# **Calculating the proton radius using lattice QCD**

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#### **Proton radius**

The distributions of charge and magnetization in a proton are probed in elastic electron-proton scattering.

Photon-proton vertex is parameterized by two form factors,

 $\langle p'|J_{\mu}|p\rangle = \bar{u}(p')\left[\gamma_{\mu}F_{1}(Q^{2}) + \frac{i\sigma_{\mu\nu}(p'-p)^{\nu}}{2m}F_{2}(Q^{2})\right]u(p), \ Q^{2} = -(p'-p)^{2}.$ 

These combine to form the electric and magnetic form factors,

 $G_E(Q^2) = F_1(Q^2) - \frac{Q^2}{4m^2}F_2(Q^2) \rightarrow$  Fourier transform of charge density  $G_M(Q^2) = F_1(Q^2) + F_2(Q^2) \rightarrow$  Fourier transf. of magnetization density. Near  $Q^2 = 0$  they contain key properties of the proton:

> electric charge  $1 = G_E(0)$ rms charge radius  $r_E^2 = -6 \frac{dG_E}{dO^2}$

# **Fitting form factors**

 $Q^2$ 

Finite-volume momentum transfers take discrete values:  $Q_{\min}^2 \approx (\frac{2\pi}{L})^2$ . Similar to scattering experiments, can determine radius by fitting. We use the *z* expansion, which conformally maps the domain for complex  $Q^2$  where  $G(Q^2)$  is analytic to |z| < 1, then uses a Taylor series:

#### e.g. R. J. Hill and G. Paz, Phys. Rev. D 84, 073006 (2011)

Rather than simply truncating the series, we impose Gaussian priors on the higher coefficients  $a_k$ , k > 1.

#### **Directly calculating at** $Q^2 = 0$

Imposing twisted boundary conditions on the quarks shifts the Fourier momenta:









Lattice quantum chromodynamics (QCD)					
	mass → ≈2.3 MeV/c <sup>2</sup>	≈1.275 GeV/c²	≈173.07 GeV/c²	0	≈126 GeV/c²
	charge $\rightarrow 2/3$	<sup>2/3</sup> C	2/3	0	• •
	spin $\rightarrow 1/2$	1/2	1/2	1 9	0
OCD, the theory of quarks and	au	charm	top	aluon	Higgs

$$q_{\theta}(\vec{x}) = e^{i\theta}q_{\theta}(\vec{x}+\hat{j}L) \implies p_{j} = \frac{2\pi n + \theta}{I}, \ n \in \mathbb{Z}.$$

For connected diagrams, we can use a vector current  $J_{\mu} = \bar{q}_{\theta'} \gamma_{\mu} q_{\theta}$  that transitions between different twist angles. This allows for arbitrary adjustment of  $p'_j - p_j$  and thus arbitrary  $Q^2 \ge 0$ . Furthermore, it has been shown how to take  $\frac{\partial}{\partial \theta}$  analytically. G. M. de Divitiis, R. Petronzio, N. Tantalo, Phys. Lett. B 718, 589 (2012) Using the derivative method, we expect that radii can be computed up to

 $O(e^{-m_{\pi}L})$  finite-volume effects. B. C. Tiburzi, Phys. Rev. D 90, 054508 (2014)

We applied this to isovector form factors:

N. Hasan *et al.*, Phys. Rev. D **97**, 034504 (2018)



Results for the radius were quite noisy. This motivated a new, mixed-derivative approach:

 $\frac{\partial^2}{\partial p'_i \partial p_i} \langle p' | J_\mu | p \rangle \longrightarrow r_1^2$ 

gluons, is the elementary theory that describes protons and neutrons.

We perform calculations using the lightest three quarks: *u*, *d*, and *s*.

The proton has *net* quark content *uud* and the neutron *udd*.





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Lattice QCD is a way of regularizing
Euclidean-space QCD on a 4d grid so that the
quantum path integral becomes finite-dimensional.
▶ Quark fields q, q live on lattice sites.

Solution Gluon field becomes gauge links  $U_{\mu}$  between adjacent sites.

Lattice action has form  $S[q, \bar{q}, U] = S_g[U] + \bar{q}D[U]q$ . We integrate quarks analytically to get  $S_{\text{eff}}[U] = S_g[U] - \log \det D[U]$ .

Path integral can have >  $10^8$  dimensions  $\rightarrow$  use Monte Carlo methods: generate stochastic samples  $U_i$  from distribution  $p[U] \propto e^{-S_{\text{eff}}[U]}$ . On a cluster, it is natural to split the lattice into sublattices, each of which is contained on one MPI rank. Much of the work is spent on repeatedly

### **Preliminary results: isovector form factors**

We use one 48<sup>4</sup> lattice ensemble with  $m_u = m_d$  and  $m_s$  set close to their physical values, and lattice spacing a = 0.116 fm.



solving  $D[U_i]\psi = \eta$  to compute the quark propagator  $\psi$  from a source  $\eta$ , on gauge fields  $U_i$  that sample the path integral.

#### **Protons and neutrons on the lattice**

Compute two-point and three-point functions.

connected: can compute straightforwardly



disconnected: require additional stochastic estimation

Using ratios  $C_{3pt}/C_{2pt}$  at large time separations (where excited states have died out), we can isolate  $\langle p'|J_{\mu}|p \rangle$  and then get  $G_E$  and  $G_M$ . Disconnected diagrams cancel out in the *isovector* (proton minus neutron) form factors.

## **Preliminary results: disconnected diagrams**

We use hierarchical probing to efficiently estimate the near-diagonal elements of  $D[U]^{-1}$ . For light quarks, we improve this by treating the low-lying modes of  $D^{\dagger}[U]D[U]$  exactly.

