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Heavy Baryons: the Υ Rooftops

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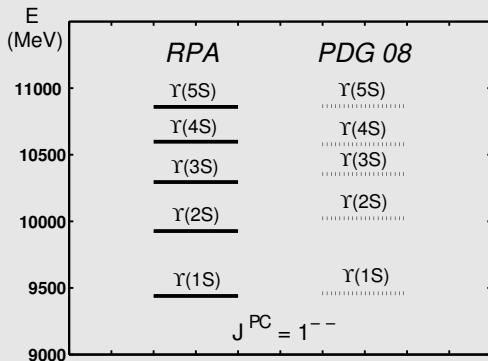
STRONG INTERACTIONS BEYOND THE STANDARD MODEL

Physikzentrum Bad Honnef, Germany, February 13-15, 2012

Introduction

Bottomonium

Corner of QCD where we understand how the hadronic bound-states are composed of quarks.

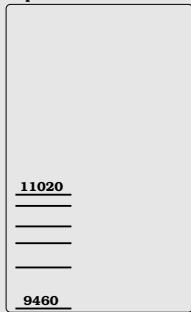




Υ -spectrum

Pure Exotics

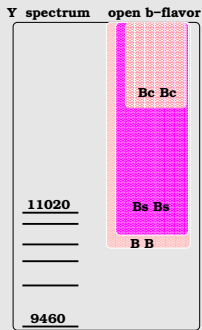
Υ spectrum



Observation Resonances in $B_c \overline{B}_c$ should be (hidden) exotic.

Υ -spectrum

Pure Exotics

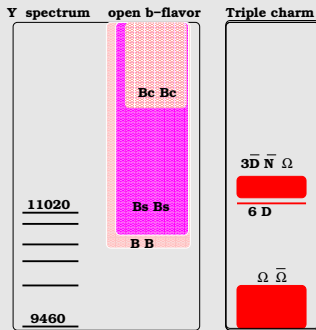


Observation Resonances in $B_c \overline{B_c}$ should be (hidden) exotic.

- Direct test of weak vs strong pNRQCD regime (Coulombic interaction: very weak resonances, bound states)
- Roof of future B -factory's energy should reach $B_c \overline{B_c}$ threshold to be exhaustive (B, B_s, B_c).

Υ -spectrum

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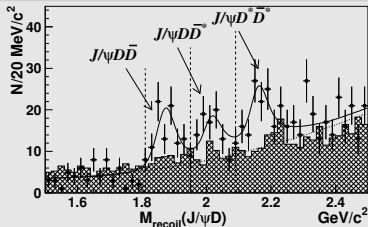
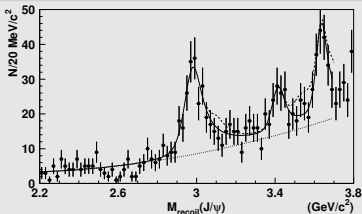
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Triple Charm

Look for Ω_{ccc} in $e^+ e^- \rightarrow \Omega_{ccc} \bar{p} \bar{D} \bar{D} \bar{D}$ by recoil method.

Double charmonium/charm production

Since 1987 (WA75 collaboration)



Using recoil method and missing mass.
Tagging D -mesons in final state.
Lots of experience.

BELLE collaboration hep-ex/0507019

Triple charm?

Thresholds

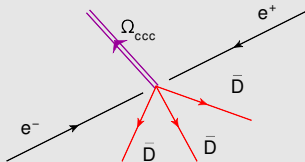
Closed charm: Triple c not reported ($3 \times M_{J/\psi} = 9300 \text{ MeV}$)

Open charm: Produce $ccc + \bar{c}\bar{c}\bar{c}$

- $DDD + \overline{DDD} \rightarrow 6 \times 1867 = 11202$
- Triply charmed baryon $\Omega_{ccc} + \overline{DDD} + \bar{N}$

No measurement of triply charmed baryons yet: theory.

Recoil





Static potential in pNRQCD

Expansion in α , $1/m$

$$V^{(0)} = V_{LO}^{(0)} + V_{NLO}^{(0)} + V_{NNLO}^{(0)} \dots$$

Leading Order (LO)

The leading order potential is just the colour Coulomb potential

$$V_{LO}^{(0)} = -\frac{4}{3} \frac{\alpha_s (r^{-2})}{r}$$



Static potential in pNRQCD

NLO and NNLO

$$V_{NLO}^{(0)} = V_{LO}^{(0)} (a_1 + 2\gamma_E \beta_0) \frac{\alpha_s (r^{-2})}{4\pi}$$

$$V_{NNLO}^{(0)} = V_{LO}^{(0)} \left(\gamma_E (4a_1 \beta_0 + 2\beta_1) + \left(\frac{\pi^2}{3} + 4\gamma_E^2 \right) \beta_0^2 + a_2 \right) \frac{\alpha_s^2 (r^{-2})}{(4\pi)^2}$$

 $V_{1/m}$ (convention dependent)

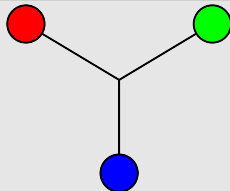
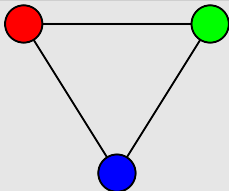
$$V_{m^{-1}} = \frac{-\alpha_s^2(\mu)}{m_r r^2} \times \left(\frac{7}{9} \right) - \frac{\alpha_s^3}{3\pi m_r r^2} \left\{ -b_2 + \log(e^{2\gamma_E} \mu^2 r^2) \left(\frac{7\beta_0}{6} + \frac{68}{3} \right) \right\}$$

(A. Vairo 2001; N. Brambilla et al. 2000)



Ω_{ccc} : Potential in pNRQCD

Δ or Y-Shaped potential





Ω_{CCC} : Potential in pNRQCD

LO Potential: Δ -shaped

$$V_{LO}^{(0)} = \frac{-2\alpha_s}{3} \sum_{i=1}^3 \frac{1}{|\mathbf{r}_i|} \quad (\text{with } \mathbf{r}_1 = \mathbf{r}_2 - \mathbf{r}_3, \text{ etc})$$

i.e. sum of two-body Coulomb interactions



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i.e. sum of two-body Coulomb interactions

NLO Potential: Still Δ -shaped

$$V_{LO}^{(0)} + V_{NLO}^{(0)} = \frac{-2}{3} \sum_i \alpha_s (|\mathbf{r}_i|^{-2}) \frac{1}{|\mathbf{r}_i|} \left[1 + \frac{\alpha_s (|\mathbf{r}_i|^{-2})}{4\pi} (2\beta_0 \gamma_E + a_1) \right]$$



Ω_{CCC} : Potential in pNRQCD

NNLO Potential: Δ -shaped Component

$$V_{NNLO-2}^{(0)} = \frac{-2}{3} \sum_i \frac{\alpha_s(\mathbf{r}_i^{-2})}{|\mathbf{r}_i|} \frac{\alpha_s(\mathbf{r}_i^{-2})^2}{(4\pi)^2} \times$$
$$\left(a_2 - 36\pi^2 + 3\pi^4 + \left(\frac{\pi^2}{3} + 4\gamma_E^2 \right) \beta_0^2 + \gamma_E(4a_1\beta_0 + 2\beta_1) \right)$$



Ω_{CCc} : Potential in pNRQCD

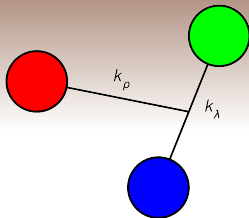
NNLO Potential: Δ -shaped Component

$$V_{NNLO-2}^{(0)} = \frac{-2}{3} \sum_i \frac{\alpha_s(\mathbf{r}_i^{-2})}{|\mathbf{r}_i|} \frac{\alpha_s(\mathbf{r}_i^{-2})^2}{(4\pi)^2} \times \\ \left(a_2 - 36\pi^2 + 3\pi^4 + \left(\frac{\pi^2}{3} + 4\gamma_E^2 \right) \beta_0^2 + \gamma_E(4a_1\beta_0 + 2\beta_1) \right)$$

NNLO Potential: Finally a Y -shaped potential

$$V_{NNLO-3}^{(0)} = 2(V_{aux}(\mathbf{r}_2, \mathbf{r}_3) + V_{aux}(\mathbf{r}_1, -\mathbf{r}_3) + V_{aux}(-\mathbf{r}_2, -\mathbf{r}_1)) \\ V_{aux}(\mathbf{r}_2, \mathbf{r}_3) = FT. \hat{V}_{aux}(\mathbf{q}_2, \mathbf{q}_3) = \frac{(-i/2)(4\pi)^3 \alpha_s^3}{8|\mathbf{q}_2|^2 |\mathbf{q}_3|^2} \times \\ \left[\frac{|\mathbf{q}_2 + \mathbf{q}_3|}{|\mathbf{q}_2| |\mathbf{q}_3|} + \frac{\mathbf{q}_2 \cdot \mathbf{q}_3 + |\mathbf{q}_2| |\mathbf{q}_3|}{|\mathbf{q}_2| |\mathbf{q}_3| |\mathbf{q}_2 + \mathbf{q}_3|} - \frac{1}{|\mathbf{q}_2|} - \frac{1}{|\mathbf{q}_3|} \right]$$

Variational



Jacobi coordinates

$$F(\mathbf{k}_\rho, \mathbf{k}_\lambda) = \psi\left(\frac{\mathbf{k}_\rho}{\alpha_\rho}\right) \psi\left(\frac{\mathbf{k}_\lambda}{\alpha_\lambda}\right)$$

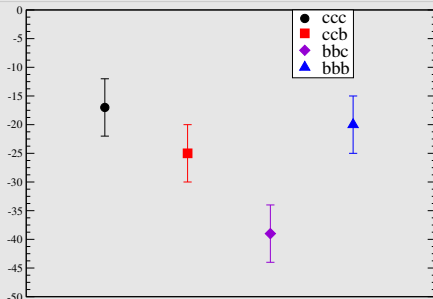
Minimize

$$\langle N|H|N\rangle(\alpha_\rho, \alpha_\lambda)$$

Parameters $\alpha_\rho, \alpha_\lambda$

Binding of three-body force

$M_Y - M_\Delta$

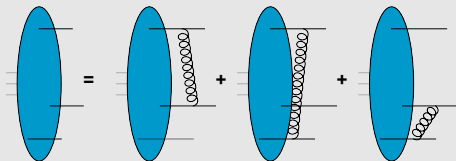


- Including three-body force \rightarrow increased binding.
- Equal flavour objects more tightly bound than mixed flavour
- Little quark mass dependence.

Faddeev equations in QCD

Covariant Faddeev equations

- Good news for Faddeev Equations (three-body force neglected).
- Nucleon studies suggest quark-diquark picture reasonable.



Also apply to triply heavy quark systems.

H. Sanchis-Alepuz *et al*, Phys. Rev. D **84** (2011) 096003.

H. Sanchis-Alepuz *et al*, PoS QCD-TNT-II (2011) 041



Results for Ω_{CCC}

Predicted Masses

NNLO pNRQCD

- 4800(250) MeV

Covariant Faddeev equations

- 4400 MeV (Gluon model MRT)
- 4900 MeV (Gluon model AFW)

Parameter Fitting

In both approaches, observables from meson sector are fit.
(see papers for details)



Summary and Outlook

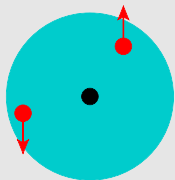
Summary

- pNRQCD static potential at NLO, NNLO. Fit experimental data.
- Covariant Faddeev equations in RL. Best fit to experiment.
- Constrain triply heavy baryon systems.
- Three-body force is small at NNLO
 - Good news for Faddeev equation.

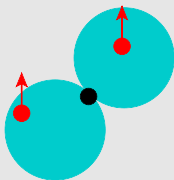
Super-B should pay attention to triple charm

Error estimate: Variational Method

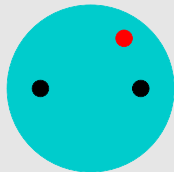
Compare to (simple) atomic/molecular systems



Para-Helium



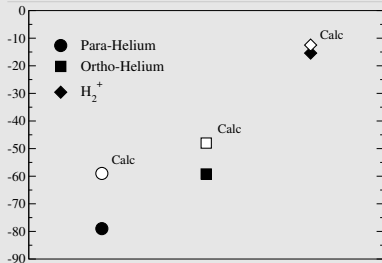
Ortho-Helium



Dihydrogen cation

Error estimate: Variational Method

Binding energies



- Para-Helium:
-79.0 eV (Calc. -59 eV)
- Ortho-Helium:
-59.2 eV (Calc. -48 eV)
- H_2^+ :
-15.4 eV (Calc. -12.5 eV)

Binding energy underestimated by 25%.
For 3-body, sign known.