The study of X(6900): fully-heavy charmonium system

2021/9/10 理论室学术报告"Weekly Theory Forum"

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Outline

- The introduction to XYZ exotic states
- The observation of X(6900) state
- Theoretical review for the fully-heavy charmonium system
- The introduction to Pomeron
- Diffractive scattering between $J/\psi \psi(2S)$ via the Pomeron exchange
- Summary

- ·传统夸克模型:很好的描述了介子(qq)、重子(qqq)能谱
- 自2003年Belle实验合作组发现奇特态X(3872)以来,各实验组观测到了越来越多的新强子结构,包括混杂态、胶球、多夸克态等

Guo et.al., RMP90(2018)015004

• 许多奇特态都靠近一些强子态的阈值附近





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介子: 3 ⊗ 3

· 多夸克态(Multi-quarks): 传统夸克模型的延申

四夸克态(Tetraquark): $3 \otimes 3 \otimes \overline{3} \otimes \overline{3}$



重子: 3 ⊗ 3 ⊗ 3 → 五夸克态(Pentaquarks): 3 ⊗ 3 ⊗ 3 ⊗ 3 ⊗ 3 §

• 强子分子态(Hadronic molecules):两体问题、三体问题...





- 理论研究方法
 - 1. 唯象层面:组分夸克势能模型、介子交换模型、运动学效应(三角奇异性等)
 - 2. 有效场论: 手征对称+重夸克对称理论

3. QCD求和规则

4. Lattice QCD

• 重要问题

1. QCD非微扰能区,低能有效常数的抽取

2. 短程相互作用动力学机制

3. 耦合道效应

• 介子交换模型(单道single-channel)

▶两个强子之间通过交换轻矢量介子产生的接触项

▶对接触项进行重求和得到T矩阵表达式

$$T = \frac{V}{1 - VG}$$

$$G(E) = \frac{1}{16\pi^2} \left\{ a(\mu) + \log \frac{m_1^2}{\mu^2} + \frac{m_2^2 - m_1^2 + s}{2s} \log \frac{m_2^2}{m_1^2} + \frac{k}{E} \log \frac{(2kE+s)^2 - m_1^2 + m_2^2}{(2kE-s)^2 - m_1^2 + m_2^2} \right\}$$

$$G(E) = \int \frac{l^2 dl}{4\pi^2} \frac{\omega_1 + \omega_2}{\omega_1 \omega_2} \frac{e^{-2l^2/\Lambda^2}}{E^2 - (\omega_1 + \omega_2)^2 + i\epsilon} \text{ with } \omega_i = \sqrt{m_i^2 + l^2}$$

• 在介子交换模型框架下, X(3872)被认为 $D\overline{D}$ *耦合形成的束缚态结构; Y(4260)可以成功被解释为 $D_1\overline{D}$ 的分子态图像



The observation of X(6900): fully-heavy charmonium system

- LHCb 在di-J/\/ 末态能谱中观测到了一些奇异结构
- 在J/ψ-ψ(2S)阈值附近看到了一个较窄的共振结构,被称 作X(6900),宽度大约为80MeV左右,其内部夸克组分为 cccc̄,属于全重粲味体系

threshold $_{J/\psi+\psi(2S)} = 6783 \text{MeV}$

• 这是实验上首次观测到可能的全重粲四夸克体系结构







LHCb Sci.Bull. 65 (2020) 23, 1983-1993

 $M(X(6900)) = 6905 \pm 11 \pm 7 \text{ MeV}$ $\Gamma(X(6900)) = 80 \pm 19 \pm 33 \text{ MeV}$

对于全重粲味的四夸克系统,长程的轻味介子交换在LO
 不贡献。而粲偶素的交换机制会受到传播子大质量压低
 以及cc产生的高能量标度压低





LHCb Sci.Bull. 65 (2020) 23, 1983-1993

从这个角度分析, X(6900)是一个好的Tetraquark候选者。
 有很多理论通过不同的势能模型在*J*/ψ-ψ(2S)阈值附近预
 言了许多cccccons波和P波结构,均可以衰变到di-*J*/ψ末态



<sup>V. R. Debastiani and F. S. Navarra Chin.Phys.C 43,013105(2018)
M. N. Anwar, J. Ferretti, F. K. Guo, E. Santopinto, and B. S. Zou, Eur. Phys. J. C 78, 647 (2018)
M. S. liu, F. X. Liu, X. H. Zhong and Q. Zhao, [arXiv:2006.11952 [hep-ph]] (2020)
Xin Jin, Yaoyao Xue, Hongxia Huang, Jialun Ping, Eur.Phys.J.C 80 (2020) 11, 1083(2020)
Jacob Sonnenschein, Dorin Weissman(Okinawa Inst. Sci. Tech.) Eur.Phys.J.C 81 (2021) 1, 25</sup>

- •然而,在LHCb的实验结果中,在6.9GeV附近值观测到X(6900)。 一个共振结构
- 探究全重粲味系统的内部动力学耦合机制尤为重要,短程相互作用在束缚全重系统中发挥着重要作用,cccc是研究非微扰QCD的良好场所。



LHCb Sci.Bull. 65 (2020) 23, 1983-1993



short-distance interaction

- 目前, X(6900)全重系统中软胶子传递强相互作用的理论工作 主要有
 - ✓ Hybrid interpretation of X(6900): Bing-Dong Wan, Cong-Feng Qiao(GUCAS) Phys.Lett.B 817 (2021) 136339
 - Two pioins interaction: Xiang-Kun Dong, Vadim Baru, Feng-Kun Guo, Christoph Hanhart, Alexey Nefediev, and Bing-Song Zou [arXiv: 2107.03946 [hep-ph]]

• X(6900): 混杂态解释 Bing-Dong Wan, Cong-Feng Qiao(GUCAS) Phys.Lett.B 817 (2021) 136339

 $[\overline{3}]_{cc} \otimes [8]_{G} \otimes [3]_{\bar{c}\bar{c}}$

▶该工作给出X(6900)的量子数为J^{PC} = 0⁺⁺

▶利用QCD求和规则计算得到m₀++ = (6.92 ± 0.14)GeV

▶预言在7.2GeV附近存在一个 $J^{PC} = 0^{-+}$ 混杂态



• 双-π交换相互作用机制 Xiang-Kun Dong, et. al [arXiv: 2107.03946 [hep-ph]]



→ $从\psi(2S) \rightarrow J/\psi \pi \pi$ 过程中抽取出 π 介子和矢量介子的相互作用强度,中性双 π 和 带电双π道是 $\psi(2S)$ 的主要衰变道,其衰变分支比分别为34.86%和18.24%

▶考虑双π末态相互作用,以及和KK道的耦合道效应,作者认为双π相互作用有可能可以提供di-J/ψ系统的吸引势

・双- π 交換相互作用机制 Xiang-Kun Dong, et. al [arXiv: 2107.03946 [hep-ph]] $\psi(2S) \rightarrow J/\psi\pi\pi$ 衰変过程中,主要由软胶子refusion到双 π 的道贡献

- •因此短程的软胶子相互作用在全重粲味系统中起着非常重要的作用
- 在我们的工作中,提出矢量介子之间交换Pomeron的相互作用机制



The Introduction To Pomeron



衍射过程是一个很很好的研究软强子-强子相互作用动力学的平台,在高能碰撞过程中,总散射截面表现出一个渐进常数的行为



P.A. Zyla et al. (Particle Data Group), Prog. Theor. Exp. Phys. 2020, 083C01 (2020)P. D. B. Collins, An Introduction to Regge Theory and High-Energy PhysicsA. Donnachie and P. V. Landshoff, Nucl. Phys. B231, 189 (1984)

- Pomeron trajectory: $\alpha_{\rm P}(t=0)=1$
- 理论认为Pomeron交换可以唯象为一对非微扰胶 子的相互作用行为

A. Donnachie and P. V. Landshoff, Nucl. Phys. B311, 509 (1988)

- 光学定理: $s\sigma_{tot} = \text{Im}T_{el}(s, t = 0)$
- 在Regge理论中, pp弹性散射截面 表示为:

$$T_{el}(s,t) \propto s^{\alpha_{\mathrm{P}}(t)} \Longrightarrow \sigma_{tot} \propto s^{\alpha_{\mathrm{P}}(0)-1}$$

The Introduction To Pomeron

• Donnachie和Landshoff在Regge理论框架下,将质子-质子弹性振幅描述为:

$$\frac{d\sigma}{dt} = \frac{1}{4\pi} (3\beta_0 F_1(t))^2 (\frac{s}{s_0})^{2\alpha_t(t)-2}$$

其中F₁(t)是质子的电磁形状因子(Pomeron相互作用在高能弹性散射过程中体现出类光子行为)

 β_0 是Pomeron和轻味夸克的耦合常数

Pomeron reggeon trajectory: $\alpha_{\rm p}(t) = 1 + 0.08 + 0.25t$

M. Diehl, Z. Phys. C 66, 181 193 (1995) A. Donnachie, J. Gravelis, and G. Shaw, Phys. Rev. D. 63, 114013(2001)

The Introduction To Pomeron

- Pomeron 交换模型很好地描述了高能强子碰撞程中小t区域的 散射行为和衍射行为
- pp弹性散射的振幅为

$$M_{pp} = (3\beta_0 F_1(t))^2 (\overline{u}\gamma_{\mu}u)(\overline{u}\gamma^{\mu}u)(\frac{s}{s_0})^{\alpha(t)-1}$$

G. A. Jaroszkiewicz and P. V. Landshoff, Phys. Rev. D10, 170 (1974)

• Pomeron 交換贡献到介子-质子弹性散射过程的微分截面大小 为 $\frac{d\sigma}{dt} = \frac{(2\beta_0 F_\pi(t))^2 (3\beta_0 F_1(t))^2}{4\pi} (\frac{s}{s_0})^{2\alpha(t)-2}$



C.W. Akerlof et. Phys.Rev.D 14 (1976) 2864

 $F_{\pi}(t)$ 是π的弹性散射形状因子, 拟合实验上πp弹性散射实验结果, 得到

$$F_{\pi}(t) = \frac{1}{1 - t / \lambda_0^2} \qquad (\lambda_0^2 = 0.5 \text{GeV}^2)$$

- 类比质子-质子、介子-质子的情形,我们可以写出介子-介子之间通过交换 pomeron 的相互作用行为
- 目前,还没有关于J/ψ-p的弹性散射实验数据,不过Pomeron相互作用仅依赖于味道,我们可以从J/ψ的光产生过程中提取出charm夸克和Pomeron相互作用的耦合系数βc







• t道和u道的散射振幅可以写为:

$$T_{t} = (2\beta_{c})^{2} (\frac{s}{s_{0}})^{\alpha(t)-1} \exp(\frac{t}{2\lambda_{J/\psi}^{2}}) \exp(\frac{t}{2\lambda_{\psi}^{2}}) \Gamma_{\alpha}^{\mu\rho} \Gamma^{\nu\alpha\sigma}$$
$$T_{u} = (2\beta_{c})^{2} (\frac{s}{s_{0}})^{\alpha(u)-1} \exp(\frac{\frac{u}{2} + \frac{m_{\psi}^{2}}{4} - \frac{m_{J/\psi}^{2}}{4}}{2\lambda_{J/\psi}^{2}}) \exp(\frac{\frac{u}{2} + \frac{m_{\psi}^{2}}{4} - \frac{m_{J/\psi}^{2}}{4}}{2\lambda_{\psi}^{2}}) \Gamma_{\alpha}^{\mu\sigma} \Gamma^{\nu\alpha\rho}$$

$$\Gamma^{\mu\alpha\rho} = (p_1 + p_3)^{\alpha} g^{\mu\rho} - 2 p_1^{\rho} g^{\alpha\mu}$$

Q. Zhao, J. P. Didelez, M. Guidal, and B. Saghai, Nucl. Phys. A660, 323 (1996) Alexander I. Titov, Yong-seok Oh, Shin Nan Yang, Tosiyuki Morii, Phys.Rev.C 58 (1998) 2429-2449

• 两个矢量介子的S波耦合量子数可能为:

 $J^{PC} = 0^{++}, 1^{++}, 2^{++}$

 我们发现Pomeron交换势是t依赖的,在近域处,我们只关注S波的贡献,因此我们 将角度依赖的部分积分掉得到一个contact顶点

$$V(s) = \frac{1}{2} \int V(s,t) d(\cos\theta)$$

▶这个顶点中包含了三种量子数的贡献,我们采用下面的投影算子将contact顶点进行投影

$$\mathcal{P}^{(0)} = \frac{1}{4} \epsilon_{\mu}(p_{1}) \epsilon^{\mu}(p_{2}) \epsilon_{\nu}(p_{3}) \epsilon^{\nu}(p_{4}),$$

$$\mathcal{P}^{(1)} = \frac{1}{2} (\epsilon_{\mu}(p_{1}) \epsilon_{\nu}(p_{2}) \epsilon^{\mu}(p_{3}) \epsilon^{\nu}(p_{4}) - \epsilon_{\mu}(p_{1}) \epsilon_{\nu}(p_{2}) \epsilon^{\nu}(p_{3}) \epsilon^{\mu})(p_{4}),$$

$$\mathcal{P}^{(2)} = \frac{1}{2} (\epsilon_{\mu}(p_{1}) \epsilon_{\nu}(p_{2}) \epsilon^{\mu}(p_{3}) \epsilon^{\nu}(p_{4}) + \epsilon_{\mu}(p_{1}) \epsilon_{\nu}(p_{2}) \epsilon^{\nu}(p_{3}) \epsilon^{\mu}(p_{4})) - \frac{1}{4} \epsilon_{\mu}(p_{1}) \epsilon^{\mu}(p_{2}) \epsilon_{\nu}(p_{3}) \epsilon^{\nu}(p_{4})$$

R.Molina, D. Nicmorus, and E. Oset Phys. Rev. D. 78, 114018 (2008)

 对于矢量介子圈的重求和,由于圈上粒子矢量求和导致张量结构较为复杂,无法 直接进行计算。但是,粲偶素介子的大质量可以提供一个很好的近似,圈积分的 三动量在近域范围内属于小量,在这种近似下,利用LS方程进行重求和

$$T = \frac{V}{1 - VG}$$

$$G(E) = \frac{1}{16\pi^2} \left\{ a(\mu) + \log \frac{m_1^2}{\mu^2} + \frac{m_2^2 - m_1^2 + s}{2s} \log \frac{m_2^2}{m_1^2} + \frac{k}{E} \log \frac{(2kE+s)^2 - m_1^2 + m_2^2}{(2kE-s)^2 - m_1^2 + m_2^2} \right\}$$

$$G(E) = \int \frac{l^2 dl}{4\pi^2} \frac{\omega_1 + \omega_2}{\omega_1 \omega_2} \frac{e^{-2l^2/\Lambda^2}}{E^2 - (\omega_1 + \omega_2)^2 + i\epsilon} \text{ with } \omega_i = \sqrt{m_i^2 + l^2}$$



Chang Gong, Meng-Chuan Du, Bin Zhou, Qiang Zhao, Xian-Hui Zhong[arXiv:2011.11374[hep-ph]]

$$V_{11}^{\Lambda} + V_{11}^{\Lambda} G_{1}^{\Lambda} + V_{12}^{\Lambda} G_{2}^{\Lambda} + V_{21}^{\Lambda} + \dots$$

$$G = \begin{pmatrix} G_{11} & 0 & 0 \\ 0 & G_{22} & 0 \\ 0 & 0 & G_{33} \end{pmatrix}$$

$$T(E) = V + VG(E)V + VG(E)VG(E)V + \dots = \frac{1}{V^{-1} - G(E)}$$

$$\mathbf{V} = \begin{pmatrix} V_{11} & V_{12} & V_{13} \\ V_{21} & V_{22} & V_{23} \\ V_{31} & V_{32} & V_{33} \end{pmatrix}$$

$$G(E) = \frac{1}{16\pi^2} \left\{ a(\mu) + \log \frac{m_1^2}{\mu^2} + \frac{m_2^2 - m_1^2 + s}{2s} \log \frac{m_2^2}{m_1^2} + \frac{k}{E} \log \frac{(2kE+s)^2 - m_1^2 + m_2^2}{(2kE-s)^2 - m_1^2 + m_2^2} \right\}$$

$$G(E) = \int \frac{l^2 dl}{4\pi^2} \frac{\omega_1 + \omega_2}{\omega_1 \omega_2} \frac{e^{-2l^2/\Lambda^2}}{E^2 - (\omega_1 + \omega_2)^2 + i\epsilon} \text{ with } \omega_i = \sqrt{m_i^2 + l^2}$$

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• We consider three coupled channel J/ψ - J/ψ , J/ψ - $\psi(2S)$ and $\psi(2S)$ - $\psi(2S)$

M. L. Du, V. Baru, F. K. Guo, C. Hanhart, U. G. Meißner, A. Nefediev and I. Strakovsky, Phys.Rev.Lett. 126, (2021) 132001

$$M_{1} = P(\sqrt{s}) \left(1 + \sum r_{i} G_{i}(s) T_{i1} \right)$$
$$\Gamma(s) = \frac{\left| \overline{P_{J/\psi}} \right|}{8 \pi s} |M_{1}|^{2}$$

• The phase difference between different channel.

 $r_1: r_2: r_3 = 1: 2e^{-i\pi/2}: 1$

• The X(6900) state favors $J^{PC} = 0^{++}, 2^{++}$



Chang Gong, Meng-Chuan Du, Bin Zhou, Qiang Zhao, Xian-Hui Zhong[arXiv:2011.11374[hep-ph]]

Diffractive scattering between η_c - η_c via the Pomeron exchange



$$\begin{split} \Gamma^{\alpha} &= \ (\frac{4m_c^2 + p_3^2}{p_3^2}) p_1^{\alpha} + (\frac{4m_c^2 - p_3^2}{p_3^2}) p_3^{\alpha} \\ &\approx \ 2p_1^{\alpha}, \end{split}$$

• The S-wave couplings between two pseudo-scalar charmonia:

 $J^{PC} = 0^{++}$



- In the case that the X(6900) being an tetra-quark state with quantum number 0^{++} , the X(6900) can also be observed in the η_c η_c spectrum.
- However, the Pomeron exchange mechanism for the fully-heavy quark system gives different pole positions for different channel.

Summary

- For fully heavy charmonium system, the pomeron exchange potential offers a short-distance interaction mechanism
- The S-wave $J/\psi \cdot \psi(2S)$ interaction via the Pomeron exchange can reproduce the X(6900) state.
- The di-pseudoscalar charmonia spectrum is a good tool to reveal the underlying dynamics in the fully heavy charmonium system.

Thanks For Your Attention.