraction or repulsion

determine a nature of J=0 Omega-Omega interaction, attractive or repulsive

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A. Nambu-Bethe-Salpeter(NBS) wave function

$$\psi_k(r) \equiv \langle 0 | \Omega(r) \Omega(0) | \overline{\Omega}(k) \overline{\Omega}(-k); in \rangle$$
  
Because

NBS wave has the same asymptotic form of the scattering wave in quantum mechanics.

Wave function↔phase shift↔S-matrix

$$\psi_k(r) \simeq e^{i\delta(k)} \frac{\sin(kr - \frac{l\pi}{2} + \delta(k))}{kr}$$

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

Energy independent potential U(r,r') is defined from NBS wave function.

because of

$$(rac{k^2}{m}+rac{1}{m}
abla^2)\psi_k(r)=\int d^3r' U(r,r')\psi_k(x')$$
 This potential reproduces the phase shift faithfully

we can extract a potential (interaction kernel) which is defined through the NBS wave function which gives the correct scattering phase shift at asymptotic state.

# Symmetry of $\Omega$ - $\Omega$

 $\boldsymbol{\Omega}$  operator is defined as

$$\Omega_{\alpha,\mathbf{k}} \equiv \varepsilon^{abc} s^a (C\gamma_{\mathbf{k}}) s^b s^c_{\alpha}$$

blue is spin1 index, red is spin $\frac{1}{2}$  index

We treat spin 3/2 made from spin 1 and spin 1/2 linear combination by using highest weight construction

•one 
$$\Omega$$
 case(spin $\frac{3}{2}$ )  
spin $\frac{1}{2} \otimes spin1 = spin \frac{3}{2} \oplus spin \frac{1}{2}$ 

•consider two  $\Omega$  case (  $\Omega$ - $\Omega$  interaction)

$${\operatorname{spin}} rac{3}{2} \otimes {\operatorname{spin}} rac{3}{2} = {\operatorname{spin3}} \oplus {\operatorname{spin2}} \oplus {\operatorname{spin1}} \oplus {\operatorname{spin0}}$$

# Symmetry of $\Omega$ - $\Omega$

### Conserved quantity $J, J_z, P$

- parity  $P = (-1)^L$
- $\cdot$  quantum spin  $(-1)^{S+1}$

fermionic condition

$$(-1)^L \times (-1)^{S+1} = -1$$

 $\psi_1\psi_2=-\psi_2\psi_1$ 

### Which L ,S is allowed at $J^P$

	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=2 L=4	S=1 L=1 , S=3 L=1 , S=1 L=3 , 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=2 L=6	S=3 L=1 , S=1 L=3 , S=3 L=3 , S=1 L=5 , S=3 L=5 , S=3 L=7

	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=2 L=4	S=1 L=1 , S=3 L=1 , S=1 L=3 , 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=2 L=6	S=3 L=1 , S=1 L=3 , S=3 L=3 , S=1 L=5 , S=3 L=5 , S=3 L=7

### at sink

	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=2 L=4	S=1 L=1 , S=3 L=1 , S=1 L=3 , 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
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### $L=0 \leftarrow We$ use wall source

at sink

	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=2 L=4	S=1 L=1 , S=3 L=1 , S=1 L=3 , S=3 L=3 , S=3 L=5
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 $L=0 \iff We \text{ use wall source}$  $J=0 \iff S=0$ 

at sink

	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
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 $L=0 \iff We \text{ use wall source}$  $J=0 \iff S=0$ 

at sink

We can extract S=0 L=0,S=2 L=2

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## Lattice set up



### We used 2 sets (Small , Large)

	Small	Large	
Lattice volume	L = 1.950(30) [fm]	L = 2.902(42) [fm]	
	$\rm K_s = 0.13710 \ \rm K_{ud} = 0.13760$	$\rm K_{s} = 0.13640 \ \rm K_{ud} = 0.13700$	
nopping parameters	$M_\Omega = 2104(8) [{\rm MeV}] \ M_\pi = 875(1) [{\rm MeV}]$	$M_\Omega = 1966(6) [{\rm MeV}] \; M_\pi = 701(5) [{\rm MeV}]$	
β	β=1.83	β=1.90	
lattice spaceing	a = 0.1219(19)[fm]	a = 0.0907(13) [fm]	

•RG improved gauge action & O(a) improved Wilson quark action

•flat wall source(P=0)

CP-PACS/JLQCD Collaborations: [T. Ishikawa, et al., Phys. Rev. D78 (2008) 011502(R)]

Experiment value of  $\Omega$  mass is [1672 MeV]

PACS-CS Collaboration: [S. Aoki, et al., Phys. Rev. D79 (2009) 034503 ]

### Results of the small volume

### $\Omega$ - $\Omega$ Potential (Small volume) @lattice2013



t is relative time between source and sink

 $t \equiv t_1 - t_0$ 

 $C_{\Omega\Omega}(\overrightarrow{x},\overrightarrow{y},t_1,t_0)\equiv \langle 0|\,\Omega(\overrightarrow{x},t_1)\Omega(\overrightarrow{y},t_1)\overline{\Omega}(0,t_0)\overline{\Omega}(0,t_0)\,|0
angle$ 

#### $\Omega$ - $\Omega$ Potential (Small volume) update



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### $\Omega$ - $\Omega$ Potential (Small volume) update



### $\Omega$ - $\Omega$ Potential (Small volume) update



### Phase shift & Binding energy (Small volume)

#### We found bound states, but binding energies are very small.

	Binding ene	ergy	Scattering length
t=7	-0.13 <u>+</u> 0.28	[MeV]	$-13.32$ $\pm$ 20.34 [fm]
t=8	-0.47 <u>+</u> 1.00	[MeV]	–7.07 ± 7.70 [fm]
t=9	-4.64 <u>+</u> 4.73	[MeV]	–2.71 ± 1.32 [fm]
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## Results of the large volume

### $\Omega$ - $\Omega$ Potential (Large volume)



### $\Omega$ - $\Omega$ Potential (Large volume)



repulsive core grow

### $\Omega$ - $\Omega$ Potential (Large volume)



## Phase shift (Large volume)

Phase Shift



## Phase shift (Large volume)

Phase Shift



### Large vs Small



stronger repulsive core at larger volume

attraction	bound	repulsion	
results	strong attraction	weak repulsion(?)	
ground sate saturation	not needed	needed	
Lattice volume	2.9[fm]	3.9[fm]	
fermion mass	heavy(π=701)	light(π=390)	
method	potential (our work)	finite volume (Buchoff et al.)	







## Conclusion

- We extended HAL method to decoupletdecouplet system.
- We showed small volume and large volume results.
- J=0 Omega-Omega interaction is strongly attractive but we can not decide whether the bound state exists or not due to large errors.

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# Back up slide

#### comparison Luscher method, HAL method(phase shift $\pi$ - $\pi$ in I=2 channel)



The result of phase shift have been found to agree well between the two methods!

It's difficult to compare these methods without calculating finite volume method at large t and more statics!

# Back up slide

Mass dependence (N-N interaction)



Fig. 5. The central potentials in the  ${}^{1}S_{0}$  channel for three different quark masses.

We expect  $\Omega$ - $\Omega$  is similar to N-N case

[Sinya Aoki et al. Prog. Theor. Phys. 123,89]

### Large vs Small



because growth of repulsive core

	boundary	effect
$V(\vec{r})$		$\tilde{V}(\vec{r}) = \sum_{\vec{n} \in \mathbb{Z}^3} V(\vec{r} + L\vec{n})$

### S=0 $\Leftarrow$ a special circumstance in $\Omega$ - $\Omega$ system

- flavor is completely symmetry
- wall source

#### source operator

$$\overline{\Omega} = \varepsilon^{abc} (\gamma_k C)_{\beta\gamma} \overline{s}^a_{\alpha} \overline{s}^b_{\beta} \overline{s}^c_{\gamma}$$

a,b,c: color index  $\alpha,\beta,\gamma$ : spin index

highest state in  $\Omega$ - $\Omega$  (spin3)

$$\overline{s}^a_{rac{1}{2}}(x)\overline{s}^b_{rac{1}{2}}(x)\overline{s}^c_{rac{1}{2}}(x)\overline{s}^a_{rac{1}{2}}(y)\overline{s}^b_{rac{1}{2}}(y)\overline{s}^c_{rac{1}{2}}(y)$$
 For simply neglect  $\epsilon$  , YC

We can make all state using lowering operator

spin3⇒spin2⇒spin1⇒spin0

For example one term of spin2 state

$$\overline{s}^a_{\frac{1}{2}}(x)\overline{s}^b_{\frac{1}{2}}(x)\overline{s}^c_{\frac{1}{2}}(x)\overline{s}^a_{\frac{1}{2}}(y)\overline{s}^b_{\frac{1}{2}}(y)\overline{s}^c_{-\frac{1}{2}}(y)$$

spin2 term is written by linear combination of these terms.

$$\overline{s}^a_{\frac{1}{2}}(x)\overline{s}^b_{\frac{1}{2}}(x)\overline{s}^c_{\frac{1}{2}}(x)\overline{s}^a_{\frac{1}{2}}(y)\overline{s}^b_{\frac{1}{2}}(y)\overline{s}^c_{-\frac{1}{2}}(y)$$

= 0 Spin2 state should be 0

$$\overline{s}^a_{\frac{1}{2}}(x)\overline{s}^b_{\frac{1}{2}}(x)\overline{s}^c_{\frac{1}{2}}(x)\overline{s}^a_{\frac{1}{2}}(y)\overline{s}^b_{\frac{1}{2}}(y)\overline{s}^c_{-\frac{1}{2}}(y)$$

$$\begin{aligned} \text{wall source} \Rightarrow \quad \left(\sum_{x} \overline{s_{\frac{1}{2}}^{a}(x)}\right) \left(\sum_{x'} \overline{s_{\frac{1}{2}}^{b}(x')}\right) \left(\sum_{x''} \overline{s_{\frac{1}{2}}^{c}(x'')}\right) \left(\sum_{y} \overline{s_{\frac{1}{2}}^{a}(y)}\right) \left(\sum_{y'} \overline{s_{\frac{1}{2}}^{b}(y')}\right) \left(\sum_{y''} \overline{s_{-\frac{1}{2}}^{c}(y'')}\right) \\ = -\left(\sum_{y} \overline{s_{\frac{1}{2}}^{a}(y)}\right) \left(\sum_{x'} \overline{s_{\frac{1}{2}}^{b}(x')}\right) \left(\sum_{x''} \overline{s_{\frac{1}{2}}^{c}(x'')}\right) \left(\sum_{x''} \overline{s_{\frac{1}{2}}^{c}(x')}\right) \left(\sum_{x''} \overline{s_{\frac{1}{2}}^{b}(y)}\right) \left(\sum_{y''} \overline{s_{\frac{1}{2}}^{b}(y')}\right) \left(\sum_{y''} \overline{s_{-\frac{1}{2}}^{c}(y'')}\right) \\ = -\left(\sum_{x} \overline{s_{\frac{1}{2}}^{a}(x)}\right) \left(\sum_{x'} \overline{s_{\frac{1}{2}}^{b}(x')}\right) \left(\sum_{x''} \overline{s_{\frac{1}{2}}^{c}(x'')}\right) \left(\sum_{y''} \overline{s_{\frac{1}{2}}^{a}(y)}\right) \left(\sum_{y''} \overline{s_{\frac{1}{2}}^{b}(y')}\right) \left(\sum_{y''} \overline{s_{-\frac{1}{2}}^{c}(y'')}\right) \\ \end{aligned}$$

= 0 Spin2 state should be 0

$$\overline{s}^a_{\frac{1}{2}}(x)\overline{s}^b_{\frac{1}{2}}(x)\overline{s}^c_{\frac{1}{2}}(x)\overline{s}^a_{\frac{1}{2}}(y)\overline{s}^b_{\frac{1}{2}}(y)\overline{s}^c_{-\frac{1}{2}}(y)$$

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## Spin0 remain

Existence of energy independent nonlocal potential

We assume linear independence of NBS wave function There is a dual bases

$$\int d^3r \widetilde{\psi}_{k'}(r) \psi_k(r) = (2\pi)^3 \delta^3(k'-k)$$

We define K

$$egin{aligned} K_k(r) &\equiv ( egin{aligned} & ( 
abla^2 + k^2 ) \psi_k(r) \ & = \int rac{d^3 k'}{(2\pi)^3} K_{k'}(r) \int d^3 r' \widetilde{\psi}_{k'}(r') \psi_k(r') \ & = \int d^3 r' \left\{ \int rac{d^3 k'}{(2\pi)^3} K_{k'}(r) \widetilde{\psi}_{k'}(r') 
ight\} \psi_k(r') \end{aligned}$$

If we define

$$U(r,r')\equiv rac{1}{m}{\intrac{d^3k'}{(2\pi)^3}K_{k'}(r)\widetilde{\psi}_{k'}(r')}$$

Then we have

$$(rac{k^2}{m}+rac{1}{m}
abla^2)\psi_k(r)=\int d^3r' U(r,r')\psi_k(r')$$

## Construction of the potential

Extraction of the NBS wave from Lattice QCD

 $C_{\Omega\Omega}(\overrightarrow{x},\overrightarrow{y},t,t_0) \equiv \langle 0 | \Omega(\overrightarrow{x},t) \Omega(\overrightarrow{y},t) \overline{\Omega}(0,t_0) \overline{\Omega}(0,t_0) | 0 \rangle$ Image  $=\sum \left<0\right| \Omega(\overrightarrow{x},t) \Omega(\overrightarrow{y},t) \left|n\right> e^{-E_n(t-t_0)} \left< n\right| \overline{\Omega}(0,t_0) \overline{\Omega}(0,t_0) \left|0\right>$  $\psi_{k_n}(\overrightarrow{x}-\overrightarrow{y},n)$ source  $=\sum A_n\psi_{k_n}(x-y,n)e^{-E_n(t-t_0)}+\cdots$  $t_0$ **Excited states are suppressed** exponentially at large  $t - t_0$ inelastic contributions We can get the NBS wave at ground state

## Construction of the potential

Extraction of the NBS wave from Lattice QCD



## Construction of the potential

Extraction of the NBS wave from Lattice QCD



Time dependent Schrodinger-type equation

$$(rac{1}{4m}rac{\partial^2}{\partial t^2}+rac{1}{m}
abla^2-rac{\partial}{\partial t})R=\int dr' U(r,r')R$$

Time dependent Schrodinger-type equation

R-correlator is defined as

$$R \equiv \frac{\Psi(r,t)}{e^{-2mt}} = \sum_{n} \phi_n(r) e^{-W_n t}$$

$$W_n\equiv 2\sqrt{m^2+ec{k}rac{2}{n}}-2m$$

$$(rac{1}{4m}rac{\partial^2}{\partial t^2}+rac{1}{m}
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Time dependent Schrodinger-type equation



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[N.Ishii et al., PLB712(2012)437.]



 $(rac{1}{4m}rac{\partial^2}{\partial t^2}+rac{1}{m}
abla^2-rac{\partial}{\partial t})R=\int dr' U(r,r')R$ 

2 NBS wave function satisfies Schorodinger eq.

$$(rac{k^2}{m}+rac{1}{m} 
abla^2)\psi_k(r)=\int d^3r' U(r,r')\psi_k(x')$$

# Time depend method

Time dependent Schrodinger-like equation

$$(rac{1}{4m}rac{\partial^2}{\partial t^2}+rac{1}{m}
abla^2-rac{\partial}{\partial t})R=\int dr' U(r,r')R$$

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[N.Ishii et al., PLB712(2012)437.]

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abla^2-rac{\partial}{\partial t})R=\int dr' U(r,r')R$$

We can calculate energy independent non-local potential without relying on the ground state saturation!

## fit function dependence

#### fit function

Gauss + Yukawa  $f(r) = b_1 e^{-b_2 r^2} + b_3 (1 - e^{-b_4 r^2}) \frac{e^{-b_5 r}}{r}$  $\lim_{r \to 0} f(r) = b_1$ Gauss + Yukawa^2  $f(r) = b_1 e^{-b_2 r^2} + b_3 (1 - e^{-b_4 r^2})^2 (\frac{e^{-b_5 r}}{r})^2$  $\lim_{r \to 0} f(r) = b_1$ 

2Gauss + Yukawa^2

$$f(r) = b_1 e^{-b_2 r^2} + b_3 e^{-b_4 r^2} + b_5 (1 - e^{-b_6 r^2})^2 (\frac{e^{-b_7 r}}{r})^2$$

 $\lim_{r \to 0} f(r) = b_1 + b_3$