

Bound states in perturbation theory

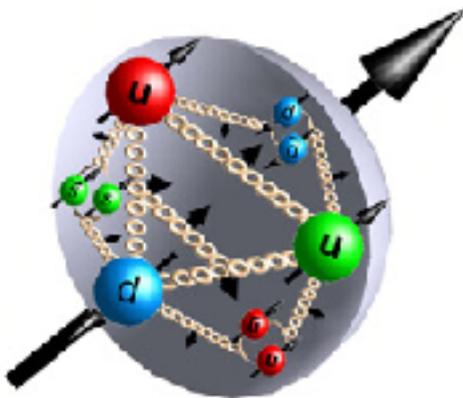
Bad Honnef, February 12-15, 2017

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Positronium has both perturbative $\mathcal{O}(\alpha)$ and non-perturbative $\mathcal{O}(\alpha^\infty)$ aspects.
Atoms are accurately described by QED.



Hadrons have confinement and CSB.
They also have perturbative features.
Is there a QCD approach analogous to QED?

arXiv:1612.09463
arXiv:1605.01532

QED works for atoms

The PQED expansion (usually) starts from the $\mathcal{O}(\alpha^\infty)$ **Schrödinger atom**.

Example: Hyperfine splitting in Positronium at $\mathcal{O}(\alpha^7 \log \alpha)$ using NRQED

Orthopositronium: $J^{PC} = 1^{--}$ Parapositronium: $J^{PC} = 0^{-+}$

$$\Delta E = E(\text{ortho}) - E(\text{para}) \qquad \Delta\nu = \Delta E / 2\pi\hbar$$

$$\begin{aligned} \Delta\nu_{QED} = m_e\alpha^4 & \left\{ \frac{7}{12} - \frac{\alpha}{\pi} \left(\frac{8}{9} + \frac{\ln 2}{2} \right) \right. \\ & + \frac{\alpha^2}{\pi^2} \left[-\frac{5}{24}\pi^2 \ln \alpha + \frac{1367}{648} - \frac{5197}{3456}\pi^2 + \left(\frac{221}{144}\pi^2 + \frac{1}{2} \right) \ln 2 - \frac{53}{32}\zeta(3) \right] \\ & \left. - \frac{7\alpha^3}{8\pi} \ln^2 \alpha + \frac{\alpha^3}{\pi} \ln \alpha \left(\frac{17}{3} \ln 2 - \frac{217}{90} \right) + \mathcal{O}(\alpha^3) \right\} = 203.39169(41) \text{ GHz} \end{aligned}$$


 $\ln \alpha$ from would-be
 IR singularities

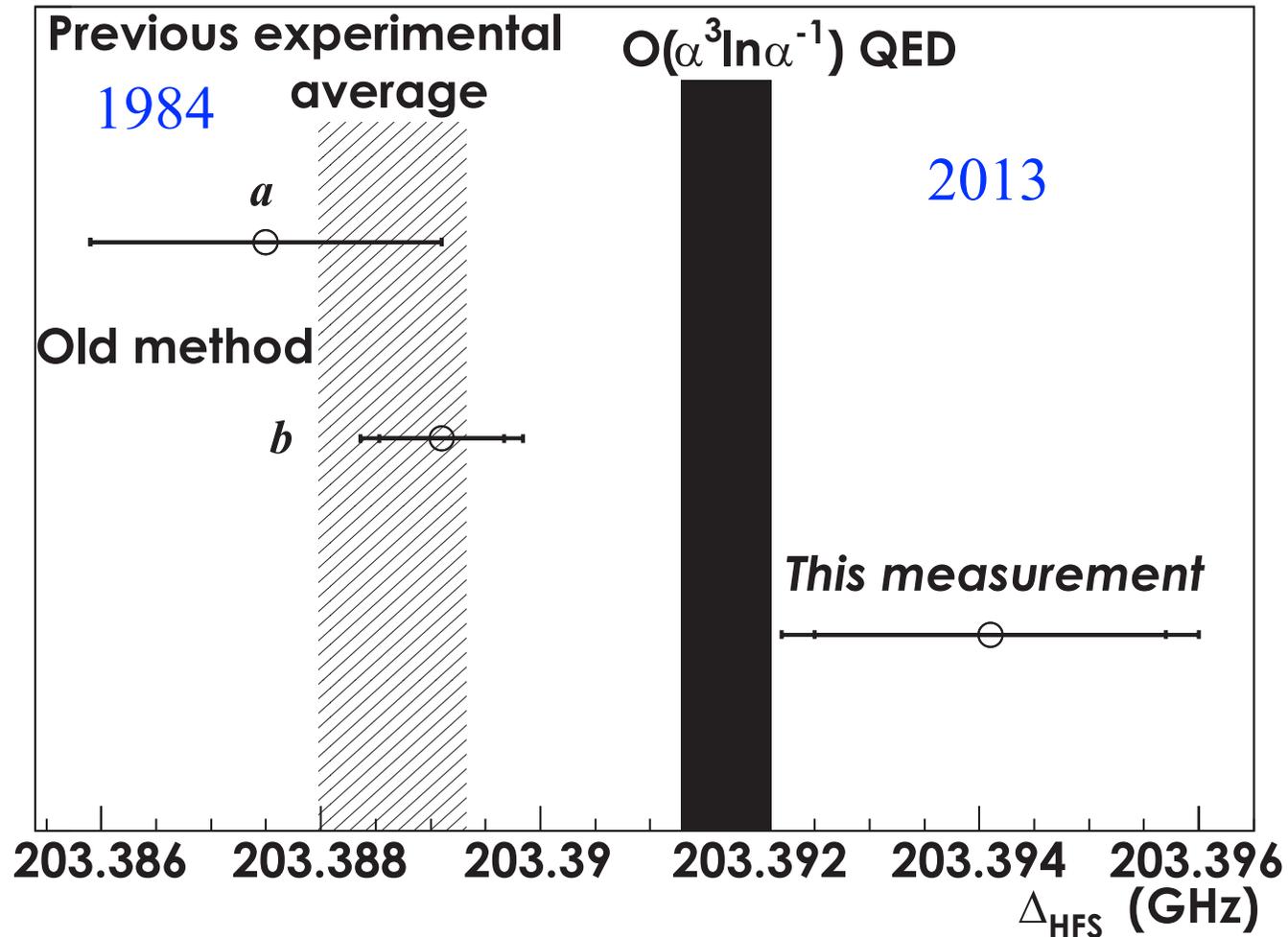
M. Baker et al, 1402.0876

QED vs Data: Hyperfine splitting in Positronium

$$\Delta\nu_{\text{QED}} = 203.39169(41) \text{ GHz}$$

$$\Delta\nu_{\text{EXP}} = 203.38865(67) \text{ GHz} \quad (1984) \quad \text{M. W. Ritter et al, Phys. Rev. A30 (1984) 1331}$$

$$\Delta\nu_{\text{EXP}} = 203.3941 \pm .003 \text{ GHz} \quad (2013) \quad \text{A. Ishida et al, PLB 734 (2014) 338 [1310.6923]}$$

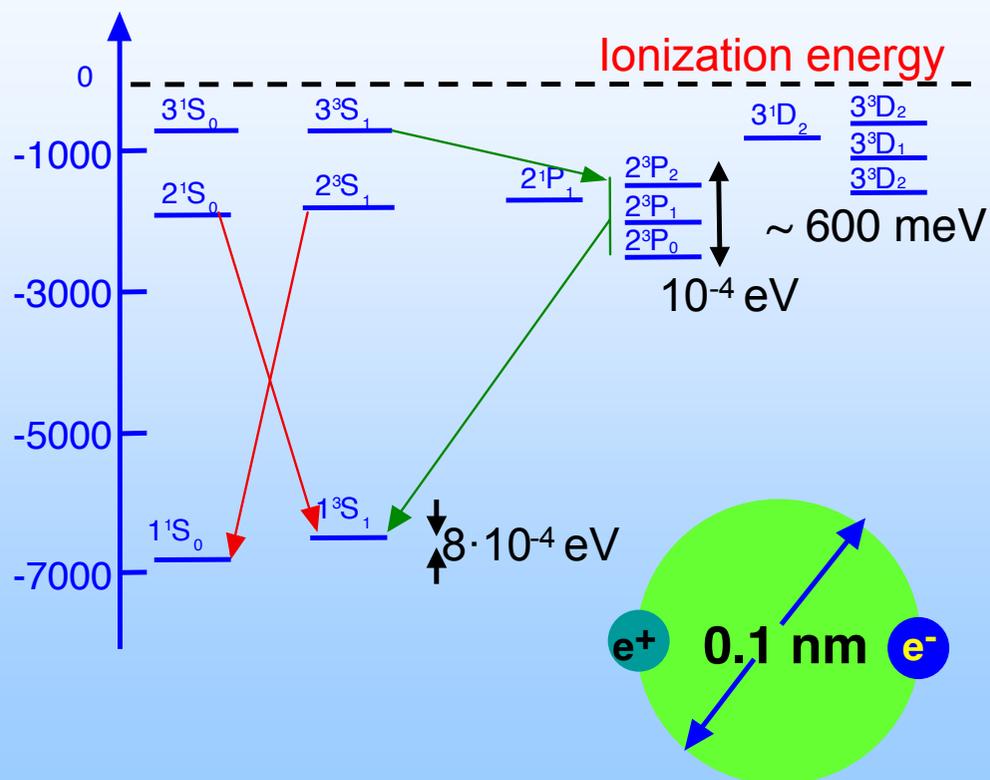


"The J/ψ is the Hydrogen atom of QCD"

QED

Binding energy [meV]

Positronium

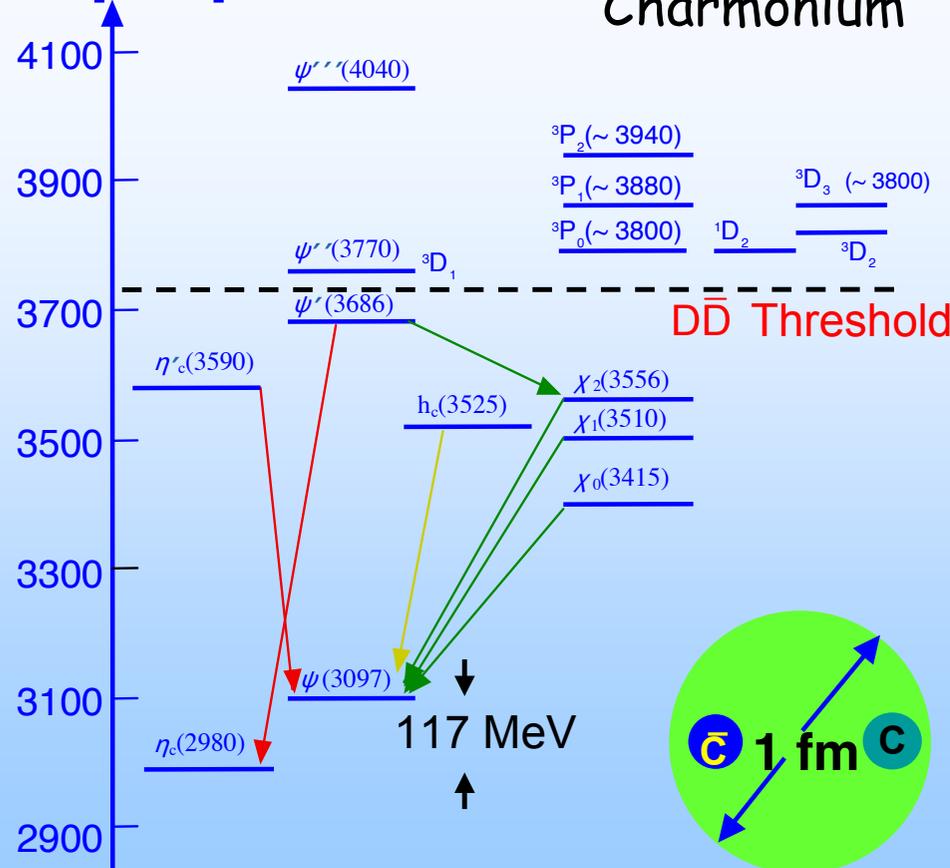


$$V(r) = -\frac{\alpha}{r}$$

QCD

Mass [MeV]

Charmonium

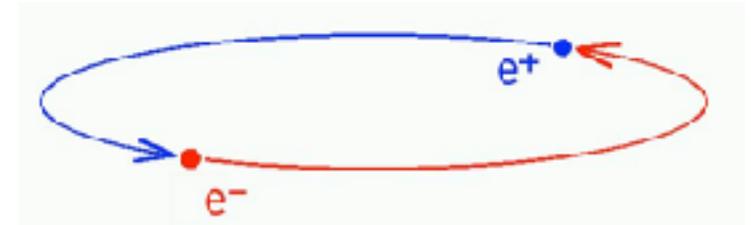


$$V(r) = cr - \frac{4}{3} \frac{\alpha_s}{r}$$

QM I: The Positronium atom

The Schrödinger equation with a classical potential is postulated:

$$\left[-\frac{\nabla^2}{2\mu} - \frac{\alpha}{|\mathbf{x}|} \right] \varphi(\mathbf{x}) = E_b \varphi(\mathbf{x})$$



Ground state
binding energy:

$$E_b = -\frac{1}{2} \mu \alpha^2 \quad \mathcal{O}(\alpha^2) \quad + \text{PQED corrections}$$

Electron velocity:

$$v = \frac{p}{m} \simeq \alpha \quad \mathcal{O}(\alpha) \quad \text{Non-relativistic}$$

Potential:

$$V(r) = -\frac{\alpha}{r} \quad \mathcal{O}(\alpha^2) \quad \text{Classical}$$

Wave function:

$$\Phi(\mathbf{x}) = N \exp(-\alpha m_e |\mathbf{x}|) \quad \mathcal{O}(\alpha^\infty)$$

QED: Atoms from Feynman diagrams

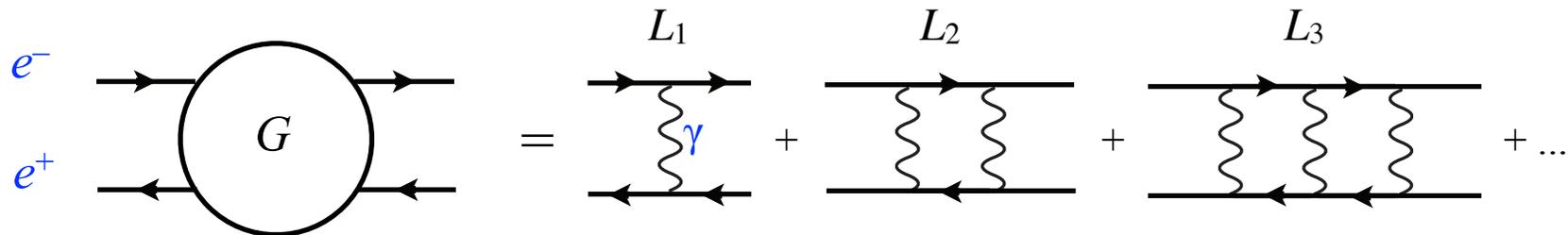
The S-matrix allows to calculate any scattering process to arbitrary order:

$$S_{fi} = \text{out} \langle f | \left\{ \text{T exp} \left[-i \int_{-\infty}^{\infty} dt H_I(t) \right] \right\} | i \rangle \text{in}$$

But: The free *in*- and *out*-states at $t = \pm\infty$ have **no overlap** with bound states.

No finite order Feynman diagram for $e^+e^- \rightarrow e^+e^-$ has a positronium pole.

Positronium poles arise from a **divergent** sum of ladder diagrams:



The divergence arises from a geometric sum of **tree diagrams** (no loops).

It gives the Schrödinger equation with **classical potential** $V(r) = -\alpha/r$.

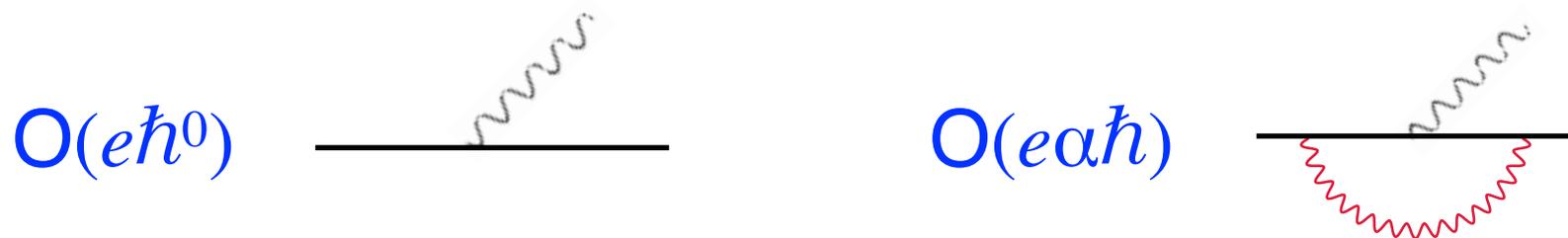
Classical fields correspond to $\hbar \rightarrow 0$

Recall functional integral formulation of QFT:

$$G(x_1, \dots, x_n) = \int [d\varphi] \exp(iS[\varphi]/\hbar) \varphi(x_1) \dots \varphi(x_n)$$

The limit $\hbar \rightarrow 0$ gives classical field eqs: $\frac{\delta S[\varphi]}{\delta \varphi} = 0$

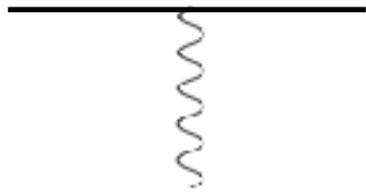
Higher orders in \hbar correspond to **loop corrections** in Feynman diagrams.



Each loop adds one power of $\alpha\hbar$: Expansions in α and \hbar are **equivalent**.

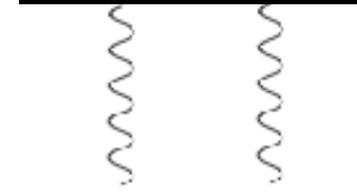
Powers of α and \hbar do not match in tree diagrams

$O(e\hbar^0)$



:

$O(e^2\hbar^0)$:



The ladders form a geometric sum of $O(\hbar^0)$ diagrams: Gives classical field.

Define Born bound states as lowest order in \hbar (no loops)

Consistent with the Schrödinger atom.

Not limited to non-relativistic dynamics.

Coupling is frozen in the absence of loops: $\alpha_s(0) \approx 0.5$

Instantaneous A^0 gauge field

The A^0 field does not propagate in time: **no** $\partial_t A^0$ term in the Lagrangian

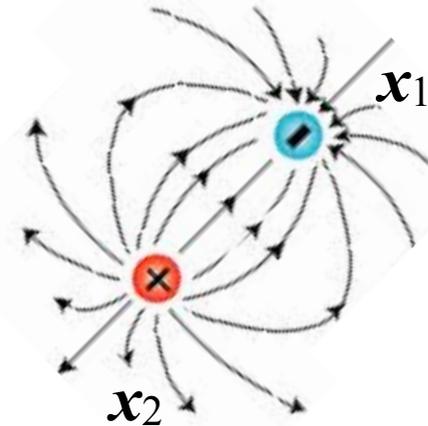
\Rightarrow Equations of motion allow to eliminate A^0 in terms of electron fields

$$\frac{\delta \mathcal{S}_{QED}}{\delta \hat{A}^0(t, \mathbf{x})} = 0 \quad \Rightarrow \quad -\nabla^2 \hat{A}^0(t, \mathbf{x}) = e \psi^\dagger(t, \mathbf{x}) \psi(t, \mathbf{x})$$

Gauss' law for the A^0 field of an $|e^-(\mathbf{x}_1) e^+(\mathbf{x}_2)\rangle$ Fock component is:

$$-\nabla^2 A^0(\mathbf{x}) = e [\delta^3(\mathbf{x} - \mathbf{x}_1) - \delta^3(\mathbf{x} - \mathbf{x}_2)]$$

$$\Rightarrow e A^0(\mathbf{x}; \mathbf{x}_1, \mathbf{x}_2) = \frac{\alpha}{|\mathbf{x} - \mathbf{x}_1|} - \frac{\alpha}{|\mathbf{x} - \mathbf{x}_2|}$$



Note: A^0 is determined **instantaneously** for all \mathbf{x} , and **depends on** $\mathbf{x}_1, \mathbf{x}_2$

The propagating vector components A^j are of higher order in α

Only fermion fields at lowest order: **No gauge fixing necessary.**

$$\begin{aligned}
 |n; \mathbf{P} = 0\rangle &= \int \frac{d\mathbf{k}}{(2\pi)^3} \varphi_n(\mathbf{k}) b_\lambda^\dagger(\mathbf{k}) d_{\bar{\lambda}}^\dagger(-\mathbf{k}) |0\rangle && \varphi_n(\mathbf{k}) : \\
 & && \text{Schrödinger wf} \\
 &= \int d\mathbf{x}_1 d\mathbf{x}_2 \bar{\psi}(t, \mathbf{x}_1)_\alpha \Phi_n^{\alpha\beta}(\mathbf{x}_1 - \mathbf{x}_2) \psi(t, \mathbf{x}_2)_\beta |0\rangle
 \end{aligned}$$

in mom. space the 4×4 wf. is:

$$\begin{aligned}
 \Phi_n^{\alpha\beta}(\mathbf{k}) &\equiv [\gamma^0 u(\mathbf{k}, \lambda)]_\alpha v_\beta^\dagger(-\mathbf{k}, \bar{\lambda}) \varphi_n(\mathbf{k}) \\
 &\propto \gamma^0 (\not{k} + m)(1 + \gamma^0) \Gamma(1 - \gamma^0) (-\not{k} + m) \varphi_n(\mathbf{k})
 \end{aligned}$$

where $(1 + \gamma^0) \Gamma(1 - \gamma^0) = \begin{pmatrix} 0 & \chi_\lambda \chi_{\lambda'}^\dagger \\ 0 & 0 \end{pmatrix}$ accounts for the 2×2 helicity d.o.f.'s of e^- and e^+

Bound state condition

$$\mathcal{H}_{QED} |n; \mathbf{P} = 0\rangle = (2m_e + E_b) |n; \mathbf{P} = 0\rangle$$

$$\mathcal{H}_{QED}(t) = \int d\mathbf{x} \psi^\dagger(t, \mathbf{x}) \left[-i\boldsymbol{\alpha} \cdot \vec{\nabla} + m\gamma^0 + \frac{1}{2}e\mathcal{A}^0(\mathbf{x}) \right] \psi(t, \mathbf{x})$$

where $\boldsymbol{\alpha} = \gamma^0\boldsymbol{\gamma}$ and $e\mathcal{A}^0(\mathbf{x}) = \int d\mathbf{y} \frac{\alpha}{|\mathbf{x} - \mathbf{y}|} \psi^\dagger(t, \mathbf{y})\psi(t, \mathbf{y})$

The factor $\frac{1}{2}$ is due to the (negative) field energy,

$$\int d\mathbf{x} \frac{1}{4} F_{\mu\nu} F^{\mu\nu} = \frac{1}{2} \int d\mathbf{x} \mathcal{A}^0 \cdot \nabla^2 \mathcal{A}^0 = -\frac{1}{2} \int d\mathbf{x} \bar{\psi}(\mathbf{x}) e\mathcal{A}^0 \psi(\mathbf{x})$$

Using $\{\psi_\alpha^\dagger(t, \mathbf{x}), \psi_\beta(t, \mathbf{y})\} = \delta_{\alpha,\beta} \delta^3(\mathbf{x} - \mathbf{y})$ and neglecting e^+e^- pair production/annihilation

the Schrödinger equation follows: $\left(-\frac{\nabla^2}{m_e} - \frac{\alpha}{|\mathbf{x}|} \right) \varphi_n(\mathbf{x}) = E_b \varphi_n(\mathbf{x})$

Positronium in motion

The Born approximation is not limited to non-relativistic dynamics.

Positronium states with CM momentum $\mathbf{P} = (0, 0, P)$ are obtained by boosts $U(\xi)$ along the z -axis, which transforms the ($t = 0$) fields as

$$U(\xi)\psi(0, \mathbf{x})U^\dagger(\xi) = e^{-\xi\alpha_3/2}\psi(z \sinh \xi, \mathbf{x})$$

Note: The transformed time depends on z .

States of any momentum \mathbf{P} are defined at **equal time**.

Boost covariance emerges via the **QED dynamics**.

The time shift can be adjusted by the generator H of time translations.

An infinitesimal boost $U(d\xi) = 1 - id\xi \mathcal{K}$ generated by

$$\mathcal{K}(t = 0) = - \int d\mathbf{x} \psi^\dagger(0, \mathbf{x}) \left[z(-i\boldsymbol{\alpha} \cdot \vec{\nabla} + m\gamma^0 + \frac{1}{2}\gamma^0 e\mathcal{A}) - \frac{1}{2}i\alpha_3 \right] \psi(0, \mathbf{x})$$

applied to a $P = 0$ Positronium state gives a state with $P = d\xi M$.

Positronium state with momentum P

In a finite boost to $P = M \sinh \xi$ the state becomes:

$$\begin{aligned}
 |n; P\rangle &\equiv U(\xi) |n; P = 0\rangle = \\
 &= \int d\mathbf{x}_1 d\mathbf{x}_2 \bar{\psi}(0, \mathbf{x}_1) e^{i\mathbf{P}\cdot(\mathbf{x}_1+\mathbf{x}_2)/2} e^{-\xi\alpha_3/2} \Phi_n^{(\xi)}(\mathbf{x}_1 - \mathbf{x}_2) e^{\xi\alpha_3/2} \psi(0, \mathbf{x}_2) |0\rangle
 \end{aligned}$$

where the wave function
is Lorentz contracted:

$$\Phi_n^{(\xi)}(x, y, z / \cosh \xi) = \Phi_n^{(0)}(x, y, z)$$

$|n; P\rangle$ is an eigenstate of the momentum \mathcal{P} and energy \mathcal{H} operators,
with the appropriate (boosted) eigenvalues.

The gauge field in the Hamiltonian applied to the boosted state,

$\mathcal{H}(\mathcal{A})U(\xi) |n; P = 0\rangle$ is the boosted version of the the rest frame A^0 field.

Summary for Positronium

- The standard PQED expression for the S -matrix in the IP is inappropriate, since Positronium has zero overlap with the free *in* and *out* states.
 - The $\mathcal{O}(\hbar^0)$ term of the PQED expansion is a (Born level) bound state.
 - The Born term involves only the instantaneous A^0 field.
 - The eqs. of motion express A^0 in terms of the electron field.
 - Eigenstates of H with any momentum \mathbf{P} can be found explicitly.
- Järvinen (2005)
- In the rest frame the result agrees with the usual Schrödinger atom.
 - The corresponding method for loop corrections remains to be developed.

Describing confinement in QCD?

The A_a^0 gluon field is instantaneous: No $\partial_t A_a^0$ term in \mathcal{L}_{QCD} .

Gauss' law for A_a^0 has **homogeneous solutions**, which do not vanish at $|\mathbf{x}| = \infty$

The boundary condition makes A_a^0 of $\mathcal{O}(\alpha_s^0)$, whereas A_a^j is of $\mathcal{O}(g)$.

At $\mathcal{O}(\alpha_s^0)$ the non-abelian contributions vanish: $f_{abc} A_a^0 A_b^0 = 0$

and Gauss' law allows to express A_a^0 in terms of the fermion fields:

$$-\nabla^2 A_a^0(t, \mathbf{x}) = g \psi_A^\dagger(t, \mathbf{x}) T_a^{AB} \psi_B(t, \mathbf{x})$$

The choice of homogeneous solutions ($\nabla^2 A_a^0(t, \mathbf{x}) = 0$) is **strongly constrained** by Poincaré invariance.

A homogeneous solution for QCD

The following is a trivial solution of $\nabla^2 \mathcal{A}_a^0(t, \mathbf{x}) = 0$ when $\kappa \neq \kappa(\mathbf{x})$:

$$A_a^0(t, \mathbf{x}) = \kappa \int d\mathbf{y} \psi_A^\dagger(t, \mathbf{y}) T_a^{AB} \psi_B(t, \mathbf{y}) \mathbf{x} \cdot \mathbf{y}$$

To preserve Poincaré invariance we need to have **color singlet** $q\bar{q}$ states:

$$|n, P = 0\rangle = \int d\mathbf{x}_1 d\mathbf{x}_2 \bar{\psi}_A(\mathbf{x}_1) \Phi_n^{AB}(\mathbf{x}_1 - \mathbf{x}_2) \psi_B(\mathbf{x}_2) |0\rangle$$

with $\Phi_n^{AB}(\mathbf{x}) = \frac{\delta^{AB}}{\sqrt{N_c}} \Phi_n(\mathbf{x})$ is a (globally) color singlet state.

Color covariance ensures that the expectation value of the field vanishes

$$\langle n, P | A_a^0(\mathbf{x}) | n, P \rangle = 0$$

Color singlet states cannot generate a color field.

The color field is **invisible to an external observer** (unlike in QED!), but confines each color component AA: $\bar{\psi}_A(t, \mathbf{x}_1) \psi_A(t, \mathbf{x}_2) |0\rangle$

Remark on $q\bar{q}$ pairs

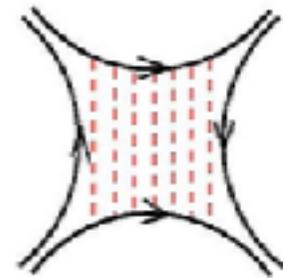
(Light) quarks in hadrons are ultra-relativistic and there is a sea of $q\bar{q}$ pairs.

Nevertheless, the hadron spectrum reflects only the valence quark dof's.

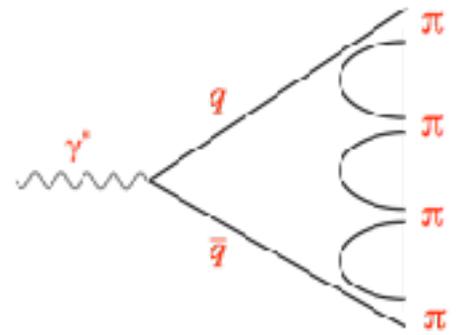
Relativistic Dirac states also have this feature (see below).

Hadron phenomenology shows that multi-particle states are dual to $q\bar{q}$:

Duality:

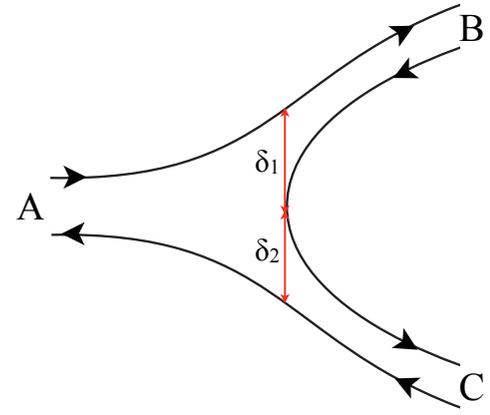


String-breaking:



In the following, the color octet $q\bar{q}$ pairs are included in the vacuum, an analogy to the Dirac case.

The singlet $q\bar{q}$ states allow to calculate (singlet) string breaking (hadron pair production, loops,...) Such effects are required by unitarity at $\mathcal{O}(\hbar^0)$.



The homogeneous A^0 solution

$$A_a^0(t, \mathbf{x}) = \kappa \int d\mathbf{y} \psi_A^\dagger(t, \mathbf{y}) T_a^{AB} \psi_B(t, \mathbf{y}) \mathbf{x} \cdot \mathbf{y} \propto \mathbf{x}$$

The field energy density is independent of \mathbf{x} for each color component AA:

$$\sum_a (\nabla A_a^0) \cdot (\nabla \hat{A}_a^0) \bar{\psi}_A(\mathbf{x}_1) \psi_A(\mathbf{x}_2) |0\rangle = C_F \kappa^2 (\mathbf{x}_1 - \mathbf{x}_2)^2 \bar{\psi}_A(\mathbf{x}_1) \psi_A(\mathbf{x}_2) |0\rangle$$

The (infinite) total field energy $\propto \int d\mathbf{x}$ must be independent of $|\mathbf{x}_1 - \mathbf{x}_2|$.

Hence

$$C_F \kappa^2 = \frac{\Lambda^4}{(\mathbf{x}_1 - \mathbf{x}_2)^2}$$

The QCD interaction Hamiltonian then gives an exactly linear potential:

$$H_{int}^{QCD} \bar{\psi}_A(\mathbf{x}_1) \psi_A(\mathbf{x}_2) |0\rangle = V(\mathbf{x}_1 - \mathbf{x}_2) \bar{\psi}_A(t, \mathbf{x}_1) \psi_A(t, \mathbf{x}_2) |0\rangle$$

where $V(\mathbf{x}) = \frac{1}{2} \sqrt{C_F} g \Lambda^2 |\mathbf{x}|$

The Fock state structure of Dirac states

Relativistically bound states have a sea of $f\bar{f}$ pairs. As an example, consider the Dirac Hamiltonian for a fixed (strong) external field $A^\mu(\mathbf{x})$,

$$H = \int d^3\mathbf{x} \bar{\psi}(\mathbf{x}) \left[-i\nabla \cdot \boldsymbol{\gamma} + m + eA(\mathbf{x}) \right] \psi(\mathbf{x})$$

There are e^- and e^+ “valence” eigenstates,

J.-P. Blaizot and PH (2015)

$$H |n\rangle = E_n |n\rangle \quad |n\rangle = \int d\mathbf{x} \psi_\alpha^\dagger(\mathbf{x}) \phi_{n\alpha}(\mathbf{x}) |\Omega\rangle \equiv c_n^\dagger |\Omega\rangle$$

$$H |\bar{n}\rangle = \bar{E}_n |\bar{n}\rangle \quad |\bar{n}\rangle = \int d\mathbf{x} \bar{\phi}_{n\alpha}^\dagger(\mathbf{x}) \psi_\alpha(\mathbf{x}) |\Omega\rangle \equiv \bar{c}_n^\dagger |\Omega\rangle$$

where the wave functions $\phi(\mathbf{x})$ and $\bar{\phi}(\mathbf{x})$ satisfy the Dirac equation

$$\left(-i\nabla \cdot \boldsymbol{\gamma} + m + eA \right) \phi_n(\mathbf{x}) = E_n \gamma^0 \phi_n(\mathbf{x}) \quad E_n > 0$$

$$\left(-i\nabla \cdot \boldsymbol{\gamma} + m + eA \right) \bar{\phi}_n(\mathbf{x}) = -\bar{E}_n \gamma^0 \bar{\phi}_n(\mathbf{x}) \quad \bar{E}_n > 0$$

The annihilation operators are Bogoliubov transforms of the free operators:

$$c_n = \sum_{\mathbf{p}} \phi_n^\dagger(\mathbf{p}) [u(\mathbf{p})b_{\mathbf{p}} + v(-\mathbf{p})d_{-\mathbf{p}}^\dagger] \equiv B_{n\mathbf{p}}b_{\mathbf{p}} + D_{n\mathbf{p}}d_{\mathbf{p}}^\dagger$$

$$\bar{c}_n = \sum_{\mathbf{p}} [b_{\mathbf{p}}^\dagger u^\dagger(\mathbf{p}) + d_{-\mathbf{p}}v^\dagger(-\mathbf{p})] \bar{\phi}_n(\mathbf{p}) \equiv \bar{B}_{n\mathbf{p}}b_{\mathbf{p}}^\dagger + \bar{D}_{n\mathbf{p}}d_{\mathbf{p}}$$

They diagonalize the Dirac Hamiltonian: $H = \sum_n [E_n c_n^\dagger c_n + \bar{E}_n \bar{c}_n^\dagger \bar{c}_n]$

The ground state $|\Omega\rangle = N_0 \exp \left[- b_{\mathbf{p}}^\dagger (B^{-1})_{\mathbf{p}\mathbf{m}} D_{\mathbf{m}\mathbf{q}} d_{\mathbf{q}}^\dagger \right] |0\rangle$

contains the e^+e^- pairs and satisfies

$$c_n |\Omega\rangle = \bar{c}_n |\Omega\rangle = H |\Omega\rangle = 0$$

Example

The Dirac states

in case of a linear potential in D=1+1 dimensions

The linear potential confines e^- , repels e^+ : $V(e^+) = -V(e^-) = -\frac{1}{2}e^2|x|$

Positrons are allowed at large $|x|$ provided they have large momenta.

The accelerating/decelerating positrons have a continuous energy spectrum.

Hence also the Dirac states have a **continuous** energy spectrum.

The Dirac Electron in Simple Fields*

By MILTON S. PLESSET

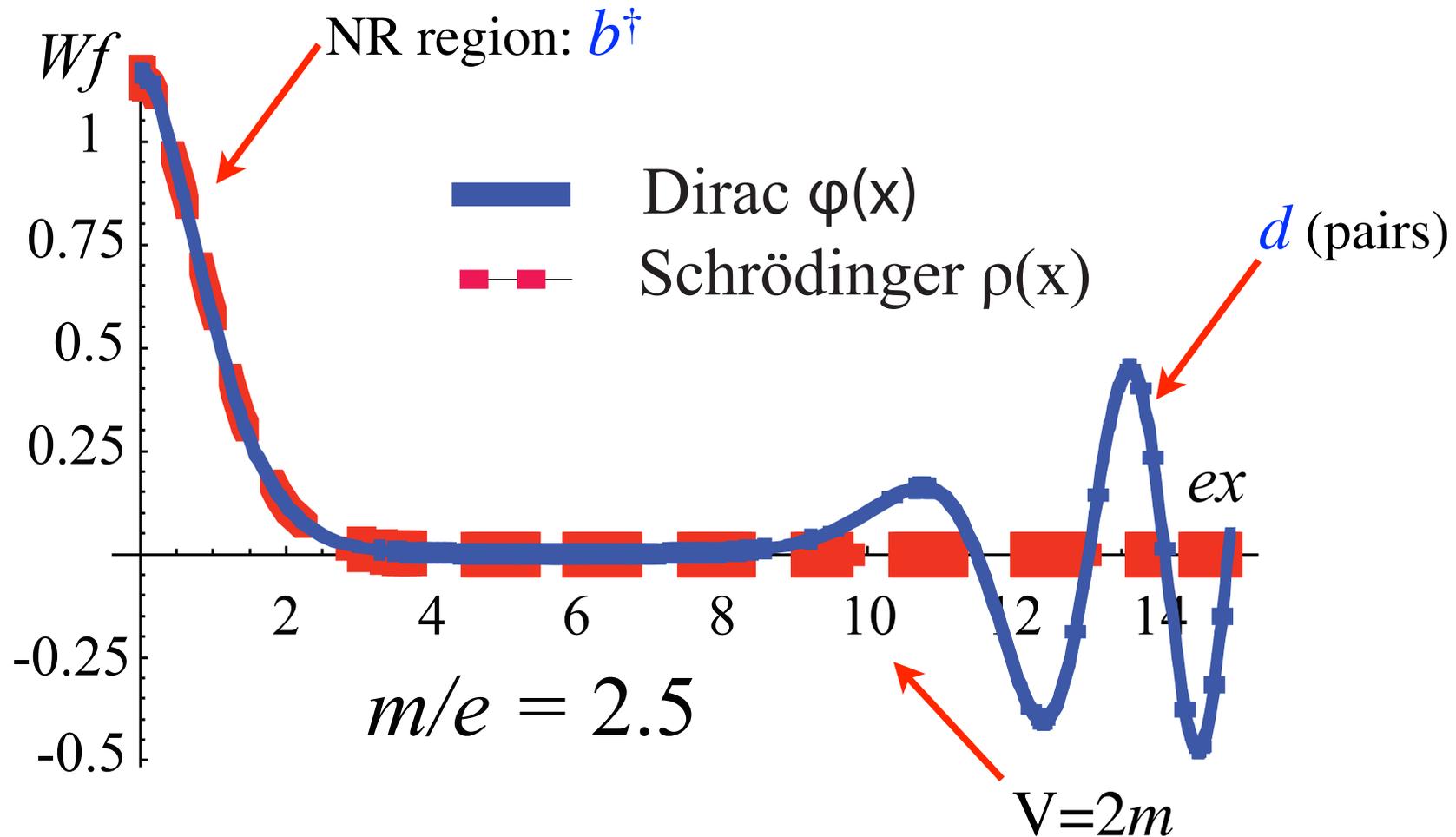
Sloane Physics Laboratory, Yale University

(Received June 6, 1932)

The relativity wave equations for the Dirac electron are transformed in a simple manner into a symmetric canonical form. This canonical form makes readily possible the investigation of the characteristics of the solutions of these relativity equations for simple potential fields. If the potential is a polynomial of any degree in x , a continuous energy spectrum characterizes the solutions. If the potential is a polynomial of any degree in $1/x$, the solutions possess a continuous energy spectrum when the energy is numerically greater than the rest-energy of the electron; values of the energy numerically less than the rest-energy are barred. When the potential is a polynomial of any degree in r , all values of the energy are allowed. For potentials which are polynomials in $1/r$ of degree higher than the first, the energy spectrum is again continuous. The quantization arising for the Coulomb potential is an exceptional case.

See also: E. C. Titchmarsh, Proc. London Math. Soc. (3) 11 (1961) 159 and 169; Quart. J. Math. Oxford (2), 12 (1961), 227.

$$|M \geq 0\rangle = \int \frac{dp}{2\pi 2E} \int dx \left[b_p^\dagger u^\dagger(p) e^{-ipx} + d_p v^\dagger(p) e^{ipx} \right] \begin{bmatrix} \varphi(x) \\ \chi(x) \end{bmatrix} |\Omega\rangle$$



The “single particle” Dirac wave function contains pair contributions (duality)

$q\bar{q}$ bound states

As previously noted, take

$$|q\bar{q}; \mathbf{P} = 0\rangle = \int d\mathbf{x}_1 d\mathbf{x}_2 \bar{\psi}_A(t, \mathbf{x}_1) \Phi^{AB}(\mathbf{x}_1 - \mathbf{x}_2) \psi_B(t, \mathbf{x}_2) |0\rangle$$

with

$$\Phi^{AB}(\mathbf{x}_1 - \mathbf{x}_2) = \frac{\delta^{AB}}{\sqrt{N_C}} \Phi(\mathbf{x}_1 - \mathbf{x}_2)$$

and

$$A_a^0(t, \mathbf{x}) = \kappa \int d\mathbf{y} \psi_A^\dagger(t, \mathbf{y}) T_a^{AB} \psi_B(t, \mathbf{y}) \mathbf{x} \cdot \mathbf{y}$$

The bound state condition $\mathcal{H}_{QCD} |q\bar{q}\rangle = M |q\bar{q}\rangle$ requires

$$i\nabla \cdot \{\gamma^0 \boldsymbol{\gamma}, \Phi(\mathbf{x})\} + m [\gamma^0, \Phi(\mathbf{x})] = [M - V(\mathbf{x})] \Phi(\mathbf{x})$$

with $V(\mathbf{x}) = \frac{1}{2} \sqrt{C_F} g \Lambda^2 |\mathbf{x}|$ (as above).

$q\bar{q}$ wave functions

The separation of angular and radial coordinates in the BSE

$$i\nabla \cdot \{\gamma^0 \boldsymbol{\gamma}, \Phi(\mathbf{x})\} + m [\gamma^0, \Phi(\mathbf{x})] = [E - V(r)] \Phi(\mathbf{x})$$

for any radial potential $V = V(r)$ and equal fermion masses $m_1 = m_2 = m$ is in:

Geffen and Suura, PRD 16 (1977) 3305

The solutions of given spin j and j_z are classified according to their charge conjugation C and parity P quantum numbers:

pion trajectory: $P = (-1)^{j+1}$ $C = (-1)^j$

a_1 trajectory: $P = (-1)^{j+1}$ $C = (-1)^{j+1}$

rho trajectory: $P = (-1)^j$ $C = (-1)^j$

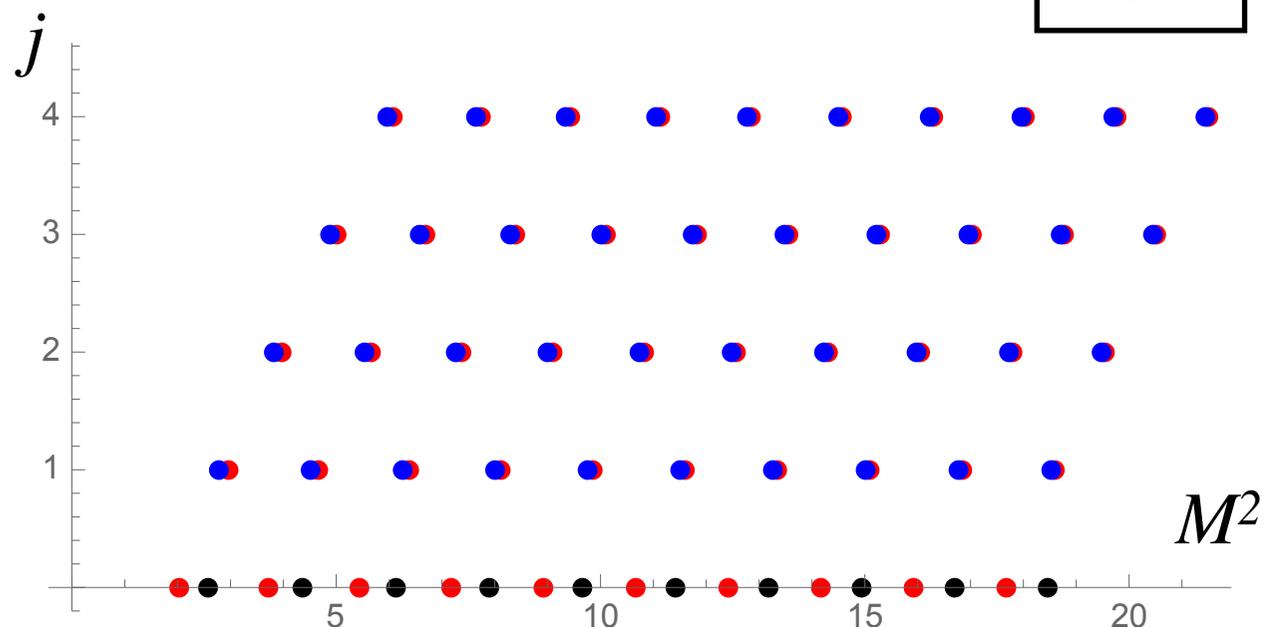
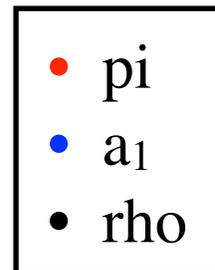
There are no “quark model exotics” with $P = (-1)^j$ and $C = (-1)^{j+1}$

π , a_1 and ρ spectra ($m = 0$)

The π , a_1 and ρ trajectories are nearly degenerate for $j \geq 1$.

There are no $j = 0$ states on the a_1 trajectory.

The $j = 0$ states are non-degenerate due to the boundary conditions



There are also $M = 0$ states.

The massless 0^{++} (σ) state has vacuum quantum numbers.

Its mixing with the chirally symmetric vacuum would cause
chiral symmetry breaking.

Summary

The similarities between quarkonia and atoms raise the question of if/how the QED bound state methods can be applied to QCD.

The perturbative expansion must be developed around an interacting state.

The Schrödinger atom is a Born level contribution: no loops, $\mathcal{O}(\hbar^0)$.

A^0 dominates A^j , and is expressed in terms of the electron field (Gauss' law).

The Hamiltonian approach allows to define atomic, equal-time states of any P .

Gauss' law has homogeneous solutions, with A^0 non-vanishing at $|\mathbf{x}| \rightarrow \infty$.

The boundary condition introduces a dimensionful constant Λ_{QCD} .

The preservation of Poincaré invariance requires to consider only color singlet states and $\nabla A^0 \neq \nabla A^0(\mathbf{x})$, giving a linear potential.

The resulting $q\bar{q}$ spectrum seems promising: Linear Regge trajectories, duality.

With Poincaré invariance, the approach is open for further studies: DIS, CSB, ..