

Tetraquarks in the Steiner tree model of confinement

available at <http://lpsc.in2p3.fr/theorie/Richard/SemConf/Talks.html>

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Introduction

Speculations on multiquarks

- low mass of $(q\bar{q})$, hence $(qq\bar{q}\bar{q})$ in S-wave perhaps lighter than $(q\bar{q})$ in P-wave. Applied to the problem of scalar mesons.
- Peculiar features of chiral dynamics. Speculations on the late pentaquarks made of light or strange quarks or antiquarks,
- Coherences in the hyperfine interaction → see next section,
- Properties of the mass dependence in a **flavour-independent** potential, → below
- Favourable 4-body interaction in QCD → below.

Symmetry breaking and tetraquarks-1

Consider

$$H = H_0(\text{even}) + \lambda H_1(\text{odd}).$$

Then for the ground state, with $\psi_0(H_0)$ as trial w.f, $\langle \psi_0 | H_1 | \psi_0 \rangle = 0$

$$E(H) \leq E(H_0),$$

i.e., H benefits of symmetry breaking.

For instance $E(p^2 + x^2 + \lambda x) = 1 - \lambda^2/4$.

This is very general.

Starting, e.g., from a symmetrical **four-body system** $(\mu, \mu, \bar{\mu}, \bar{\mu})$ breaking **particle identity** or **charge conjugation** lowers the ground state, but has different consequences on **stability**

Breaking particle identity

$H(M, m, M, m)$, where V does not change if M and/or m is modified, can be rewritten as

$$H = \underbrace{\left(\frac{1}{4M} + \frac{1}{4m} \right) [\mathbf{p}_1^2 + \cdots + \mathbf{p}_4^2]}_{H_0} + V + \underbrace{\left(\frac{1}{4M} - \frac{1}{4m} \right) [\mathbf{p}_1^2 - \mathbf{p}_2^2 + \mathbf{p}_3^2 - \mathbf{p}_4^2]}_{H_1}$$

Thus $E(H) \leq E(H_0)$. But in general, the threshold *also* benefits from this symmetry breaking, and actually benefits **more**, so that four-body binding **deteriorates**.

For instance, in atomic physics (e^+, e^+, e^-, e^-) and any equal-mass ($\mu^+, \mu^+, \mu^-, \mu^-$) weakly bound below the atom–atom threshold, but (M^+, m^+, M^-, m^-) unstable for $M/m \gtrsim 2.2$, see Dario.

Then: breaking the symmetry of **identical particles** does not help

Breaking charge conjugation

$H(M, M, m, m)$ written as

$$H = \underbrace{\left(\frac{1}{4M} + \frac{1}{4m} \right) [\mathbf{p}_1^2 + \dots + \mathbf{p}_4^2]}_{H_0} + V + \underbrace{\left(\frac{1}{4M} - \frac{1}{4m} \right) [\mathbf{p}_1^2 + \mathbf{p}_2^2 - \mathbf{p}_3^2 - \mathbf{p}_4^2]}_{H_1}$$

still benefits to the four-body system, $E(H) \leq E(H_0)$, but H and H_0 have the **same** threshold $(M^+, m^-) + (M^+, m^-)$. Hence **binding improves**. Indeed, H_2 more bound than P_{S_2} and has even a rich spectrum of excitations.

Quark model analogs

For a central, **flavour-independent**, confining interaction V ,

- Equal mass case (q, q, \bar{q}, \bar{q}) hardly bound
- Hidden-flavour case (Q, q, \bar{Q}, \bar{q}) even farther from binding,
- $(QQ\bar{q}\bar{q})$ with flavour = 2 bound if M/m large enough
See Ader et al. (then at CERN), Heller et al. (Los Alamos), Zouzou et al. (Grenoble), D. Brink et al. (Oxford), Rosina et al. (Slovenia), Lipkin, Vijande et al., etc.

$(QQ\bar{q}\bar{q})$ expected at least in the limit of **large** or **very large** M/m .
As compared to the “colour-chemistry” (late 70’s and early 80’s)

- **no exotic colour** configuration
- for large M/m , almost pure $\mathbf{3} \rightarrow \bar{\mathbf{3}}$ for (QQ) as in every baryon,
- and then $\bar{\mathbf{3}} \times \bar{\mathbf{3}} \times \bar{\mathbf{3}} \rightarrow \mathbf{1}$ for $[(QQ) - \bar{q}\bar{q}]$ as in every antibaryon:
well probed colour structure.

The additive model of tetraquark confinement

Questions:

- What is V for tetraquarks?
- Even earlier: what is the link from mesons to baryons?

The additive model By analogy with QED,

$$V(1, 2, \dots) = -\frac{3}{16} \sum_{i < j} \tilde{\lambda}_i^{(c)} \cdot \tilde{\lambda}_j^{(c)} v(r_{ij}) ,$$


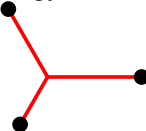
- $v(r)$ is the quarkonium potential fitted to mesons,
- $\lambda^{(c)}$ is the non-abelian colour operator

Consequences

- A reasonable simultaneous phenomenology of baryon and meson spectra
- Multiquarks **unbound**, except $(QQ\bar{q}\bar{q})$ with large M/m .
- Hence multiquark binding was based on other mechanisms: chromomagnetism (Jaffe, Lipjkin, Gignoux et al.), chiral dynamics (cf. the late pentaquark), etc.

The Steiner-tree model of baryons

Y-shape potential:

- Proposed by Artru, Dosch, Merkuriev, etc., proposed a better ansatz, often verified and rediscovered (strong coupling, adiabatic bag model (Kuti et al.), flux tube, lattice QCD, etc.)
- The linear $q - \bar{q}$ potential of mesons interpreted as minimising the gluon energy in the flux tube limit
 
- The $q - q - q$ potential of baryons is with the junction **optimised**, i.e., fulfilling the conditions of the well-known Fermat-Torricelli problem.
 

This potential was used for baryons (Taxil et al., Semay et al., Kogut et al.), but it does not make much difference as compared with the additive ansatz.

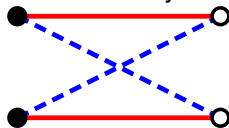
The Steiner tree model of tetraquarks

A tempting generalisation to **tetraquarks** is the combination

$$V_4 = \min(V_f, V_S) \text{ of}$$

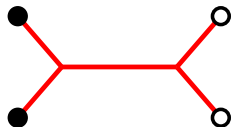
- **flip-flop** V_f (already used in its quadratic version by Lenz et al.)

$$V_f = \lambda \min(r_{13} + r_{24}, r_{23} + r_{14})$$



- and **Steiner-tree** V_S

$$V_S = \lambda \min_{k,l} (r_{1k} + r_{2k} + r_{kl} + r_{l3} + r_{l4}) .$$



- J. Carlson and V.R. Pandharipande concluded that this potential does not bind, **but**

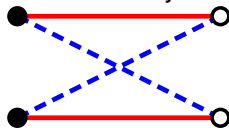
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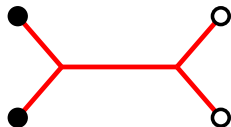
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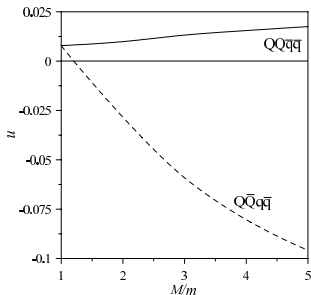
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- J. Carlson and V.R. Pandharipande concluded that this potential does not bind, **but**
- they used **too simple** trial wave functions for the 4-body problem, and did not consider **unequal masses**.

Tetraquarks in the minimal-path model-1

Vijande, Valcarce and R. revisited the calculation of Carlson et al. with a basis of correlated Gaussians (matrix elements painfully calculated numerically), and obtained **stability** for $(QQ\bar{q}\bar{q})$ even for $M/m = 1$, but better stability for $M/m \gg 1$.



$$u = (E_{th} - E_4)/E_{th}$$

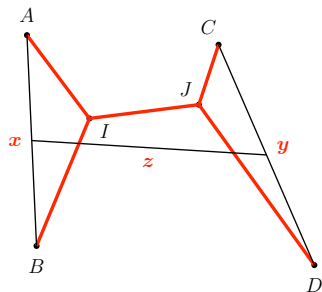
Tetraquarks in the minimal-path model-2

More recently, Cafer Ay, Hyam Rubinstein (Melbourne) and R.: **rigorous** proof of stability within the minimal-path model if $M \gg m$.

Obviously,

$$V_4 \leq V_S \leq |\mathbf{x}| + |\mathbf{y}| + |\mathbf{z}|$$

where $\mathbf{x} = \overrightarrow{AB}$, $\mathbf{y} = \overrightarrow{CD}$,
and \mathbf{z} links the middles.



Then

$$H \leq \left[\frac{p_x^2}{M} + |\mathbf{x}| \right] + \left[\frac{p_y^2}{m} + |\mathbf{y}| \right] + \left[\frac{p_z^2}{2\mu} + |\mathbf{z}| \right]$$

exactly solvable, but not does **not** demonstrate binding of $(QQ\bar{q}\bar{q})$

Better bound

- A better bound **demonstrates** stability for large M/m :

$$H \leq \left[\frac{p_x^2}{M} + \frac{\sqrt{3}}{2} |\mathbf{x}| \right] + \left[\frac{p_y^2}{m} + \frac{\sqrt{3}}{2} |\mathbf{y}| \right] + \left[\frac{p_z^2}{2\mu} + |\mathbf{z}| \right]$$

- $p^2 + |\mathbf{x}| \implies e_0 = 2.3381\dots$ (Airy function)
- by **scaling** $p^2/m + \lambda|\mathbf{x}| \implies e_0 \lambda^{2/3} m^{-1/3}$.
- Threshold** $2(Q\bar{q})Q\bar{q}$ at $E_{\text{th}} = 2e_0\mu^{-1/3}$, $\mu = Mm/(M+m)$.
- The **tetraquark** energy has an upper bound

$$E_4 \leq E_4^{\text{up}} = e_0 \left\{ \left(\frac{3}{4} \right)^{1/3} \left[M^{-1/3} + m^{-1/3} \right] + (2\mu)^{-1/3} \right\}$$

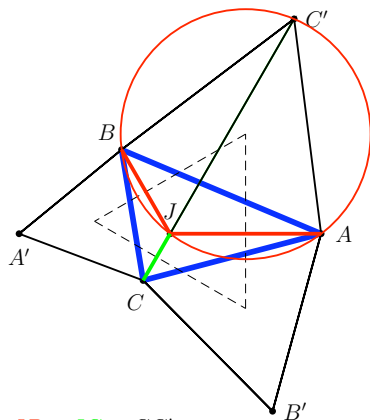
- Straightforward to check that $E_4^{\text{up}} < E_{\text{th}}$ for $M/m < 6403$
- Thus $E_4 < E_{\text{th}}$ at large M/m demonstrated rigorously
- Actually $\forall M/m$ from solving numerically the 4-boby pb.

Proof-1

A flavour of the proof. In the 3-body case, Steiner tree linked to [Napoleon's theorem](#).

$JA + JB + JC = CC'$ where C' makes an external equilateral triangle associated to the side AB .

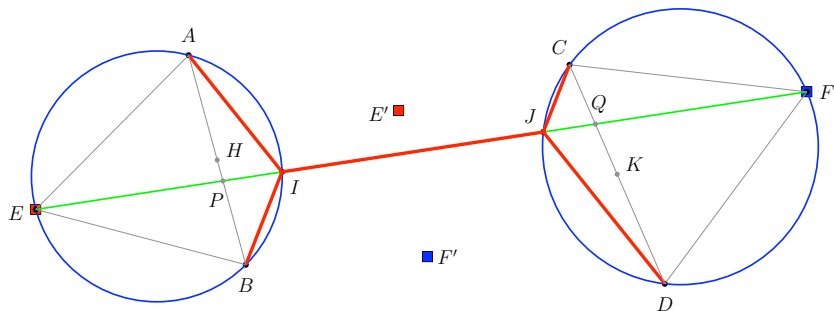
Well-known property of the Fermat-Torricelli problem. (C' belongs to the torroïdal domain associated to AB)



$$JA + JB + JC = CC'$$

Proof-2

The analogue for the planar tetraquark is



$$V_S = JA + JB + JK + KC + KD = EF$$

The **minimal** network linking (A, B, C, D) is the **maximal** distance between $\{E, E'\}$ and $\{F, F'\}$, which are the torroïdal domains associated to (A, B) and (C, D) (= points completing an equilateral tr.)

Conclusions

- New models of **confinement** beyond naive additive models, in better agreement with QCD in the strong coupling limit
- New inequality on the combined flip-flop and Steiner-tree paths,
- Drastic revision of the four-body spectrum within this model
- Analogous to the Wheeler (1945) – Ore (1946) – Hyllerras & Ore (1947) views on the Ps_2 molecule.

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Binding Energy of Polyelectrons

ADNE ORE
Sloane Physics Laboratory, Yale University, New Haven, Connecticut
June 10, 1946

THE question as to the existence of groups of electrons and positrons having temporary stability has recently been raised by J. A. Wheeler,¹ who shows that clusters of

Although the evidence here presented against the stability of the polyelectron composed of two electrons and two positrons is not conclusive in a strict mathematical sense, it counsels against the assumption that clusters of this (or even of higher) complexity can be formed.

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Binding Energy of the Positronium Molecule

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AND

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(Received December 26, 1946)

A system of two electrons and two positrons is shown to possess dynamic stability. The variational calculation performed leads to a binding energy of at least 0.11 eV for this cluster. The approximate wave function which yields this value depends on the four electron-positron distances only. Neglect of the two distances between particles of the same kind permits an essential mathematical simplification which might be of interest in other problems.