

On quarks and the origin of QCD: Partons and baryons from intrinsic states

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Abstract – We create quarks from baryons in stead of constituting baryons from quarks. The quantum fields of QCD are generated via the exterior derivative (momentum form) of baryon wave functions on an intrinsic configuration space, the Lie group $U(3)$. Local gauge transformations correspond to coordinate translations in the intrinsic space. A proton spin structure function and a proton magnetic moment are derived. We show how the spectrum of unflavoured baryons, the N and Delta resonances, can be understood from a mass Hamiltonian on the intrinsic space and note how our model resolves the problem of colour confinement. We calculate an approximate value for the relative neutron-to-proton mass shift and give an exact value for the neutron mass. We predict neutral charge singlets that may be interpreted as neutral pentaquarks at LHCb.



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Introduction. – Nature, at the present level of our understanding, exhibits degrees of freedom that we call quarks. Quarks are described as fractionally charged spin-one-half particles of different kinds, labelled by flavour and colour. The constituent quark model for baryons like the proton and the neutron carries with it for many years a *missing resonance problem* [1,2]: many more resonances are predicted than observed¹. Quantum chromodynamics [1] carries with it a *confinement problem*: colour degrees of freedom are confined. Confinement has not yet been shown to follow from the QCD Lagrangian density. In the present work we try to solve these two seemingly independent problems by a common idea: we consider colour to live in a compact, intrinsic space, the Lie group $U(3)$ —much like a generalized spin variable with QCD as a projection. Compactness confines colour per construction.

In 1925 George Uhlenbeck and Samuel Goudsmit [4] realised that the intrinsic angular momentum, the spin, of

¹The constituent quark model was successful in predicting, *e.g.*, the Omega minus baryon at 1685 MeV [3]. Today however, the multiplet idea is mostly used *post festum* to label resonances when they are already discovered. The present model uses a compact configuration space which gives a periodic potential that lifts higher lying levels out of the “resonance domain”.

the electron represents a new, intrinsic, degree of freedom. We generalise this insight to cover also colour, flavour and electric charge. This is possible by combining the strong and electroweak interactions into a description of baryons as stationary states on an intrinsic configuration space, the Lie group $U(3)$. This space contains all three gauge groups of the standard model as subspaces which is a first motivation to choose it as a configuration space for baryons. It is compact and has nine generators equivalent to the nine kinematic generators in the laboratory space: momentum, angular momentum and Laplace-Runge-Lenz generators [5] which could explain its origin. Baryons feel both the strong interactions with $SU(3)$ symmetry and electroweak interactions with $U(1) \times SU(2)$ symmetry. We show that coordinate translations in the intrinsic space equate local gauge transformations in the laboratory space. We also show that the fundamental fields of quantum chromodynamics, quarks and gluons, are generated by the exterior derivative, the momentum form, on the intrinsic wave function. To give an intuitive picture, imagine playing “ducks and drakes” where a stone thrown at a small glancing angle scatters on a water surface and creates ripples on the surface where it hits. The

ripple patterns are quantised to fit the compact intrinsic space for mass eigenstates. The three Abelian momentum generators excite toroidal orbits which we interpret as colour degrees of freedom. The off-toroidal generators excite non-commuting degrees of freedom taking care of spin and flavour via off-toroidal derivatives in the Laplacian on $U(3)$.

The quarks in the present model are not fundamental [6] but share gauge groups with standard model quarks. Quarks are intrinsic orbits in the baryon excited in scattering by flavour generators which use quantum numbers for quark charge and baryon hypercharge as coefficients on their colour generators. The quark masses may be related to the curvature of such orbits when embedded in laboratory space, see footnote ³.

The creation of unit electric charge, *e.g.*, in the $n \rightarrow p$ decay is interpreted topologically to originate in period doublings in the intrinsic nucleon wave function. The decay relates the strong and electroweak sectors to yield equations for the electroweak energy scale and the Higgs mass in closed forms [7] and relates the electroweak mixing angle θ_W and the Cabibbo angle θ_C via quark flavour generators T_u, T_d and T_s to have $\sin^2 \theta_W \approx 1 - \text{Tr } T_u^\dagger T_d = 2/9, |\sin \theta_C| \approx |\text{Tr } T_u^\dagger T_s| = 2/9$ [8]. The u, d flavour generators also enter our treatment of the electroweak sector via a slight change in the electroweak energy scale $v = v_{\text{SM}}/\sqrt{|V_{ud}|}$. It may be a matter of definition whether this should be ascribed to a change in the electroweak interactions of quarks. We prefer to see it as a shift in interpretation by redefining electroweak coupling constants [8] even though signal strengths $\mu_{HVV}/\mu_{HVV,\text{SM}} = \frac{1}{|V_{ud}|} = 1.03$ will distinguish the viewpoints [7]. In the present work, however, we focus our attention on the strong interactions.

Baryons as intrinsic configuration states. – Baryons are considered as stationary states on an intrinsic configuration space, the Lie group $U(3)$. We identify baryon masses as the energy eigenvalues of [9]

$$\frac{\hbar c}{a} \left[-\frac{1}{2} \Delta + \frac{1}{2} d^2(e, u) \right] \Psi(u) = \mathcal{E}_0 \Psi(u), \quad u \in U(3), \quad (1)$$

with $mc^2 = \mathcal{E}_0$. The length scale a is related to the classical electron radius $r_e = \frac{e^2}{4\pi\epsilon_0 m_e c^2}$ [10,11] by the expression $r_e = \pi a$ [9] from mapping of the intrinsic baryon dynamics to the laboratory space. The factor π reflects the toroidal shape of the intrinsic configuration space $U(3)$. The baryonic energy scale $\Lambda = \hbar c/a = \frac{\pi}{\alpha} m_e c^2$ is close to the scale of non-perturbative quantum chromodynamics $\Lambda_{\text{MS}}^{(5)} \approx 210(14) \text{ MeV}$ [12]. The Hamiltonian in (1) is radically reinterpreted from lattice gauge theory [13] with a potential inspired by [14]. The unitary configuration variable $u \in U(3)$ now concerns the full baryon configuration and has eigenvalues $e^{i\theta_j}, j = 1, 2, 3$ given by three dynamical *eigenangles* $\theta_j \in \mathbb{R}$ which we interpret as colour degrees of freedom, confined per construction by the compactness of $U(3)$. These eigenangles can be used for a

polar decomposition of the Laplacian [15]

$$\Delta = \sum_{j=1}^3 \frac{1}{J^2} \frac{\partial}{\partial \theta_j} J^2 \frac{\partial}{\partial \theta_j} - \sum_{\substack{1 \leq i < j \leq 3 \\ k \neq i, j}} \frac{(S_k^2 + M_k^2)/\hbar^2}{8 \sin^2 \frac{1}{2}(\theta_i - \theta_j)}, \quad (2)$$

where [16]

$$J = \prod_{1 \leq i < j \leq 3} 2 \sin \frac{1}{2}(\theta_i - \theta_j) \quad (3)$$

and the generators S_k and M_k take care of spin and flavour, respectively. The potential in (1) is constructed from the shortest geodesic $d(e, u)$ from the neutral element e , the *origo* in the configuration space. The potential is periodic in the eigenangles and depends only on these [17]

$$\frac{1}{2} d^2(e, u) = \frac{1}{2} \text{Tr } \chi^2 = \sum_{j=1}^3 w(\theta_j), \quad u = e^{i\chi}, \quad (4)$$

where the generator

$$\chi = \theta_j T_j + (\alpha_j S_j + \beta_j M_j)/\hbar, \quad iT_j = \frac{\partial}{\partial \theta_j}; \quad \alpha_j, \beta_j \in \mathbb{R} \quad (5)$$

and (see fig. 1)

$$w(\theta) = \frac{1}{2}(\theta - n \cdot 2\pi)^2, \quad \theta \in [(2n-1)\pi, (2n+1)\pi], \quad n \in \mathbb{Z}. \quad (6)$$

The periodicity of the potential reflects the compactness of the configuration space. We see that (1) with (2) inserted is analogous to solving the hydrogen atom [18]. Now however, there are three “radial” degrees of freedom, the three eigenangles θ_j that span the Abelian (maximal) torus (7) of $U(3)$. The independence of the potential upon the remaining six dynamical variables α_j, β_j follows from the invariance of the trace on similarity operations $u \rightarrow v^{-1}uv, v \in U(3)$, in particular the ones that diagonalise u , thus

$$u \sim v^{-1}uv = \begin{pmatrix} e^{i\theta_1} & & \\ & e^{i\theta_2} & \\ & & e^{i\theta_3} \end{pmatrix}. \quad (7)$$

The length scale a in (1) can be used to map from the laboratory space to the intrinsic space by the identification, see fig. 1,

$$\theta_j = x_j/a. \quad (8)$$

We imagine the mapping in (8) to take place in scattering experiments where the intrinsic degrees of freedom are excited by the nine kinematic generators from laboratory space, namely momentum

$$p_j = \frac{-i\hbar}{a} \frac{\partial}{\partial \theta_j} = \frac{\hbar}{a} T_j, \quad (9)$$

angular momentum, *e.g.*,

$$S_1 = a\theta_2 p_3 - a\theta_3 p_2 = \hbar\lambda_7, \quad (10)$$

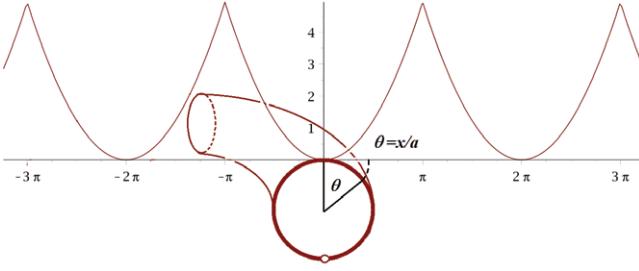


Fig. 1: Trace potential (6) in dynamical variables θ and mapping into toroidal eigenangles $\theta_j = x_j/a, j = 1, 2, 3$ from laboratory space. The length scale a is taken from the classical electron radius r_e [10,11] as $\pi a = r_e$ [9]. Figure elaborated from [19].

and Laplace-Runge-Lenz operators, *e.g.*,

$$M_1/\hbar = \theta_2\theta_3 + \frac{a^2}{\hbar^2}p_2p_3 = \lambda_6. \quad (11)$$

Here, the lambdas are the Gell-Mann generators [5].

The quantisation inherent in (9) generalises to all of the configuration space by the global concepts of left (or right) invariant coordinate fields ∂_j and corresponding coordinate forms $d\theta_j$ [20]

$$\begin{aligned} \partial_j &\equiv \frac{\partial}{\partial\theta}ue^{\theta iT_j}|_{\theta=0} = uiT_j, \quad (12) \\ \left[\theta_i, \frac{\partial}{\partial\theta_j}\right]_e &= \delta_{ij} \rightarrow d\theta_i(\partial_j) = \delta_{ij}. \end{aligned}$$

This generalises the well-known commutation relations

$$[a\theta_i, p_j] = -i\hbar\delta_{ij}. \quad (13)$$

From the coordinate representations [5] in (10) and (11)

$$[M_i, M_j] = [S_i, S_j] = -i\hbar\varepsilon_{ijk}S_k. \quad (14)$$

Note the minus sign for the spin commutators as in body fixed intrinsic coordinate systems in nuclear physics [21].

Quarks and gluons as scattering states. – In solving (1) we introduce the measure-scaled wave function [9]

$$\Phi(u) \equiv J\Psi(u) \equiv R(\theta_1, \theta_2, \theta_3)\Upsilon(\alpha_1, \alpha_2, \alpha_3, \beta_1, \beta_2, \beta_3). \quad (15)$$

The two functions R and Υ are analogues of the radial wave function and the spherical harmonics used in describing the Euclidean three-dimensional case of the hydrogen atom [18].

We consider quarks and gluons as scattering states on baryons. We create their corresponding fields ψ_j and $G^{(k)}$ via the momentum forms dR and $d\Phi$ acting on the toroidal generators iT_j and the basis $i\lambda_k$ of the adjoint representation of the $su(3)$ algebra, respectively,

$$\psi_j(u) = dR_u(\partial_j), \quad G^{(k)}(u) = d\Phi_u(\partial_k). \quad (16)$$

Here

$$\partial_k = uit_k, \quad k = 1, 2, \dots, 8 \quad (17)$$

are left invariant coordinate fields corresponding to the gluon field generators $t_k = \lambda_k/2$ [1], six of which are proportional to S_j and M_j and the remaining two are diagonal linear combinations of the T_j 's. The momentum form becomes “operational” by derivation of Φ along the direction given by the generator. For a generator Z we have the general definition of a derivation as

$$Z[\Phi](u) \equiv d\Phi_u(Z) = \frac{d}{dt}\Phi(ue^{tZ})|_{t=0}, \quad Z \in u(3), \quad (18)$$

where $u(3)$ is the algebra of $U(3)$.

From (16) applied to the ground state of (1) we have derived exemplar parton distribution functions for the valence u and d quarks of the proton [9] and we have recently derived exemplar energy-momentum tensor components of the proton too [22]. We add here as a further motivation a proton spin structure function and the proton magnetic moment. For an approximate calculation we use a Slater determinant R with period doubling in the eigenangles interpreted as a topological origin of the proton's electric charge [9]

$$R = \frac{1}{N} \begin{vmatrix} 1 & 1 & 1 \\ \sin \frac{1}{2}\theta_1 & \sin \frac{1}{2}\theta_2 & \sin \frac{1}{2}\theta_3 \\ \cos \theta_1 & \cos \theta_2 & \cos \theta_3 \end{vmatrix}. \quad (19)$$

Here N with $N^2 = \frac{3}{2}\pi^3 - \frac{44}{3}\pi$ normalises R on $\theta_j \in [0, \pi]$. The quark distribution functions derived in [9] are obtained along one-parameter curves generated by the momentum form dR applied to flavour generators T_q (23) while summing over colours and squaring to get probability distributions for the parton momentum fraction $x \in [0, 1]$

$$f_q(x)dx = \left(\sum_{j=1}^3 dR_{u=\exp(\theta iT_q)}(iT_j) \right)^2 d\theta. \quad (20)$$

Here $\theta = \pi\xi$ and the *boost* variable ξ is [9]

$$\xi \equiv \frac{\mathcal{E} - \mathcal{E}_0}{\mathcal{E}} = \frac{2 - 2x}{2 - x}, \quad \mathcal{E}_0 = mc^2, \quad (21)$$

from impacting a massless energy momentum $q = (\mathcal{E} - \mathcal{E}_0, \mathbf{q})$ on a parton xP acquiring mass $x\mathcal{E}$ in a proton at rest $(\mathcal{E}_0, \mathbf{0})$, *i.e.*,

$$(xP_\mu + q_\mu)(xP^\mu + q^\mu) = x^2\mathcal{E}^2. \quad (22)$$

For u, d, s quarks we use, respectively,

$$T_u = \frac{2}{3}T_1 - T_3, \quad T_d = -\frac{1}{3}T_1 - T_3, \quad T_s = -\frac{1}{3}T_1, \quad (23)$$

to get the unpolarised proton spin structure function averaging over three colours [19] ($f_s(x) \equiv 0$ for R in (19))

$$g_1^p(x) = \frac{1}{2} \left[e_u^2 \frac{1}{3} f_u(x) + e_d^2 \frac{1}{3} f_d(x) \right]. \quad (24)$$

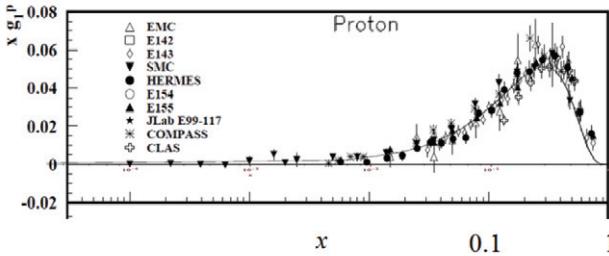


Fig. 2: Spin structure function of the proton (solid) from an exemplar calculation (24) based on flavour quarks derived from an intrinsic $U(3)$ configuration overlaid on experimental data [12] spanning four orders of magnitude. Figure taken from [8].

This represents rather well the experimentally extracted values [12] in fig. 2 without fitting parameters.

From the constituent quark model, we have an expression for the proton magnetic moment [23]

$$\mu_p = \frac{4}{3}\mu_u - \frac{1}{3}\mu_d, \quad \mu_q = \frac{e_q \hbar/2}{4\pi \varepsilon_0 m_q}. \quad (25)$$

We use [19]

$$m_p = 2m_u + m_d, \quad \frac{m_u}{m_d} = \frac{\int_0^1 x f_u(x) dx}{\int_0^1 x f_d(x) dx} = \frac{0.2722}{0.1432} \quad (26)$$

and get $m_u c^2 = 371$ MeV, $m_d c^2 = 195$ MeV from which

$$\mu_p = 2.779 \dots \mu_N, \quad \mu_N = \frac{e \hbar/2}{4\pi \varepsilon_0 m_p}, \quad (27)$$

to compare with experiment $2.7928473446(8)\mu_N$ [1].

Basic building blocks for QCD construction.

We require local gauge invariance of a field Hamiltonian [24] constructed from the colour quark fields generated in (16)²

$$H = \int \psi^\dagger (-i\hbar c \boldsymbol{\alpha} \cdot \boldsymbol{\nabla} + \beta mc^2) \psi dx^3, \quad \psi^\dagger = (\psi_1^*, \psi_2^*, \psi_3^*). \quad (28)$$

Here we suppressed spinor indices which are mixed by the 4×4 Dirac matrices $\boldsymbol{\alpha} = (\alpha_1, \alpha_2, \alpha_3)$ and β . The spinor indices commute with the colour indices. Using left invariance of the coordinate fields $\partial_j|_u = uiT_j = u\partial_j|_e$ from (12) we get in the mass term of (28)³

$$\begin{aligned} \psi'(u')^\dagger \psi'(u') &= (u'iT_j[R])^\dagger (u'iT_j[R]) = \\ &= (iT_j[R])^\dagger (u')^\dagger u' (iT_j[R]) = \psi(u)^\dagger \psi(u), \end{aligned} \quad (29)$$

²This section is edited from [25].

³Quark masses are *ad hoc*. One may speculate that the finite values of current masses m_u, m_d represent a topological tension from mapping the curved intrinsic group to the flat algebra in laboratory space, somewhat like our interpretation of the electron as a peel-off from the neutron in the $n \rightarrow p$ decay [9] which scales (1) by $\pi a \equiv r_e$ via the strength of the electric charge in the classical electron radius, $m_e c^2 = e^2/(4\pi \varepsilon_0 r_e) = \alpha \frac{\hbar c}{r_e}$. With flavour distributed over three colour charges, by analogy, $m_q c^2 = (g_s/3)^2/(4\pi) \frac{\hbar c}{r_q} = \frac{\alpha_s}{9} \frac{\hbar c}{r_q}$. An interpretation of $1/r_q$ as a curvature is supported by treating the u, d, s quark mass matrix M as an external field, cf. sect. 59 in [1]

provided the configuration variables u', u are unitary, *i.e.*, $(u')^\dagger u' = u^\dagger u = \mathbf{1}$. Next we impose the local gauge transformation

$$\psi \rightarrow \psi' = g(x)\psi, \quad g(x) \in SU(3), \quad \partial_\mu \rightarrow D_\mu = \partial_\mu + \mathcal{G}_\mu, \quad (30)$$

with colour gauge field, $\mathcal{G} = ig_s \mathbf{G}$ [1] containing a strong coupling g_s

$$\mathcal{G}_\mu = ig_s G_\mu^k t_k, \quad k = 1, 2, \dots, 8, \quad G^k \sim G^{(k)}(e) = d\Phi_e(it_k) \quad (31)$$

and transforming (when $e \rightarrow g(x)$ in $G^{(k)}$) like in [27]

$$\begin{aligned} \mathcal{G}'_\mu &= g(x)\mathcal{G}_\mu g(x)^{-1} - \partial_\mu(g(x))g(x)^{-1} \rightarrow \\ &= (D'_\mu \psi')^2 = (D_\mu \psi)^2. \end{aligned} \quad (32)$$

We thus have the basic ingredient colour fields for setting up QCD. Spin degrees of freedom enter from (10) and flavours enter from (11) and are extracted by (23). Choosing $u = g(x)$ in (30) and in (12) equates local gauge transformation in laboratory space to left translation of the intrinsic coordinate fields

$$\psi_j(u) = \partial_j|_u[\Phi] = u\partial_j|_e[\Phi] = u\psi_j(e). \quad (33)$$

Unflavoured baryons: N and Δ states. – In (1) we multiply by J and integrate over the off-toroidal degrees of freedom to get [9]

$$\left[-\frac{1}{2} \sum_{j=1}^3 \frac{\partial^2}{\partial \theta_j^2} + W(\boldsymbol{\theta}) \right] R(\boldsymbol{\theta}) = ER(\boldsymbol{\theta}), \quad \boldsymbol{\theta} = (\theta_1, \theta_2, \theta_3), \quad (34)$$

for the dimensionless eigenvalues $E = \mathcal{E}_0/\Lambda$. The total potential

$$W(\boldsymbol{\theta}) = -1 + \sum_{1 \leq i < j \leq 3}^3 \frac{4/3}{16 \sin^2 \frac{1}{2}(\theta_i - \theta_j)} + \sum_{j=1}^3 w(\theta_j) \quad (35)$$

is periodic with w from (4). The constant term -1 follows from differentiating through J in the Laplacian (2) [15]; Dowker calls it the constant global curvature potential [28]. The centrifugal term has the numerator $4/3$ from the minimum value $(\mathbf{S}^2 + \mathbf{M}^2)/\hbar^2 = 4$ [19] common for N and Δ in integrating the second term in (2) over the six off-toroidal degrees of freedom contained in Υ from (15). Here the arbitrary labelling of the three eigenangles is exploited to make an average over the non-commuting components of \mathbf{S} and \mathbf{M} in (2). For a Wilson

—quite analogous to the relation in general relativity between the curvature of the metric field and the energy-momentum density in spacetime. In stead of (8), we may embed the 2D tori generated by $T_{u,d}$ from (23) in laboratory space to have $(x, y, z) = (a(1 + \cos e_q \theta_1) \cos(-\theta_3), a(1 + \cos e_q \theta_1) \sin(-\theta_3), a \sin e_q \theta_1)$. Using [26] we find Gaussian curvatures $K = \frac{\cos e_q \theta_1}{a e_q (a + a \cos e_q \theta_1)}$ with e_q in units of e .

Integration with R^2 from (19) over $[0, \pi]^3$ yields $\frac{1/r_d}{1/r_u} = \sqrt{\frac{\langle K_d \rangle}{\langle K_u \rangle}} = 2.19$ to compare with $m_d/m_u = 2.16(40)$ [1].

analogue [29], which is more commonly used, w would read $w_{\text{Wilson}}(\theta) = (1 - \cos \theta)$. However, the Wilson analogue does *not* represent well even the lowest-lying states if used in (34) in stead of the Manton-inspired potential we use here.

The total potential W has periodicity 2π in all three eigenangles. This opens for the introduction of Bloch wave degrees of freedom. The measure-scaled toroidal wave function R is antisymmetric under interchange of the three colour degrees of freedom θ_j and we may therefore expand it on Slater determinants [19] (see Supplementary Material `Supp1_1Deigenvalues_Matlab.m` and `Supp1a_1Deigenvalues_Matlab.pdf` (SM1), `Supp2-1DbaseFunctionSet-Maple.mw` and `Supp2a-1DbaseFunctionSet-Maple.pdf` (SM2), `Supp3-3DchargedN-Mathcad.mcd` and `Supp3a-3DchargedN-Mathcad` (SM3))

$$g_{lmn} = \begin{vmatrix} \phi_l(\theta_1) & \phi_l(\theta_2) & \phi_l(\theta_3) \\ \phi_m(\theta_1) & \phi_m(\theta_2) & \phi_m(\theta_3) \\ \phi_n(\theta_1) & \phi_n(\theta_2) & \phi_n(\theta_3) \end{vmatrix}, \quad l, m, n \in \mathbb{Z}^+, \quad (36)$$

constructed from solutions ϕ_i with periodicity either 2π or 4π to the one-dimensional problem

$$\left[-\frac{1}{2} \frac{d^2}{d\theta^2} + w(\theta) \right] \phi_i(\theta) = e_i \phi_i(\theta). \quad (37)$$

These correspond to Bloch wave functions [30]

$$\phi(\theta) = e^{i\kappa\theta} b(\theta), \quad (38)$$

with b carrying the 2π periodicity of the potential and $\kappa = 0$, $\kappa = \pm\frac{1}{2}$, respectively. No other values of κ are allowed since the square of the wave function must be single-valued on the configuration space to keep up its probability interpretation. Note that in (37) there is no summation over i . Table 1 shows the first eigenvalues e_i for real-valued wave functions emulating Bloch wave vectors $\kappa = 0$ (2π periodicity) and e'_i for $\kappa = \pm\frac{1}{2}$ (4π periodicity). The eigenvalues are found by collocation (see SM1). We get an eigenvalue $E = e_l + e_m + e_n$ for (36) under an approximate, separable Hamiltonian⁴

$$\left[\sum_{j=1}^3 \left(-\frac{1}{2} \frac{\partial^2}{\partial\theta_j^2} + w(\theta_j) \right) \right] g_{lmn} = E g_{lmn}. \quad (39)$$

Combining three different eigenvalues from table 1 gives the resonance structure shown in fig. 3 for maximally charged states, *i.e.*, N^+ -like and Δ^{++} -like states. We note in passing the resulting value for the relative neutron-to-proton mass shift [31],

$$\frac{m_n - m_p}{m_p} \approx \frac{e_1 + e_2 + e_3 - (e'_1 + e'_2 + e'_3)}{e'_1 + e'_2 + e'_3} = 0.13847(14)\%, \quad (40)$$

⁴The eigenvalues of this Hamiltonian are closer to those involving the full potential (35) than one might expect. This is because the total curvature term and the centrifugal terms have opposite signs and more or less cancel each other when integrated [31].

Table 1: 1D eigenvalues (37) to construct the approximate baryon spectrum in fig. 3 from (36). The eigenvalues are calculated with 1500 collocation points (sm1). The lowest eigenvalues, as expected, are close to those of the ordinary harmonic oscillator. Moving up to higher levels, the eigenvalues increase quadratically as indicated in fig. 4. Table extract from [19].

i Level	e_i Eigenvalue	e'_i Diminished	e'_i Augmented
1	0.499804708		0.5001727904
2	1.502988968	1.496433950	
3	2.471378779		2.522629649
4	3.600509000	3.377236032	
5	4.218515963		4.803947527
...

to compare with the observed value $0.13784190(5)\%$ [1]. Note that we combine two period doublings in the choice of κ 's for the protonic state. The choices are shown with black dots in fig. 4. The same kind of coupling between Bloch wave vectors is necessary when the full potential W in (35) is to be used for exact solutions of (1) since then the singularity of the centrifugal term can be cancelled. Unfortunately we have not succeeded in finding a complete base on which to expand protonic states for solving with analytical integrals. Our best suggestion so far is [19,25]

$$f_{pqr} = \begin{vmatrix} \cos p\theta_1 & \cos p\theta_2 & \cos p\theta_3 \\ \sin q\theta_1 & \sin q\theta_2 & \sin q\theta_3 \\ \cos r\theta_1 & \cos r\theta_2 & \cos r\theta_3 \end{vmatrix} \in R_n \quad \text{and} \\ \frac{1}{2} \left(e^{\frac{i}{2}(\theta_1+\theta_2+\theta_3)} + e^{-\frac{i}{2}(\theta_1+\theta_2+\theta_3)} \right) f_{pqr} \in R_p \quad (41)$$

with non-negative integers $0 \leq p < r$ and $0 < q$. For R_n the set is complete and yields the very accurate neutron-to-electron mass ratio [9,32] (see Supplementary Material `Supp4-3DneutralN-Mathcad.mcd` and `Supp4a-3DneutralN-Mathcad.pdf` (SM4))

$$\frac{m_n}{m_e} = \frac{\pi}{\alpha(m_n)} E_n = 1839(1), \quad E_n = 4.382(2), \quad (42)$$

in agreement with the observed value [1]. However, for the protonic state R_p we have not settled how to select the p, q, r -triples to keep completeness and avoid overcompleteness. An abstract crystallographic approach is needed; Jones [33] seems appropriate.

Charge singlet states: neutral pentaquarks? – A particularly interesting set of states with no charged partners and parity opposite to those from (41) are

$$f_{pqr}^0 = \begin{vmatrix} \cos p\theta_1 & \cos p\theta_2 & \cos p\theta_3 \\ \cos q\theta_1 & \cos q\theta_2 & \cos q\theta_3 \\ \cos r\theta_1 & \cos r\theta_2 & \cos r\theta_3 \end{vmatrix}, \quad 0 \leq p < q < r \in \mathbb{N}. \quad (43)$$

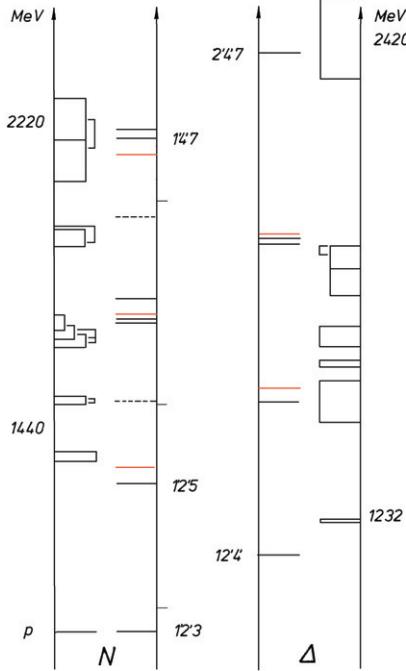


Fig. 3: Unflavoured baryon spectra. The dashed lines are singlet states. The red lines mark states with augmented contribution in level 3. All lines are approximate predictions based on (39). The boxes represent baryons observed with certainty [1]. The box widths represent the uncertainty in the mass pole peaks, not resonance widths, which are much larger. Figure updated from [31].

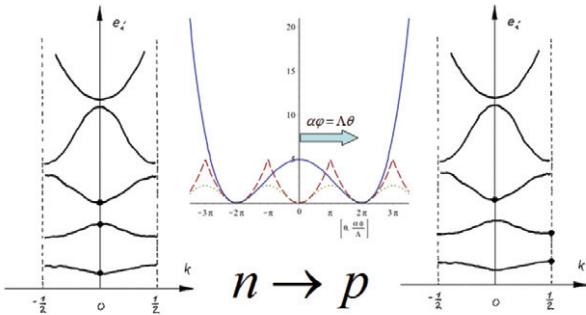


Fig. 4: Left and right: reduced zone schemes, cf. [30], for Bloch wave numbers for the neutron state (left) and the proton state (right). Middle: Higgs potential (solid, blue) matching the Manton-inspired potential [14] (dashed, red) and the Wilson-inspired potential [29] (dotted, green). The Manton and Wilson inspired potentials yield the same value for the Higgs mass and the electroweak energy scale [19] whereas only the Manton inspired potential (6) gives a satisfactory reproduction of the baryon spectrum seen in fig. 3. Figure adapted from [34].

Like for the neutral states R_n in (41) also the Hamiltonian in (1) for (43) can be diagonalized with a Rayleigh-Ritz method [19,32] where the integrals needed for the Hamiltonian matrix elements can be found analytically and the

Table 2: Scarce singlet states. Eigenvalues based on Slater determinants (43) of three cosines up to order 20 (see SM5). The first column shows eigenvalues from the approximate Hamiltonian (39) and the third column shows eigenvalues of the exact equation (34). The rest masses are predicted from a common fit of the neutron state 939.57 MeV to the ground state 4.382 of (34) with no period doublings. The four resonances marked by an asterisk (*) lie within the observational window in Λ_b^0 decays (44) at LHCb. Table updated from [31].

Singlet approximate (39)	Toroidal label	Singlet exact (34)	Rest mass MeV/ c^2
7.1895	1 3 5	7.1217	1527
...
18.9214	1 5 11	19.7327	4231
20.3774	5 7 9	20.9940	4501*
20.8910	3 5 11	21.7110	4655*
21.0766	1 7 11	22.0409	4726*
23.0609	3 7 11	23.7887	5101*
24.4575	1 9 11	23.9981	5146
...

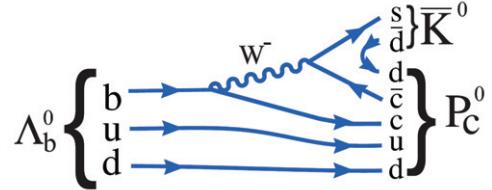


Fig. 5: Feynman diagram for neutral pentaquark formation (44). Figure adapted from [35].

eigenvalues therefore be found with high accuracy. The eigenvalues for these cases are given in table 2 (see Supplementary Material `Supp5-3Dsinglets-Mathcad.mcd` and `Supp5a-3Dsinglets-Mathcad.pdf` (SM5)).

We have previously suggested to look for neutral charge resonances in [31] and had the opportunity to discuss the possibilities at LHCb with Sheldon S. Stone at the EPS-HEP 2015 after Marta Calvi had been so kind as to forward our request. Sheldon Stone suggested the following channel (when enough data are acquired) [36]:

$$\Lambda_b^0 \rightarrow \bar{K}^0 + P_c^0 \rightarrow \bar{K}^0 + J/\Psi + \Delta^0 \rightarrow \bar{K}^0 + J/\psi + p + \pi^- \quad (44)$$

Figure 5 shows a quark structure interpretation for P_c^0 production in Λ_b^0 decay which can be reached at LHCb. Other channels could be narrow resonances in photoproduction on neutrons, in $\pi^- p$ scattering and in invariant mass spectra of $\Sigma_c^+(2455)D^-$ from decays.

Conclusion. – We derived quark and gluon fields for QCD from baryonic states on an intrinsic $U(3)$ Lie group configuration space with a mass Hamiltonian. We have shown in general that the intrinsic variable must be unitary for the mass term of the field Hamiltonian to be

invariant under translations in the intrinsic space and that local gauge transformations in laboratory space correspond to left translations in configuration space. As applications of the momentum form, we derived a proton spin structure function and a proton magnetic moment. We used the conjugacy of coordinate fields and coordinate forms as a generalisation of action-angle quantisation in ordinary quantum mechanics. We hinted at a topological origin of quark masses which should be better understood.

We have shown how to derive unflavoured baryon spectra without explicit introduction of quarks and gluons. Colour degrees of freedom are carried by the toroidal degrees of freedom and spin and flavour by off-toroidal derivatives in the Laplacian on $U(3)$. The unflavoured spectrum from approximate calculations compares well with observed four star resonances. We have given an approximate value for the relative neutron to proton mass shift and an accurate value for the neutron mass. We predict unflavoured neutral charge singlets that might be interpreted as neutral pentaquarks.

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