

Triton-like molecules

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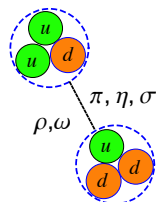
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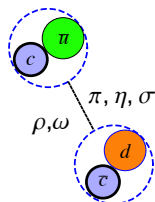
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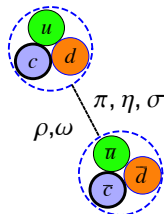
The deuteron-like molecules



Deuteron



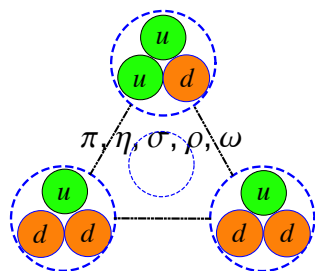
Di-meson molecule



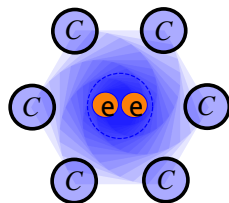
Di-baryon molecule

- $X(3872) \rightarrow D\bar{D}^*$, $D_{s1}(2460) \rightarrow D^*K$, $Z_b(10610) \rightarrow B\bar{B}^*$
- $Y(4260) \rightarrow D_1\bar{D}$, $Z_c(3900) \rightarrow D\bar{D}^*$, $Z_c(4020) \rightarrow D^*\bar{D}^*$
- $P_c(4380) \rightarrow \Sigma_c\bar{D}^*$, $P_c(4450) \rightarrow \Sigma_c^*\bar{D}^*$

The triton-like molecules



Triton



Benzene ring

- $NDK, \bar{K}DN, NDD\bar{D}, N\bar{K}K, DKK, DK\bar{K}, BBB^*, DDK, DD^*K, \Omega NN$ and $\Omega\Omega N$
- Faddeev equation, or FCA

Dimer or isobar formalism

Gaussian

expansion method (GEM)

Efimov effect

- $B_2 = \frac{\hbar^2}{ma^2}(1 + \mathcal{O}(\frac{r_0}{a}))$
- $B_3(1 + \mathcal{O}(\frac{r_0}{a})) = -\frac{\hbar^2}{ma^2} + [e^{-2\pi n} f(\xi)]^{\frac{1}{s_0}} \frac{\kappa_*^2}{m}$
- $s_0 = 1.00624\dots$, ξ is defined by $\tan \xi = -(mB_3)^{1/2} a/\hbar$, $f(\xi)$ is a universal function with $f(-\pi/2) = 1$,
 $K = \text{sgn}(E)\sqrt{m|E|}$.

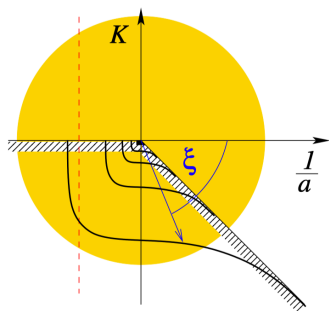


Figure: Efimov plot

Tri-baryon systems $\Lambda\Lambda\Lambda$, $\Xi\Xi\Xi$ and $\Sigma\Sigma\Sigma$

The OBE interaction indicates that there is only one virtual boson exchanged by any two constituents as shown in the following.

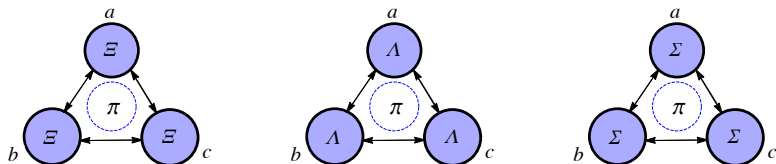
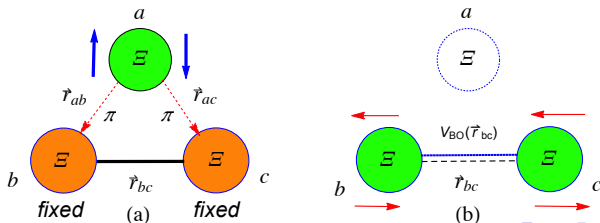


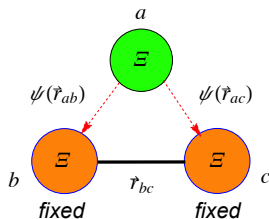
Figure: Dynamical illustration of the $\Lambda\Lambda\Lambda$, $\Xi\Xi\Xi$ and $\Sigma\Sigma\Sigma$ systems with a circle describing the delocalized π bond inside.

Monopole form factor $\mathcal{F}(q) = \frac{\Lambda^2 - m_\alpha^2}{\Lambda^2 - q^2}$, ($\alpha = \pi, \eta, \rho, \omega, \sigma, \phi$) with q the four-momentum of the pion and Λ the cutoff parameter.

Born-Oppenheimer potential

- Considering that the particle b and c are static with the separation r_{bc} , one can separate the degree of freedom of a from the three-body system.
- We assume the distance r_{bc} is a parameter. The mesons b and c are static, and have one-pion interactions with meson a , which can be viewed as two static sources.
- We explore the dynamics for the meson a in the limit $r_{bc} \rightarrow \infty$, and subtract the binding energy for the break-up state which is trivial for the three-body bound state.



Interpolating wave function of meson a 

The zero order of the final wave function for the meson a could be the superposition of these two components

$$\psi(\vec{r}_{ab}, \vec{r}_{ac}) = C[\psi(\vec{r}_{ab}) \pm \psi(\vec{r}_{ac})]|\Xi\Xi\Xi\rangle$$

Accordingly, one can obtain the energy eigenvalue of the meson a

$$E_a(\Lambda, \vec{r}_{bc}) = \langle \psi(\vec{r}_{ab}, \vec{r}_{ac}) | H_a | \psi(\vec{r}_{ab}, \vec{r}_{ac}) \rangle$$

BO potential and Its physical meaning (Intensity of "glue")

We define the BO potential as

$$V_{BO}(\Lambda, \vec{r}_{bc}) = E_a(\Lambda, \vec{r}_{bc}) - E_2(\Lambda).$$

The BO potential can describe the contribution for the one meson on the dynamics of the two remaining mesons. The meson a here works like a mass of "glue".

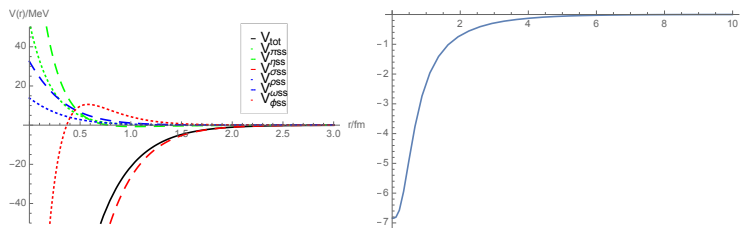
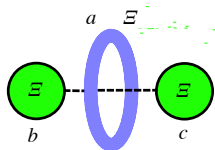


Figure: Here we chose the parameter $\Lambda = 900$ MeV. $E_{I=0}^{\Xi\Xi\Xi} = -3.09$ MeV.

$$\Psi_T = \alpha \Phi(\vec{r}_{bc}) \psi(\vec{r}_{ab}, \vec{r}_{ac}).$$

The configurations of the three-body systems

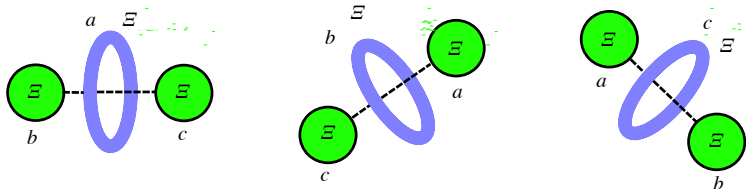


Figure: Every meson can be considered to be a lighter one and separated from the three-body system. Each of them can generate the "glue" for the remaining mesons.

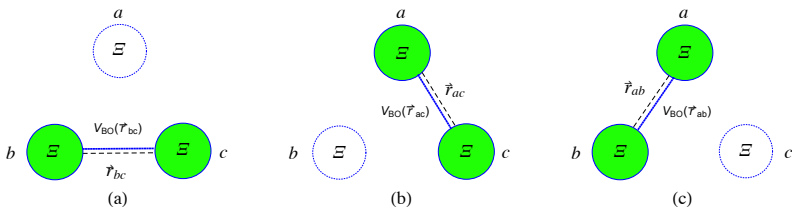


Figure: (a), (b) and (c) correspond to the wave functions $\psi_{\mathcal{A}} = \Phi(\vec{r}_{bc})\psi(\vec{r}_{ab}, \vec{r}_{ac})$, $\psi_{\mathcal{B}} = \Phi(\vec{r}_{ac})\psi(\vec{r}_{ab}, \vec{r}_{bc})$ and $\psi_{\mathcal{C}} = \Phi(\vec{r}_{ab})\psi(\vec{r}_{bc}, \vec{r}_{ac})$, respectively.

Interpolating wave functions

The basis constitute a configuration space $\{\psi_{\mathcal{A}}, \psi_{\mathcal{B}}, \psi_{\mathcal{C}}\}$.

$$\begin{aligned}\Psi_T &= \alpha\Phi(\vec{r}_{bc})\psi(\vec{r}_{ab}, \vec{r}_{ac}) + \beta\Phi(\vec{r}_{ac})\psi(\vec{r}_{ab}, \vec{r}_{bc}) + \gamma\Phi(\vec{r}_{ab})\psi(\vec{r}_{bc}, \vec{r}_{ac}) \\ &= \alpha\psi_{\mathcal{A}} + \beta\psi_{\mathcal{B}} + \gamma\psi_{\mathcal{C}} = \begin{pmatrix} \alpha \\ \beta \\ \gamma \end{pmatrix},\end{aligned}$$

Expand $\Phi(\vec{r}_{bc})$, $\Phi(\vec{r}_{ac})$ and $\Phi(\vec{r}_{ab})$ as a set of Laguerre polynomials

$$\chi_{nl}(r) = \sqrt{\frac{(2\lambda)^{2l+3}n!}{\Gamma(2l+3+n)}} r^l e^{-\lambda r} L_n^{2l+2}(2\lambda r), \quad n = 1, 2, 3\dots$$

$$\psi_{\mathcal{A}} = \sum_i \phi_i(\vec{r}_{bc})\psi(\vec{r}_{ab}, \vec{r}_{ac}), \psi_{\mathcal{B}} = \sum_i \phi_i(\vec{r}_{ac})\psi(\vec{r}_{ab}, \vec{r}_{bc}), \psi_{\mathcal{C}} = \sum_i \phi_i(\vec{r}_{ab})\psi(\vec{r}_{bc}, \vec{r}_{ac})$$

Orthonormalization

We orthonormalize the $\{\psi_{\mathcal{A}}, \psi_{\mathcal{B}}, \psi_{\mathcal{C}}\}$ into a new basis $\{\tilde{\psi}_{\mathcal{A}}, \tilde{\psi}_{\mathcal{B}}, \tilde{\psi}_{\mathcal{C}}\}$.

$$\tilde{\psi}_{\mathcal{A}}^i = \frac{1}{N_i} [(\psi_{\mathcal{A}}^i + \psi_{\mathcal{B}}^i + \psi_{\mathcal{C}}^i) - \sum_j x_{ij} \psi_{\mathcal{A}}^j],$$

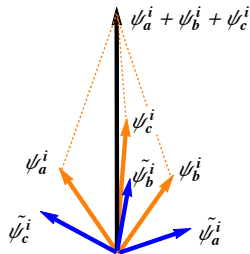
$$\tilde{\psi}_{\mathcal{B}}^i = \frac{1}{N_i} [(\psi_{\mathcal{A}}^i + \psi_{\mathcal{B}}^i + \psi_{\mathcal{C}}^i) - \sum_j x_{ij} \psi_{\mathcal{B}}^j],$$

$$\tilde{\psi}_{\mathcal{C}}^i = \frac{1}{N_i} [(\psi_{\mathcal{A}}^i + \psi_{\mathcal{B}}^i + \psi_{\mathcal{C}}^i) - \sum_j x_{ij} \psi_{\mathcal{C}}^j],$$

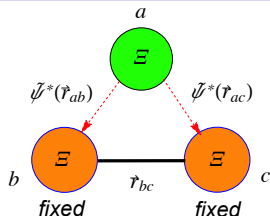
where the x_{ij} is a parameter matrix which will be determined later. The N_i are normalization coefficients.

Then the eigenvector for the three-body system $B_a^{(*)} B_b^{(*)} B_c^{(*)}$ can be written as a vector in the configuration space $\{\tilde{\psi}_{\mathcal{A}}, \tilde{\psi}_{\mathcal{B}}, \tilde{\psi}_{\mathcal{C}}\}$. Therefore, we have

$$\Psi_T = \sum_i \tilde{\alpha}_i \tilde{\psi}_{\mathcal{A}}^i + \sum_j \tilde{\beta}_j \tilde{\psi}_{\mathcal{B}}^j + \sum_k \tilde{\gamma}_k \tilde{\psi}_{\mathcal{C}}^k,$$



First order correction



The first order of the final wave function for the meson a could be the superposition of these two components

$$\tilde{\psi}^*(\vec{r}_{ab}, \vec{r}_{ac}) = C^*[\tilde{\psi}^*(\vec{r}_{ab}) \pm \tilde{\psi}^*(\vec{r}_{ac})]|\Xi\Xi\Xi\rangle$$

Accordingly, one can obtain the energy eigenvalue of the meson a

$$E_a^*(\Lambda, \vec{r}_{bc}) = \langle \tilde{\psi}(\vec{r}_{ab}, \vec{r}_{ac}) | H_a | \tilde{\psi}(\vec{r}_{ab}, \vec{r}_{ac}) \rangle$$

We define the BO potential as

$$V_{BO}^*(\Lambda, \vec{r}_{bc}) = E_a^*(\Lambda, \vec{r}_{bc}) - E_2(\Lambda).$$

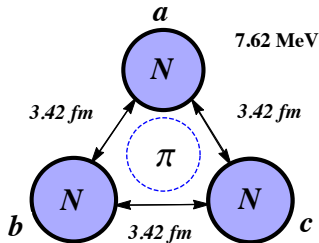
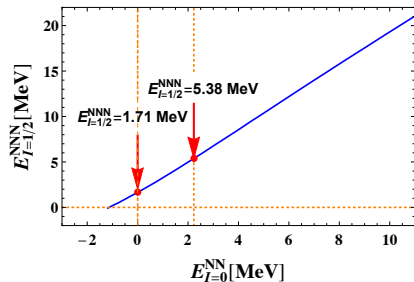
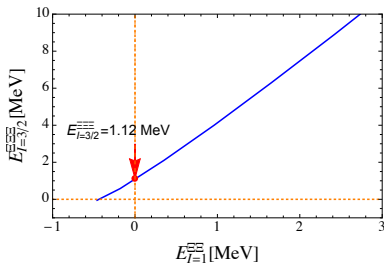
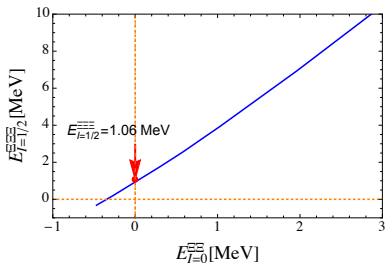
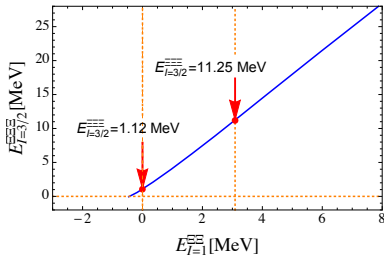
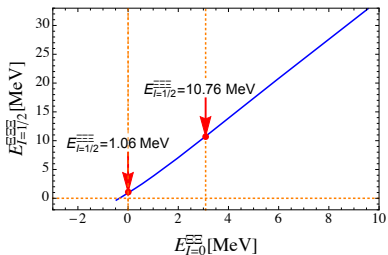
Application to the NNN system (Triton or Helium-3 nucleus)

Figure: Dependence of the reduced three-body binding energy on the binding energy of its two-body subsystem (the deuteron). The result is comparable with the empirical binding energies of the triton (8.48 MeV) and helium-3 (7.80 MeV) nuclei.

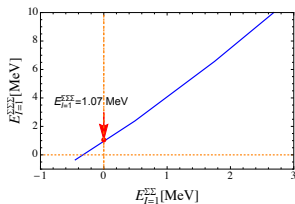
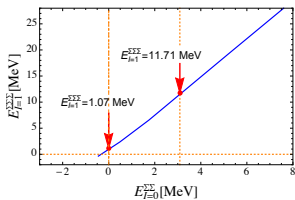
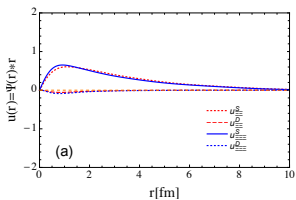
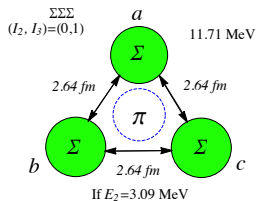
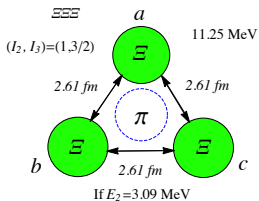
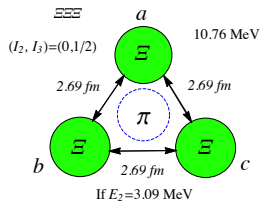
Numerical results for the NNN system (Triton or Helium-3 nucleus)

Λ (MeV)	E_2 (MeV)	E_3 (MeV)	E_T (MeV)	$V_{BO}(0)$ (MeV)	S wave(%)	D wave(%)	r_{rms} (fm)
830.00	-0.18	-1.93	-2.11	-4.54	94.01	5.99	4.21
850.00	-0.67	-2.71	-3.38	-5.36	93.36	6.64	4.00
870.00	-1.23	-3.65	-4.88	-6.32	92.68	7.32	3.78
890.00	-1.88	-4.77	-6.66	-7.42	91.99	8.01	3.54
899.60	-2.23	-5.38	-7.62	-8.00	91.66	8.34	3.42
900.00	-2.25	-5.41	-7.66	-8.03	91.64	8.36	3.42
920.00	-3.05	-6.85	-9.90	-9.35	90.97	9.03	3.18
940.00	-3.98	-8.51	-12.49	-10.83	90.35	9.65	2.95
960.00	-5.03	-10.42	-15.45	-12.46	89.76	10.24	2.74
980.00	-6.21	-12.57	-18.78	-14.23	89.23	10.77	2.54
1000.00	-7.55	-14.97	-22.51	-16.14	88.73	11.27	2.37
1020.00	-9.04	-17.61	-26.65	-18.19	88.27	11.73	2.23
1040.00	-10.69	-20.51	-31.20	-20.37	87.84	12.16	2.10

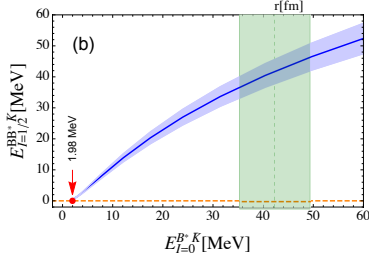
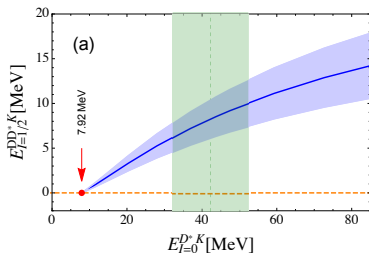
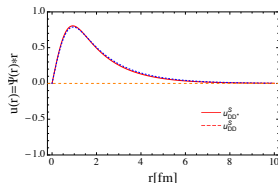
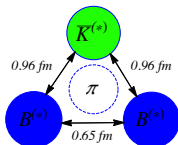
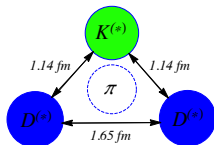
Table: Bound state solutions for the NNN system with isospin $I_3 = 1/2$. E_2 is the energy eigenvalue of its subsystem. E_3 is the reduced three-body energy eigenvalue relative to the break-up state of the NNN system. E_T is the total three-body energy eigenvalue relative to the NNN threshold.

Numerical results for the $\Lambda\Lambda\Lambda$, $\Xi\Xi\Xi$ and $\Sigma\Sigma\Sigma$ 

Numerical results for the $\Lambda\Lambda\Lambda$, $\Xi\Xi\Xi$ and $\Sigma\Sigma\Sigma$



Numerical results for the Doubly heavy tri-meson bound states



(MeV)	DD^*K	BB^*K	DD^*K	BB^*K
Binding energy	$8.29^{+4.32}_{-3.66}$	$41.76^{+8.84}_{-8.49}$	$8.29^{+6.55}_{-6.13}$	$41.76^{+9.02}_{-8.68}$
Mass	$4317.92^{+3.66}_{-4.32}$	$11013.65^{+8.49}_{-8.84}$	$4317.92^{+6.13}_{-6.55}$	$11013.65^{+8.68}_{-9.02}$

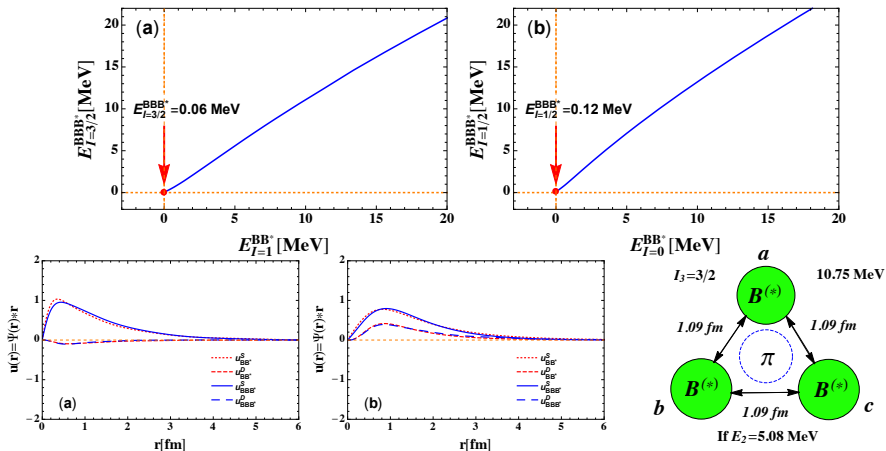
Numerical results for the tri-meson bound state BBB^* 

Figure: Here we chose the parameter $\Lambda = 1440$ MeV in (a) and $\Lambda = 1107.7$ MeV in (b) for a better comparison of all the cases, since they have the same two-body binding energy of 5.08 MeV.

Summary

- We predict that the triton-like molecular states for the $\Xi\Xi\Xi$ and $\Sigma\Sigma\Sigma$ systems are probably existent as long as the molecular states of their two-body subsystems exist.
- In our calculations, we use the Born-Oppenheimer potential method to construct our interpolating wave functions, which can be regarded as a version of the variational principle which always gives an upper limit of the energy of a system.
- Other configurations $N\Lambda\Lambda$, $N\Lambda\Xi$, $N\Xi\Xi$, $N\Lambda\Sigma$, $N\Sigma\Sigma$ will be explored in a future work.

Thank you very much!