Interactions between Decuplet Baryons from Lattice QCD

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HAL(Hadrons to Atomic nuclei from Lattice) QCD Collaboration

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SG, K.Sasaki + (HAL QCD Coll.), Phys.Rev.Lett. 120 (2018) 212001

 $\Omega\Omega$ interaction with J=0 at almost physical point



"di-Omega" Most Strange dibaryon

Introduction



Dibaryon = Bound (or resonance) two baryon states

Introduction: SU(3) classification for Baryon (B=1)



All octet baryons are stable under strong decay

Introduction: SU(3) classification for Baryon (B=1)



All octet baryons are stable under strong decay In Decuplet baryons, only Ω -baryon is stable



Introduction: SU(3) classification for Dibaryon

candidates (B=2)



1) octet-octet system $8 \otimes 8 = 27 \oplus 8_{s} \oplus 1 \oplus \overline{10} \oplus 10 \oplus 8_{a}$ Deuteron(J=1) 2) decuplet-octet system $10 \otimes 8 = 35 \oplus 8 \oplus 10 \oplus 27$ Goldman et al (1987) 3) decuplet-decuplet system $10 \otimes 10 = 28 \oplus 27 \oplus 35 \oplus \overline{10}$ by CELSIUS/WASA, 2009

ΩΩ system (J=0)

Zhang et al(1997)

ΔΔ system (J=3)

Dyson, Xuong (1964) Oka, Yazaki(1980)

Previous model works on $\Omega\Omega$ in J=0

SU(3) chiral quark model Z.Y. Zhang et al(1997)

$$\Delta M_{\Omega\Omega} = -166 \mathrm{MeV}$$

 $\Delta M_{\Omega\Omega} \equiv E_{\Omega\Omega} - 2M_{\Omega}$

Quark Disloc/Color-screen model F. Wang et al(1992)

$$\Delta M_{\Omega\Omega} = 43 \pm 18 \mathrm{MeV}$$

- Bound/unbound problem highly depends on models and their parameters.
- To clarify $\Omega\Omega$ interaction in our world, first-principle calculations are needed.

Baryon-Baryon interaction from lattice QCD -HAL method-

Aoki, Hatsuda, Ishii, PTP123, 89 (2010)

c.f. anothor method: Luscher's direct method

Nambu-Bethe-Salpeter (NBS) w.f. $\Psi_n(\vec{r}) e^{-E_n t}$

$$= \sum_{\vec{x}} \langle 0 | B_1(t, \vec{r} + \vec{x}) B_2(t, \vec{x}) | E_n \rangle$$

Local operators B_1 and B_2 for Ω baryon

()

S

SS

S

S

S

$$\Omega_{\alpha,k}(x) = \epsilon^{abc} \left[s_a^T(x) C \gamma_k s_b(x) \right] s_{c,\alpha}(x)$$

In asymptotic region (r>>R)
Helmholtz eq. is satisfied:
$$\left(\nabla^2 + k^2 \right) \Psi(\vec{r}) = 0_{\epsilon^0} \Psi(\vec{r}) \simeq A \frac{\sin(kr - l\pi/2 + \delta(k))}{kr}$$

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$$\Omega_{\alpha,k}(x) = \epsilon^{abc} \left[s_a^T(x) C \gamma_k s_b(x) \right] s_{c,\alpha}(x)$$

In interacting region,

Schroedinger type equation is satisfied

$$\left(\vec{p}_{n}^{2} + \nabla^{2}\right)\Psi_{n}\left(\vec{r}\right) = 2\mu \int d\vec{r'} U(\vec{r}, \vec{r'})\Psi_{n}(\vec{r'})$$

Aoki, Hatsuda, Ishii, PTP123, 89 (2010) Nonlocal potential U(r,r')

$$\left(\vec{p}_n^2 + \nabla^2\right)\Psi_n\left(\vec{r}\right) = 2\mu \int d\vec{r'} U(\vec{r},\vec{r'})\Psi_n(\vec{r'})$$

- The potential is energy-independent but non-local.
- The local leading potential can be obtained by its derivative expansion (c.f. Okubo-Marshak expansion):

$$U(\vec{r}, \vec{r'}) = V_c(r) + V_\sigma(r)(\vec{S}_1 \cdot \vec{S}_2) + S_{12}V_{T_1}(r) + O(\nabla^2) = V_C^{eff}(r) + O(\nabla^2)$$

- The convergence of the expansion can be checked.
- The NLO term is explicitly determined by utilizing two source functions (Iritani et. al, arXiv:1805.02365)

N. Ishii et al, PLB712, 437 (2010)

Time-dependent HAL method

• original (t-indep) HAL method=> applicable for each NBS w.f. $G_{BB}(\vec{r},t) = \langle 0 | B(\vec{y},t)B(\vec{x},t)\overline{\mathcal{J}}(t_0;J^P) | 0 \rangle$ $\mathcal{R}(\vec{r},t;J^P) = G_{BB}(\vec{r},t)/G_B(t)G_B(t) = \sum_{n} A_n \psi_n(\vec{r})e^{-(W_i-2m_B)t}$ $\int d\mathbf{r}' U(\mathbf{r},\mathbf{r}')\underline{\psi}_{W_0}(\mathbf{r}') = (\underline{E}_{W_0} - H_0)\underline{\psi}_{W_0}(\mathbf{r})$ $\int d\mathbf{r}' U(\mathbf{r},\mathbf{r}')\underline{\psi}_{W_1}(\mathbf{r}') = (\underline{E}_{W_1} - H_0)\underline{\psi}_{W_1}(\mathbf{r})$

- Many states contribute to the R-correlator
- As lattice size increases, the extraction of g.s. becomes difficult

$$E_n \sim 2\sqrt{m_B^2 + (2\pi n/L)^2} \sim E_0 \ (L \gg 1)$$

$$L = 3 \text{fm} \qquad L = 6 \text{fm} \qquad L = 8 \text{fm} \qquad L = \infty$$
Inelastic
Inelasti

The same problem appears for the direct method (Iritani et al. JHEP(2016), PRD(2017)) ₉

N. Ishii et al, PLB712, 437 (2010)

potential

Time-dependent HAL method

new (t-dep) HAL method=> directly applicable for R-correlator



Weaker condition as $L \rightarrow \infty$



 More strange quarks, more difficult experiment due to short life time
 Less strange quarks, more difficult numerical simulation due to increasing statistical noise

$\Omega\Omega$ system is the best S/N ratio calculation on lattice

Interactions in $\Omega\Omega$ (J=0) system

1) Nf = 2+1, L = 1.93fm, m_{π} =1015MeV, SU(3) limit 2) Nf=2+1, L = 3fm, m_{π} =700MeV, SU(3) breaking 3) Nf=2+1, L = 8.1fm, m_{π} =146MeV, almost physical mass

SG and K. Sasaki (HAL QCD)

1)Nf=2+1 full QCD with L = 1.93fm m_{π} = 1015MeV, SU(3) limit $\Omega\Omega$ in J =0



- Short range repulsive core and attractive pocket are found
- Phase shift shows the system is in the unitary limit

let's consider the lighter quark masses with SU(3) breaking

M. Yamada et. al.(HAL QCD), PTEP2015

2)Nf=2+1 full QCD with L = 3fm, m_{π} = 700MeV w. SU(3) breaking $\Omega\Omega$ in J=0m_Ω= 1970MeV



- Short range repulsive core and attractive pocket are found
- Potential is nearly independent on "t" within error
- Phase shift shows rapid changes depending on "t"
- The system may appear close to the unitary limit

c.f. Direct method by Buchoff et al., PRD(2012): L=4fm, m_{π} = 390MeV $a=0.16 \pm 0.22$ fm <= unitary limit

Numerical Setup at (almost) physical mass

2+1 flavor gauge configurations

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

- a= 0.0846 [fm], a⁻¹ = 2333 [MeV]
- 96³x96 lattice, L = 8.1[fm]
- 400 confs x 48 source positions x 4 rotations



Wall source is employed. only S-wave state is produced.

	[MeV]	phys.
π	146	8%
K	525	6%
Ν	964	3%
Ω	1712	2%



SG and K. Sasaki et.al.(HAL), PRL(2018)

"most strange dibaryon"

Nf=2+1 full QCD with L = 8.1fm, m_{π} = 146MeV

 $\Omega\Omega$ in J=0



- Short range repulsive core and attractive pocket are found
- Phase shift shows the presence of a bound state
- The state is very close to the unitary region (r/a<1)

SG and K. Sasaki et.al.(HAL), PRL(2018)

$\Omega\Omega$ in J = 0Binding energy and the Coulomb effect

"most strange dibaryon"



Conservative estimate at exact phys. pt.

 $m_{\pi=}146 \text{ MeV} \rightarrow 135 \text{ MeV}, m_{\Omega}= 1712 \text{MeV} \rightarrow 1672 \text{ MeV}$



conservative estimate:

only change the mass of kinetic term

 $(B_{\Omega\Omega}^{(\text{QCD})}, B_{\Omega\Omega}^{(\text{QCD+Coulomb})}) = (1.6(6) \text{MeV}, 0.7(5) \text{MeV})$ $\rightarrow (1.3(5) \text{MeV}, 0.5(5) \text{MeV})$ These changes are within errors

Kenji Morita (Wroclaw/RIKEN)

How HIC Can Tell Us Interaction?



 $C_{AB}(Q) = \frac{N_{AB}^{\text{pair}}(Q)}{N_A N_B(Q)} = \begin{cases} 1 & \text{No Correlation} \\ \text{others Interaction} \end{cases}$

Deviation from "1", tells us the behavior of the interaction **No Correlation**

TS Interaction Interference etc

Kenji Morita (Wroclaw/RIKEN)

ΩΩ Correlation@LHC





ΩΩ Correlation: Statistics?

Kenji Morita (Wroclaw/RIKEN)

of pair A(Q)



Summary

- We have investigated $\Omega\Omega$ interaction (J=0) from lattice QCD
- (almost) physical pion masses: $\Omega\Omega$ interaction in ${}^{1}S_{0}$
 - short range repulsive and attractive pocket
 - a very shallow bound state
 - [Most strange dibaryon, di-Omega]

Dibaryon

Deuteron



found in 1930s



will be found by J-PARC or FAIR?

Back Slides

Estimate NLO contribution for $\Omega\Omega$ system at almost physical pt.

 $U(\boldsymbol{r},\boldsymbol{r}') = V_0(r)\delta(\boldsymbol{r}-\boldsymbol{r}') + \sum_{n=1}V_{2n}(r)\nabla^{2n}\delta(\boldsymbol{r}-\boldsymbol{r}')$

• Determining the higher order potentials explicitly by utilizing multiple quark sources is the best way to estimate their contributions.

c.f. Iritani et. al (HAL QCD), arXiv:1805.02365

$$V(r) = R^{-1}(\boldsymbol{r}, t) \left(\frac{\nabla^2}{m_{\Omega}} - \frac{\partial}{\partial t} + \frac{1}{4m_{\Omega}} \frac{\partial^2}{\partial t^2} \right) R(\boldsymbol{r}, t)$$
$$= V_0(r) + \sum_{n=1}^{\infty} V_{2n}(r) R^{-1}(\boldsymbol{r}, t) \nabla^{2n} R(\boldsymbol{r}, t)$$

- Instead, we have estimated in two alternative ways:
 - 1. their contributions are estimated from its t-dependence
 - 2. pertubative estimate on the binding energy

 $|V_2 / m_{2\pi}| \sim |V_0|$ + several functional forms such as square-well form...

=> B.E. changes less than 20% in all cases

(within systematic errors from the t-dependence)