Introduction to Lattice QCD

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Standard model

The Standard Model (SM) is a synthesis of three of the four forces of nature described by gauge theories with coupling constants:

- Strong Interactions: $\alpha_s \sim 1$
- Electromagnetic interactions: $\alpha_{em} \approx 1/137$
- Weak interactions: $G_F \approx 10^{-5} \text{ GeV}^{-2}$.

Basic constituents of matter:

- Six quarks, u, d, s, c, b, t, each in 3 colors, and six leptons e, ν_e , μ , ν_μ , τ , ν_τ
- The quarks and leptons are classified into 3 generations of families.
- The interactions between the particles are mediated by vector bosons: the 8 gluons mediate strong
 interactions, the W[±] and Z mediate weak interactions, and the electromagnetic interactions are carried
 by the photon γ.
- The weak bosons acquire a mass through the Higgs mechanism.
- The SM is a local gauge field theory with the gauge group SU(3) × SU(2) × U(1) specifying the interactions among these constituents.
 Three Generations

			of Matter (Fermions)				
Masses in the Standard Model				1	Ш	III	
Parameters	Number	Comments	mass- charge-	2.4 MeV 3/3	1.27 GeV 3/3	171.2 GeV 2/3	0 17
Masses of quarks	6	u, d, s light	spin-	, _{1/2} U	- ^{4/3} C	1/2	1 Y
•		c, b heavy	name	• up	charm	top	photon
		$t = 175 \pm 6 \; GeV$		4.8 MeV	104 MeV	4.2 GeV	0
			ş	1/3 C	1/3 S	-1/3 b	$\bigcup_{i=1}^{6} \mathbf{G}$
Masses of leptons	6	e, μ, au	Quarks	down	strange	bottom	gluon
		$M_{ u_{\theta},\ u_{\mu},\ u_{ au}}$ non-zero		<2.2 eV	<0.17 MeV	<15.5 MeV	91.2 GeV O
		σ, μ, ι		° V _e	$^{\circ}_{\frac{1}{2}}V_{\mathfrak{u}}$	$V_{1/2}$	$^{\circ}_{1}$ 7
Mass of W^{\pm}	1	80.3 GeV		electron neutrino	muon neutrino	tau neutrino	weak force
Mass of Z	1	91.2 GeV		0.511 MeV	105.7 MeV	1.777 GeV	80.4 GeV +
Mass of gluons, γ		0 (Gauge symmetry)	eptons	1/2 e	⁻¹ μ	¹½ T	
		.0.00		electron	muon	tau	weak force
Mass of Higgs	1	$125.03^{+0.26}_{-0.27} \text{ (stat)}^{+0.13}_{-0.15} \text{ (sys)}$ GeV disco	overed	d at LH	PC, 20	12	

QCD - Gauge theory of the strong interaction

Lagrangian: formulated in terms of quarks and gluons

$$\begin{split} \mathcal{L}_{\text{QCD}} &= -\tfrac{1}{4} F^a_{\mu\nu} F^{a\,\mu\nu} + \sum_f \overline{\psi}_f \left(i \gamma^\mu D_\mu - m_f \right) \psi_f, \quad f = u, d, s, c, b, t \\ D_\mu &= \partial_\mu - i g(\tfrac{1}{2} \lambda^a) A^a_\mu \end{split}$$



Harald Fritzsch



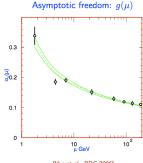
Murray Gell-Mann



Heinrich Leutwyler

Phys.Lett. B47 (1973) 365

Properties of QCD



[Yao et al., PDG 2006]

continuum limit $\beta = 6.92$

Confinement

 $$\rm r/r_{\rm 0}$$ [Necco & Sommer, Nucl Phys B622 (2002) 328]

1.5

0.5

Nobel Prize in Physics 2004

"...for the discovery of asymptotic freedom in the theory of the strong interaction"

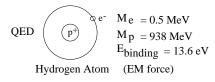


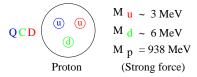


Frank Wilczek David Politzer

QCD versus QED

- QCD is the theory of strong interactions formulated in terms of quarks and gluons as the basic degrees of freedom of hadronic matter.
- Conventional perturbative approach cannot be applied for hadronic process at scales \lesssim 1 GeV since the strong coupling constant $\alpha_s \sim$ 1
 - \implies we cannot calculate the masses of mesons and baryons from QCD even if we are given α_s and the masses of quarks.
- ullet Bound state in QCD very different from QED e.g. the binding energy of a hydrogen atom is to a good approximation the sum of it constituent masses. Similarly for nuclei the binding energy is $\mathcal{O}(\text{MeV})$. For the proton almost all the mass is attributed to the strong non-linear interactions of the gluons.

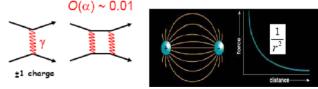




QCD versus QED

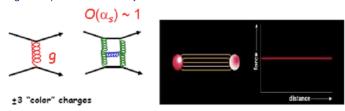
Quantum Electrodynamics (QED): The interaction is due to the exchange of photons. Every time there is an exchange of a photon there is a correction in the interaction of the order of 0.01.

ightarrow we can apply perturbation theory reaching whatever accuracy we like



QCD: Interaction due to exchange of gluons. In the energy range of $\,^{\sim}$ 1GeV the coupling constant is $^{\sim}1$

→ We can no longer use perturbation theory



QCD on the lattice

Why Lattice QCD?

- Discrete space-time lattice acts as a non-perturbative regularization scheme with the lattice spacing a providing an ultraviolet cutoff at π/a → no infinities. Furthermore, renormalized physical quantities have a finite well behaved limit as a → 0.
- Can be simulated on the computer using methods analogous to those used for Statistical Mechanics systems. These simulations allow us to calculate correlation functions of hadronic operators and matrix elements of any operator between hadronic states in terms of the fundamental quark and gluon degrees of freedom.

Like continuum QCD lattice QCD has as unknown input parameters the coupling constant α_s and the masses of the up, down, strange, charm and bottom quarks (the top quark is too short lived).

⇒ Lattice QCD provides a well-defined approach to calculate observables non-perturbative starting directly from the QCD Langragian.

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Walking down the street one night, someone found Nasruddin crawling on his hands and knees under a lamp post near his own front door..

Nasruddin, what's wrong? What are you doing here?

Looking for my house key. I can't get in until I find it..

Oh? Let me help you.

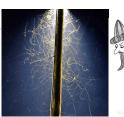
The friend joined Nasruddin in the search, and together they combed all the area under the lamp, but could find nothing.

That's strange, Nasruddin. It's just not here. Are you sure this is where you dropped it?

Here? No, of course not. I dropped the key over there in the shadows by the door.

But - Nasruddin, if you lost the key there, why are you looking here under the lamp?

Well, it's very simple. Over there it's so dark I can't see a thing, and here it's as bright as day!





Path Integrals-An overview

Path integrals in one-dimensional quantum mechanics

Time evolution of a quantum mechanical system:

$$|\psi(t_f)\rangle=e^{-iH(t_f-t_i)}\,|\psi(t_i)\rangle$$
 , where H is the Hamiltonian and \hbar is set to one.

Evolution of a quantum position eigenstate $|x_i\rangle$ from time t_i to time t_i :

$$\langle x_f, t_f | x_i, t_i \rangle = \langle x_f | e^{-iH(t_f - t_i)} | x_i \rangle$$

Break $t \equiv t_f - t_i$ into a large number N of time slices separated by $\Delta t = (t_f - t_i)/N$:

$$e^{-iHt} = e^{-iH\Delta t} \int dx_{N-1} \left| x_{N-1} \right\rangle \left\langle x_{N-1} \right| e^{-iH\Delta t} \int dx_{N-2} \left| x_{N-2} \right\rangle \left\langle x_{N-2} \right| \cdots e^{-iH\Delta t}$$

$$\begin{split} &\langle x_{k+1} | \, e^{-i\Delta t \left(\frac{\hat{p}^2}{2m} + V(\hat{x})\right)} \, | \, x_k \rangle \overset{\Delta t \to 0}{\sim} \, \int \frac{dp}{2\pi} \, \langle x_{k+1} | \, p \rangle \, e^{-i\Delta t} \frac{p^2}{2m} \, e^{-i\Delta t V(x_k)} \, \langle p \, | \, x_k \rangle \\ &\sim \int \frac{dp}{2\pi} \, e^{jp(x_{k+1} - x_k) - i\Delta t} \, \frac{p^2}{2m} - i\Delta t \, V(x_k) \, = \, \sqrt{\frac{2m\pi}{\Delta t}} \, e^{i\Delta t} \bigg[\frac{m}{2} \left(\frac{(x_{k+1} - x_k)}{\Delta t}\right)^2 - V(x_k) \bigg] \, + \, \mathcal{O}(\Delta t^2) \end{split}$$

$$\langle x_{f} | e^{-iH(t_{f}-t_{f})} | x_{i} \rangle = \int \mathcal{D}(x_{1}, \cdots, x_{N-1}) e^{i\Delta t \sum_{k} \left[\frac{m}{2} \left(\frac{x_{k+1}-x_{k}}{\Delta t} \right)^{2} - V(x_{k}) \right] }$$

$$\rightarrow \int_{x(0)=x_{f}}^{x(t)=x_{f}} \mathcal{D}[x(t)] e^{iS_{\text{classical}}[x(t)]}$$

i.e. the evolution operator is the sum over all paths weighted by the exponential of the classical action.

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$$\begin{split} &\langle x_{k+1} \,|\, e^{-i\Delta t \left(\frac{\hat{p}^2}{2m} + V(\hat{x})\right)} \,|\, x_k \rangle \overset{\Delta t \to 0}{\sim} \int \frac{dp}{2\pi} \,\, \langle x_{k+1} \,|\, p \rangle e^{-i\Delta t} \frac{p^2}{2m} e^{-i\Delta t V(x_k)} \,\langle p \,|\, x_k \rangle \\ &\sim \int \frac{dp}{2\pi} e^{ip(x_{k+1} - x_k) - i\Delta t} \, \frac{p^2}{2m} e^{-i\Delta t} \,V(x_k) = \sqrt{\frac{2m\pi}{\Delta t}} e^{i\Delta t} \left[\frac{m}{2} \left(\frac{(x_{k+1} - x_k)}{\Delta t}\right)^2 - V(x_k)\right] + \mathcal{O}(\Delta t^2) \end{split}$$

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$$\rightarrow \int_{x(0)=x_{i}}^{x(t)=x_{f}} \mathcal{D}[x(t)] e^{iS_{classical}[x(t)]}$$

i.e. the evolution operator is the sum over all paths weighted by the exponential of the classical action.

Many degrees of freedom:

$$\left\langle x_{f}^{1}, \cdots, x_{f}^{N} \middle| e^{-iH(t_{f}-t_{i})} \middle| x_{i}^{1}, \cdots, x_{i}^{N} \right\rangle \int_{x_{i}^{1}(0)=x_{f}^{1}, \cdots, x_{i}^{N}(0)=x_{f}^{N}}^{x_{i}^{1}(t)=x_{f}^{1}, \cdots, x_{i}^{N}(t)=x_{f}^{N}} \mathcal{D}[x^{1}(t), \cdots, x^{N}(t)] e^{iS_{\text{classical}}[x^{1}(t), \cdots, x^{N}(t)]}$$

Time-ordered product

$$\begin{split} \hat{T}\mathcal{O}(t_1)\mathcal{O}(t_2)e^{-iHt} &= e^{-iH(t_1-t_2)}\mathcal{O}(t_2)e^{-iH(t_2-t_1)}\mathcal{O}(t_1)e^{-iH(t_1-t_1)} \\ &\to \int \mathcal{D}[x(t)]\,\mathcal{O}(x(t_1))\,\mathcal{O}(x(t_2))\,e^{iS_{cl}[x(t)]} \end{split}$$

i.e. we automatically get the time-ordered product.

Classical limi

Include \hbar in the expressions and consider $\hbar \to 0$

$$\int \mathcal{D}[X(t)] e^{iS_{cl}[X(t)]/\hbar} \stackrel{SPA}{\Rightarrow} e^{iS_{cl}[X_{cl}(t)]/\hbar} \left(\det \left(m \frac{d^2}{dt^2} + V^{\prime\prime}(x) \right) \right)^{-1/2} + \mathcal{O}(\hbar)$$

i.e. the leading correction are quadratic fluctuations around the classical path.

Many degrees of freedom:

$$\left\langle x_{f}^{1}, \cdots, x_{f}^{N} \middle| e^{-iH(t_{f}-t_{i})} \middle| x_{i}^{1}, \cdots, x_{i}^{N} \right\rangle \int_{x_{i}^{1}(0)=x_{f}^{1}, \cdots, x^{N}(t)=x_{f}^{N}}^{x^{1}(t)=x_{f}^{1}, \cdots, x^{N}(t)=x_{f}^{N}} \mathcal{D}[x^{1}(t), \cdots, x^{N}(t)] e^{iS_{\text{classical}}[x^{1}(t), \cdots, x^{N}(t)]}$$

Time-ordered product

$$\begin{split} \hat{T}\mathcal{O}(t_1)\mathcal{O}(t_2)e^{-iHt} &= e^{-iH(t_1-t_2)}\mathcal{O}(t_2)e^{-iH(t_2-t_1)}\mathcal{O}(t_1)e^{-iH(t_1-t_1)} \\ &\to \int \mathcal{D}[x(t)]\,\mathcal{O}(x(t_1))\,\mathcal{O}(x(t_2))\,e^{iS_{cl}[x(t)]} \end{split}$$

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$$\int \mathcal{D}[x(t)] e^{iS_{cl}[x(t)]/\hbar} \stackrel{\text{SPA}}{\Rightarrow} e^{iS_{cl}[x_{cl}(t)]/\hbar} \left(\det \left(m \frac{d^2}{dt^2} + V''(x) \right) \right)^{-1/2} + \mathcal{O}(\hbar)$$

i.e. the leading correction are quadratic fluctuations around the classical path.

Rotation to Euclidean time

Under Wick rotation

$$x_0 \equiv t \rightarrow -ix_4 \equiv -i\tau$$
,
 $p_0 \equiv E \rightarrow ip_4$.

The Euclidean convention is

$$x_E^2 = \sum_{i=1}^4 x_i^2 = \vec{x}^2 - t^2 = -x_M^2,$$

$$p_E^2 = \sum_{i=1}^4 p_i^2 = \vec{p}^2 - E^2 = -p_M^2.$$

$$e^{-\tau H} \to \int \mathcal{D}[x_1, \cdots, x_{N-1}] e^{-\Delta \tau \sum_k \left[\frac{m}{2} \left(\frac{x_k - x_{k-1}}{\Delta \tau}\right)^2 + V(x_{k-1})\right]}$$

i.e. the Lagrangian is effectively replaced by the Hamiltonian in the exponent

- Euclidean path purely real
- Similar to statistical mechanics: $Z = Tre^{-\tau H} = \int dx \langle x|e^{-H\tau}|x\rangle = \int \mathcal{D}(x_1, \dots, x_N)e^{-S(x_1, \dots, x_N)}$

$$S[x] \equiv \int_{\tau_i}^{\tau_f} d\tau \, L(x, \dot{x}) \equiv \int_{\tau_i}^{\tau_f} dt \, \left[\frac{m \dot{x}(\tau)^2}{2} + V(x(\tau)) \right],$$

forms the basis of numerical simulations

Low lying spectrum

Setting

$$x_i = x_f \equiv x$$
 $t_f - t_i \equiv t = \tau$

the propagator is written as

$$\langle x| e^{-Ht} |x\rangle = \sum_{n} \langle x| \psi_{n} \rangle e^{-E_{n}t} \langle \psi_{n}| x \rangle$$

where $|\psi_n\rangle$ is the energy eigenstate with eigenvalue E_n . The sum is dominated by the lowest-energy states when T is large, because of the exponentials, and in the limit of very large T only the groundstate, $|\psi_0\rangle$, contributes:

$$\langle x | e^{-Ht} | x \rangle \stackrel{t \to \infty}{\longrightarrow} e^{-E_0 t} |\langle x | \psi_0 \rangle|^2$$
.

We extract the groundstate energy E_0 by integrating over x,

$$\int dx \langle x| e^{-Ht} |x\rangle \stackrel{t\to\infty}{\longrightarrow} e^{-E_0 t},$$

and then, going back to the previous equation, we determine the groundstate wave-function $\psi_{E_0}(x) \equiv \langle x | E_0 \rangle$.

Numerical evaluation

Develop a numerical procedure for evaluating the propagator using a path integral in its discrete version

$$t_i = t_i + j \Delta t$$
 for $j = 0, 1 \dots N \Delta t = t$

 $\to x = \{x(t_0), x(t_1) \dots x(t_N)\}$: refer to such a path as a "configuration". Integral over all paths becomes an ordinary integral over all possible values for each of the $x(t_j)$'s:

$$\int \mathcal{D}x(t) \to A \int_{-\infty}^{\infty} dx_1 dx_2 \dots dx_{N-1}, \quad x_j \equiv x(t_j), \quad A = \left(\frac{2\pi m}{\Delta t}\right)^{N/2}$$

We don't integrate over the endpoints since they are held fixed; for example, in the Tre^{-tH} the boundary conditions are: $x_0 = x_N = x$.

⇒ we have reduce Quantum Mechanics to a problem in numerical integration.

Comment: $\frac{(x_j - x_{j-1})}{\Delta t}$ can be arbitrarily large in our path integral; i.e. paths can be arbitrarily rough. While not so important for our one-dimensional problem, this becomes a crucial issue for four-dimensional field theories. It is dealt with using renormalization theory.

Evaluation of excited states

To analyze excited states using path integrals, we interrupt the propagation of the groundstate by introducing new operators at intermediate times. Consider, for example,

$$\langle x(t_2)x(t_1)\rangle \equiv \frac{1}{Z} {\rm Tr} \hat{x}(t_2) \hat{x}(t_1) \, e^{-H(t_f-t_f)} = \frac{\int \mathcal{D} x(t) \, x(t_2) x(t_1) \, e^{-S[x]}}{\int \mathcal{D} x(t) \, e^{-S[x]}} \, , \quad t_2 > t_1$$

with $x_i = x_f = x$.

The numerator on the right-hand side equals

$$\int dx \langle x| e^{-H(t_{\bar{1}}-t_{\bar{2}})} \hat{x} e^{-H(t_{\bar{2}}-t_{\bar{1}})} \hat{x} e^{-H(t_{\bar{1}}-t_{\bar{1}})} |x\rangle.$$

Setting $T = t_{\rm f} - t_{\rm i}$ and $t = t_{\rm 2} - t_{\rm 1}$ we can rewrite the full expression as

$$\langle \hat{x}(t_2)\hat{x}(t_1)\rangle = \frac{1}{Z}\sum_n \langle \psi_n|\, e^{Ht_2}\hat{x}e^{-Ht_2}\, e^{Ht_1}\hat{x}e^{-Ht_1}\, e^{-HT}\, |\psi_n\rangle = \frac{\sum_n e^{-E_nT}\, \langle \psi_n|\, \hat{x}\, e^{-(H-E_n)t}\, \hat{x}\, |\psi_n\rangle}{\sum e^{-E_nT}}.$$

If $T\gg t$ and large, then the groundstate $|\psi_0\rangle$ dominates:

$$G(t) \equiv \langle x(t_2)x(t_1)\rangle \rightarrow \sum_m \langle \psi_0 | \hat{x} | \psi_m \rangle \langle \psi_m | e^{-(H-E_0)t} \hat{x} | \psi_0 \rangle.$$

In our harmonic oscillator example, the state propagating between the two x's cannot be $|\psi_0\rangle$ since x switches the parity of the state. Thus if we now make t large (but still $\ll T$)

$$G(t) \stackrel{t \text{ large}}{\longrightarrow} |\langle \psi_0 | x | \psi_1 \rangle|^2 e^{-(E_1 - E_0)t}$$

where $|\psi_1\rangle$ is the first excited state. Consequently we can extract the first excitation energy from the large-t dependence of G(t).

$$\log\left(\frac{G(t)}{G(t+a)}\right)\to (E_1-E_0)a,$$

and then, going back to G(t), we can determine the quantum mechanical transition matrix element $\langle \psi_0 | x | \psi_1 \rangle$.

Monte Carlo averages

Here we focus on the zero-temperature limit of large T.

For an arbitrary functional $\Gamma[x]$ employ a Monte Carlo procedure:

$$\langle \Gamma[x] \rangle = \frac{\int \mathcal{D}x(t) \, \Gamma[x] \, \mathrm{e}^{-S[x]}}{\int \mathcal{D}x(t) \, \mathrm{e}^{-S[x]}},$$

is a weighted average over paths with weight $\exp(-S[x])$. We generate a large number, N_{cf} , of random paths or configurations,

$$x^{(\alpha)} \equiv \{x_0^{(\alpha)} x_1^{(\alpha)} \dots x_{N-1}^{(\alpha)}\} \qquad \alpha = 1, 2 \dots, N_{cf},$$

on our grid in such a way that the probability $P[x^{(\alpha)}]$ for obtaining any particular path $x^{(\alpha)}$ is

$$P[x^{(\alpha)}] \propto e^{-S[x^{(\alpha)}]}.$$

Then an unweighted average of $\Gamma[x]$ over this set of paths approximates the weighted average over uniformly distributed paths:

$$\langle \Gamma[x] \rangle \approx \overline{\Gamma} \equiv \frac{1}{N_{cf}} \sum_{\alpha=1}^{N_{cf}} \Gamma[x^{(\alpha)}].$$

 $\overline{\Gamma}$ is our "Monte Carlo estimator" for $\langle \Gamma[x] \rangle$ on our lattice.

The Monte Carlo uncertainty $\sigma_{\overline{\Gamma}}$ in our estimate is a potential source of error; it is estimated in the usual fashion:

$$\sigma_{\overline{\Gamma}}^2 \approx \frac{1}{N_{\rm cf}} \left\{ \frac{1}{N_{\rm cf}} \sum_{\alpha=1}^{N_{\rm cf}} \Gamma^2[x^{(\alpha)}] - \overline{\Gamma}^2 \right\}.$$

This becomes

$$\sigma_{\overline{\Gamma}}^2 = \frac{\langle \Gamma^2 \rangle - \langle \Gamma \rangle^2}{N_{cf}}$$
 for large N_{cf} .

The statistical uncertainties vanish as $1/\sqrt{N_{cf}}$ when N_{cf} increases.

Metropolis Algorithm

We need an algorithm to to create our set of random paths $x^{(\alpha)}$ with probability $\frac{e^{-S[x]}}{Z}$, where $Z = \int \mathcal{D}[x(t)]e^{-S[x]}$.

 \implies a simple procedure, though not always the best, is the Metropolis Algorithm:

- Start with an arbitrary path $x^{(0)}$
- Modify by visiting each of the sites on the lattice, and randomizing the x_i's at those sites, one at a time, in a particular fashion as described below → generate a new random path from the old one: x⁽⁰⁾ → x⁽¹⁾. This is called "updating" the path.
- Apply to $x^{(1)}$ to generate path $x^{(2)}$, and so on until we have N_{cf} random paths.

The algorithm for randomizing x_i at the j^{th} site is:

- Generate a random number $-\epsilon < \zeta \le \epsilon$, with uniform probability;
- Let $x_i \to x_i + \zeta$ and compute the change ΔS in the action;
- If $\Delta S < 0$ retain the new value for x_i , and proceed to the next site;
- If $\Delta S > 0$ accept change with probability $\exp(-\Delta S)$ i.e. generate a random number η uniformly distributed between 0 and 1; retain the new value for x_j if $\exp(-\Delta S) > \eta$, otherwise restore the old value; proceed to the next site.

Comments

- Choice of ϵ : should be tuned so that 40%–60% of the x_j 's are changed on each pass (or "sweep") through the lattice. Then ϵ is of order the typical quantum fluctuations expected in the theory. Whatever the ϵ , successive paths are going to be quite similar and so contain rather similar information about the theory. Thus when we accumulate random paths $x^{(\alpha)}$ for our Monte Carlo estimates we should keep onlevery N_{cor} -th path; the intervening sweeps erase correlations, giving us configurations that are statistically independent. The optimal value for N_{cor} depends upon the theory, and can be found by experimentation. It also depends on the lattice spacing a.
- Initial configuration: Guess the first configuration \rightarrow discard some number of configurations at the beginning, before starting to collect $x^{(\alpha)}$'s. This is called "thermalizing the lattice."

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 \Longrightarrow a simple procedure, though not always the best, is the Metropolis Algorithm:

- Start with an arbitrary path $x^{(0)}$
- Modify by visiting each of the sites on the lattice, and randomizing the x_i's at those sites, one at a time, in a particular fashion as described below → generate a new random path from the old one: x⁽⁰⁾ → x⁽¹⁾. This is called "updating" the path.
- Apply to $x^{(1)}$ to generate path $x^{(2)}$, and so on until we have N_{cf} random paths.

The algorithm for randomizing x_i at the j^{th} site is:

- Generate a random number $-\epsilon < \zeta \le \epsilon$, with uniform probability;
- Let $x_i \to x_i + \zeta$ and compute the change ΔS in the action;
- If $\Delta S < 0$ retain the new value for x_i , and proceed to the next site;
- If $\Delta S > 0$ accept change with probability $\exp(-\Delta S)$ i.e. generate a random number η uniformly distributed between 0 and 1; retain the new value for x_j if $\exp(-\Delta S) > \eta$, otherwise restore the old value; proceed to the next site.

Comments:

- Choice of ϵ : should be tuned so that 40%–60% of the x_j 's are changed on each pass (or "sweep") through the lattice. Then ϵ is of order the typical quantum fluctuations expected in the theory. Whatever the ϵ , successive paths are going to be quite similar and so contain rather similar information about the theory. Thus when we accumulate random paths $x^{(\alpha)}$ for our Monte Carlo estimates we should keep only every N_{cor} -th path; the intervening sweeps erase correlations, giving us configurations that are statistically independent. The optimal value for N_{cor} depends upon the theory, and can be found by experimentation. It also depends on the lattice spacing a.
- Initial configuration: Guess the first configuration → discard some number of configurations at the beginning, before starting to collect x^(α)'s. This is called "thermalizing the lattice."

Statistical errors

An important part of any Monte Carlo analysis is the estimate of the statistical errors.

The "statistical bootstrap," method:

The bootstrap procedure provides new, almost *zero-cost* random ensembles of measurements by synthesizing them from the original ensemble of N_{cf} measurements.

Consider an ensemble $\{G^{(\alpha)}, \alpha = 1 \dots N_{cf}\}$ of Monte Carlo measurements

- ▶ Construct a "bootstrap copy" of that ensemble by selecting $G^{(\alpha)}$'s at random from the original ensemble, taking $N_{\rm cf}$ in all while allowing duplications and omissions
 - ightarrow resulting ensemble of G's may have two or three copies of some $G^{(\alpha)}$'s, and no copies of others
- Use new ensemble to obtain a new estimate of some the quantity of interest.
- Repeat this procedure to generated as many bootstrap copies of the original ensemble as one wishes, and from each we can generate a new estimate.

The distribution of these estimates approximates the distribution of the quantity that would have been obtained from the original Monte Carlo, and so can be used to estimate the statistical error in our original estimate.

- The "Jackknife" method: Similar to boostrap but remove a set of measurements at a time from the sample set. In general easier to use than boostrap.
- The "binning" method:

At the end of a simulation we have set of configurations $x^{(\alpha)}$, and for each a set of measurements like $G^{(\alpha)}$, our propagator. We partially average or bin the measurements: For example, instead of storing each of

$$G^{(1)}$$
 $G^{(2)}$ $G^{(3)}$ $G^{(4)}$ $G^{(5)}$.

we might instead store

$$\overline{G}^{(1)} \equiv \frac{G^{(1)} + G^{(2)} + G^{(3)} + G^{(4)}}{4}$$
 $\overline{G}^{(2)} \equiv \frac{G^{(5)} + G^{(6)} + G^{(7)} + G^{(8)}}{4}$

Binning reduces or can even remove the effects of correlations between different configurations

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Binning reduces or can even remove the effects of correlations between different configurations.

Scalar field theory

Let the continuum \vec{r} be defined on lattice points i.e. $\vec{r} \to \vec{n} \equiv (n_1, n_2, n_3)a$ where a is the lattice spacing. \Longrightarrow equivalent to many-body problem where:

$$\hat{x}_i, \hat{p}_i \rightarrow \hat{\phi}(\vec{n}), \hat{\pi}(\vec{n})$$
 $\hat{x}_i | x \rangle = x_i | x \rangle \rightarrow \hat{\phi}(\vec{n}) = \phi(\vec{n}) | \phi \rangle$.

We then have

$$\int d^3r \left\{ \frac{1}{2} \pi^2(\vec{r}) + \frac{1}{2} |\vec{\nabla}\phi(\vec{r})|^2 + V(\phi(\vec{r})) \right\} \rightarrow \sum_{\vec{n}} a^3 \left\{ \frac{1}{2} \pi^2(\vec{n}) + \frac{1}{2a^2} \sum_{i=1}^3 |\phi(\vec{n} + a\mu_i) - \phi(\vec{n})|^2 + V(\phi(\vec{n})) \right\}$$

where μ_i denotes a displacement by one lattice site in the i^{th} direction.

The evolution operator in Euclidean time:

$$e^{-t\sum_{\vec{n}}a^3\left\{\frac{1}{2}\pi^2(\vec{n})+F(\phi(\vec{n}))\right\}}=\int\mathcal{D}[\phi(\vec{n})]e^{-\Delta ta^3\sum_{\vec{n},k}\left[\frac{1}{2\Delta t^2}\left(\phi_{k+1}(\vec{n})-\phi_k(\vec{n})\right)^2+F(\phi_k(\vec{n}))\right]}$$

Take isotropic lattice i.e. $\Delta t = a$

 \implies time slicing replaces $\hat{\pi}^2(\vec{n})$ by $\frac{1}{\Delta t}(\phi_{k+1}(\vec{n}) - \phi_k(\vec{n}))^2 \equiv \frac{1}{\Delta t}(\phi(n+a\mu_0) - \phi(n))^2$ which has the same structure as the discrete spatial derivative and where $n=(n_0,n_1,n_2,n_3)a$.

$$\Longrightarrow \mathcal{O}(\phi) e^{-t \int \sigma^3 r \left\{ \frac{1}{2} \, \pi^2(\vec{r}) + \frac{1}{2} \, |\vec{\nabla} \, \phi(\vec{r})|^2 + V(\phi(\vec{r})) \right\}} \to \int \mathcal{D}[\phi(n)] \, O(\phi) e^{-S_{cl}[\phi]}$$

where $S_{cl}[\phi] = \sum_{n} a^4 \left\{ \sum_{i=0}^{3} \frac{\left(\phi(n+a\mu_i) - \phi(n)\right)^2}{a^2} + V(\phi(n)) \right\}.$

Note that S_{cl} is completely symmetric in time and space \rightarrow if we choose periodic b.c. then the shortest dimension acts as a finite temperature.

Coherent States for bosons

Goal: generalize the scalar field example for second quantized boson and fermion fields. For the Feynman path integral of our 1-d QM example we needed:

- Eigenstates of \hat{x} , $\hat{x} | x \rangle = x | x \rangle$ and
- Unity: $\int dx |x\rangle \langle x| = 1$

The analogs for creation and annihilation operators are provided by boson coherent states.

Consider a creation operator a^{\dagger} then

$$\left[\hat{a},\hat{a}^{\dagger}\right]=1 \qquad \hat{a}^{\dagger}\left|n\right\rangle = \sqrt{n+1}\left|n+1\right\rangle \qquad \hat{a}\left|n\right\rangle = \sqrt{n}\left|n-1\right\rangle \qquad \left|n\right\rangle = \frac{1}{\sqrt{n!}}\left(\hat{a}^{\dagger}\right)^{n}\left|0\right\rangle$$

Define the coherent state $|z\rangle$ by

$$|z\rangle \equiv e^{za^{\dagger}} |0\rangle = \sum_{n} \frac{z^{n}}{n!} (\hat{a}^{\dagger})^{n} |0\rangle = \sum_{n} \frac{z^{n}}{\sqrt{n!}} |n\rangle$$

Properties:

$$\hat{a}|z\rangle = \sum_{n} \frac{z^{n}}{\sqrt{n!}} \hat{a}|n\rangle = z \sum_{n} \frac{z^{n-1}}{\sqrt{(n-1)!}} |n-1\rangle = z|z\rangle$$

$$\langle z|z'\rangle = \sum_{mn} \langle m| \frac{z^{*m}}{\sqrt{m!}} \frac{z'^{n}}{\sqrt{n!}} |n\rangle = e^{z^{*}z'}$$

$$\langle z| : A(\hat{a}^{\dagger}\hat{a}) : |z\rangle' = e^{z^{*}z'} A(z^{*}, z')$$

$$1 = \int \frac{dzdz^{*}}{2\pi i} e^{-z^{*}z'} |z\rangle \langle z|$$

$$(1)$$

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Generalize for a set of creation operators $\hat{a}^{\dagger}_{\alpha}$

$$\begin{array}{rcl} |z\rangle & = & e^{\sum_{\alpha}z_{\alpha}\hat{a}^{\dagger}_{\alpha}}|0\rangle \\ \hat{a_{\alpha}}|z\rangle & = & z_{\alpha}|z\rangle \\ \langle z|:A(\hat{a}^{\dagger}\hat{a}):|z\rangle' & = & e^{\sum_{\alpha}z_{\alpha}Z'_{\alpha}}A(z^{*},z') \\ 1 & = & \int\prod_{\alpha}\frac{dz_{\alpha}z^{*}_{\alpha}}{2\pi i}e^{-z^{*}_{\alpha}z'_{\alpha}}|z\rangle\langle z|\equiv\int d\mu(z)|z\rangle\langle z| \end{array}$$

Path integral using coherent states

Time slicing the evolution operator:

$$\left\langle z_{f}\right|e^{-tH}\left|z_{i}\right\rangle =\left\langle z_{f}\right|e^{-\Delta tH}\int d\mu(z_{N-1})\left|z_{N-1}\right\rangle \left\langle z_{N-1}\right|e^{-\Delta tH}\int d\mu(z_{N-2})\cdots e^{-\Delta tH}\left|z_{i}\right\rangle$$

The matrix element of the infinitesimal evolution operator is

$$d\mu(z_k) \langle z_k | e^{-tH} | z_{k-1} \rangle = \prod_{\alpha} \frac{dz_{k,\alpha}^* dz_{k,\alpha}}{2i\pi} e^{-\sum_{\alpha} z_{k,\alpha}^* (z_{k,\alpha} - z_{k-1,\alpha}) - \Delta t H(z_{k,\alpha}^*, z_{k-1,\alpha})}$$

resulting in

$$\begin{split} \langle z_f | \, e^{-tH} \, | z_i \rangle &= \int \mathcal{D}[z_{k,\alpha}^*, z_{k,\alpha}] e^{-S(z_{k,\alpha}^*, z_{k,\alpha})} \\ S(z^*, z) &= \sum_k \Delta t \left\{ \sum_\alpha z_{k,\alpha}^* \left(\frac{z_{k,\alpha} - z_{k-1,\alpha}}{\Delta t} \right) + H(z_{k,\alpha}^*, z_{k-1,\alpha}) \right\} \end{split}$$

Coherent states for fermions

Fermions are represented by anti-commuting creation and annihilation operators c^\dagger_α and $c_\alpha \to$ need to introduce anti-commuting Grassmann variables ξ such that

$$\hat{c}_{lpha}\ket{\xi}=\xi_{lpha}\ket{\xi} \qquad \hat{c}_{lpha}\hat{c}_{eta}\ket{\xi}=\xi_{lpha}\xi_{eta}\ket{\xi}=-\xi_{eta}\xi_{lpha}\ket{\xi}=-\hat{c}_{eta}\hat{c}_{lpha}\ket{\xi}$$

Since $\xi_{\alpha}^2=0$ (Pauli principle) the only functions allowed are monomials. The rules for integration over a Grassmann variable ξ and ξ^* are

$$\int d\xi_{\alpha} = \int d\xi_{\alpha}^* = 0, \int d\xi_{\alpha} \xi_{\alpha} = \int d\xi_{\alpha}^* \xi_{\alpha}^* = 1$$

A fermion coherent state is defined by

$$|\xi\rangle \equiv e^{-\sum_{\alpha} \xi_{\alpha} c_{\alpha}^{\dagger}} |0\rangle$$

with similar properties to bosons. The path integral have similar form to that for bosons with some minus signs that distinguish between bosons and fermions.

Integration over fermions

For numerical evaluation we can not have the path integrals in terms of Grassmann variables. Fortunately for normalizable field theories we can integrate analytically over the fermionic degrees of freedom Recall Gaussian integral

$$\int \prod_i \frac{dz_i^* \, dz_i}{2i\pi} \, \mathrm{e}^{-z_i^* \, H_{ij} z_j + J_i^* \, z_i + z_i^* \, J_i} = [\det\! H]^{-1} \, \mathrm{e}^{J_i^* \, H_{ij}^{-1} \, J_j}$$

An analogous result is obtained for Grassmann "Gaussian": For one pair of Grassmann variables we have

$$\int d\xi^* d\xi \, e^{-\xi^* a\xi} = \int d\xi^* d\xi (1 - \xi^* a\xi) = a$$

This generalizes to

$$\int \prod_i \, d\xi_i^* \, d\xi_i \, e^{-\xi_i^* H_{ij} \xi_j + \eta_i^* \xi_i + \xi_i^* \, \eta_i} = [\textit{detH}] \, e^{\eta_i^* H_{ij}^{-1} \, \eta_j}$$

i.e. the only difference is that detH appears in the numerator \to accounts for the minus sign of fermion loops. If our action is of the form $S(\xi^*, \xi, \phi) = \xi_i^* M(\phi)_{ij} \xi_j + S_B(\phi)$ then

$$\int d\xi^* d\xi d\phi \, e^{\xi_i^* M(\phi)_{ij} \xi_j + S_B(\phi)} = \int d\phi \, det M(\phi) \, e^{S_B(\phi)}$$

i.e. $S_{\text{eff}}(\phi) = \ln \text{detM}(\phi) + S_{\mathcal{B}}(\phi)$

Fermion propagators

Consider the time ordered product of field creation and annihilation operators at space-time points $j = (x_j, t_j)$ and $i = (x_i, t_i)$ respectively:

$$\langle \hat{T}\psi_i\bar{\psi}_j\rangle = \text{Tr}\hat{T}\psi_i\bar{\psi}_j\,e^{-\bar{\psi}\textit{M}(\phi)\psi + \textit{S}_{\textit{B}}(\phi)} = \int \mathcal{D}[\phi]\mathcal{D}[\bar{\xi}\xi]\xi_i\bar{\xi}_j\,e^{-\bar{\xi}\textit{M}(\phi)\xi + \textit{S}_{\textit{B}}(\phi)} = \int \mathcal{D}[\phi]\textit{M}_{ij}^{-1}(\phi)\,e^{\textit{S}_{eff}(\phi)}$$

In general for n pairs of creation and annihilation operators

$$\begin{split} &\int \mathcal{D}(\xi^*,\xi)\,\xi_{i_1}\cdots\xi_{i_n}\xi_{j_n}^*\cdots\xi_{j_1}\;e^{-\xi^*M\xi}\\ = &\;\;\frac{\delta^{2n}}{\delta\eta_{i_1}^*\cdots\delta\eta_{i_n}^*\delta\eta_{j_n}\cdots\delta\eta_{j_1}}\;\int \mathcal{D}(\xi^*,\xi)\,e^{-\xi_i^*M_{ij}\xi_j+\eta_i^*\xi_i+\xi_i^*\eta_i\,\big|\eta=\eta^*=0}\\ = &\;\;\frac{\delta^{2n}}{\delta\eta_{i_1}^*\cdots\delta\eta_{i_n}^*\delta_{j_n}\cdots\delta\eta_{j_1}}\;detM\;e^{\eta_i^*M_{ij}^{-1}\eta_j\,\big|\eta=\eta^*=0}\\ = &\;\;\sum_P(-1)^PM_{i_P_nj_n}^{-1}\cdots M_{i_{P_1}j_1}^{-1}\;e^{\ln detM} \end{split}$$

where *P* denotes a permutation of the indices. This is nothing else but Wick's theorem.

⇒ fermions can be integrated out and we are left only with an effective action with the bosonic degrees of freedom.

Boundary conditions:

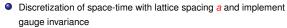
$$\textit{Tr} e^{-tH} = \int dz_0^* \, dz_0 \, e^{-z_0^* \, z_0} \, \left\langle \pm z_0 \right| \, e^{-tH} \, |z_0\rangle = \int dz_0^* \, dz_0 \, e^{-z_0^* \, z_0} \, \int d\mathcal{D}[z^*,z] \, e^{-S(z^*,z)}$$

where the plus is for bosons and minus for fermions and

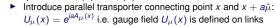
$$S(z^*,z) = \pm z_0^* (\pm z_0 - z_{N-1}) + H_{0,N-1} + z_{N-1}^* (z_{N-1} - z_{N-2}) + H_{N-1,N-2} + \cdots + z_1^* (z_1 - z_0) H_{1,0}.$$

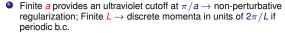
QCD on the lattice

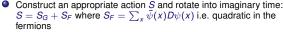
Lattice QCD: K. Wilson, 1974 provided the formulation; M. Creutz, 1980 performed the first numerical simulation











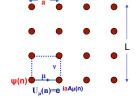
---- can be integrated out

Path integral over gauge fields:

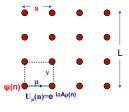
 $Z \sim \int \mathcal{D}U_{\mu}(x) \prod_{f} det(D_{f}[U]) e^{-S_{G}[U]}$

o Monte Carlo simulation to produce a representative ensemble of $\{U_{\mu}(x)\}$ using the largest supercomputers o

Observables: $\langle \mathcal{O} \rangle = \sum_{\{U_{\mu}\}} O(D^{-1}, U_{\mu})$

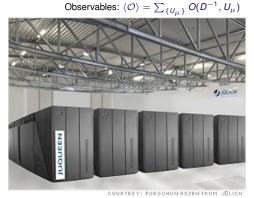


QCD on the lattice



Lattice QCD: K. Wilson, 1974 provided the formulation; M. Creutz, 1980 performed the first numerical simulation

- Discretization of space-time with lattice spacing a and implement gauge invariance
- Finite a provides an ultraviolet cutoff at π/a → non-perturbative regularization; Finite L → discrete momenta in units of 2π/L if periodic b.c.
- Path integral over gauge fields: $Z \sim \int \mathcal{D}U_{\mu}(x) \prod_{I} \det(D_{I}[U]) e^{-S_{G}[U]}$ → Monte Carlo simulation to produce a representative ensemble of $\{U_{\mu}(x)\}$ using the largest supercomputers → Observables: $\{U_{\mu}(x)\}$ = $\sum_{I} \mathcal{D}(D_{I}^{-1}, U_{I})$



Hadron mass

First goal: reproduce the low-lying masses

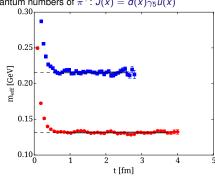
Use Euclidean correlation functions:

$$\begin{split} G(\vec{q},t_s) &= \sum_{\vec{x}_S} e^{-i\vec{x}_S \cdot \vec{q}} \langle J(\vec{x}_s,t_s) J^{\dagger}(0) \rangle \\ &= \sum_{n=0,\cdots,\infty} A_n e^{-E_n(\vec{q})t_S} \stackrel{t_S \to \infty}{\longrightarrow} A_0 e^{-E_0(\vec{q})t_S} \end{split}$$



Interpolating field with the quantum numbers of π^+ : $J(x) = \bar{d}(x)\gamma_5 u(x)$

- Large Euclidean time evolution gives ground state for given quantum numbers ⇒ enables determination of low-lying hadron properties



 $N_f=2+1+1$ TM fermions at $m_\pi=210$ MeV $N_f=2$ TM plus clover fermions at physical pion mass

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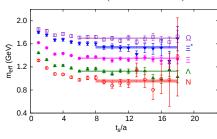
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Interpolating field with the quantum numbers of p: $J(x) = \epsilon^{abc} \left(u^{a \top}(x) C \gamma_5 d^b(x) \right) u^c(x)$

- Large Euclidean time evolution gives ground state for given quantum numbers ⇒ enables determination of low-lying hadron properties
- $\mathbf{0} \quad aE_{\text{eff}}(\vec{q}, t_s) = \ln \left[G(\vec{q}, t_s) / G(\vec{q}, t_s + a) \right]$ $= aE_0(\vec{q}) + \text{excited states}$ $\rightarrow aE_0(\vec{q}) \stackrel{\vec{q}=0}{\rightarrow} am$

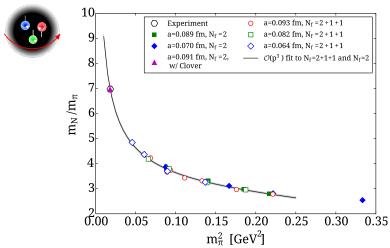


 $N_f=2$ TM plus clover fermions at physical pion mass Noise to signal increases with $t_s:\sim e^{(m_h-\frac{3}{2}m_\pi)t_s}$

Simulations with physical quark masses

A number of collaborations are producing simulations with physical values of the quark mass

European Twisted Mass Collaboration (ETMC): The nucleon

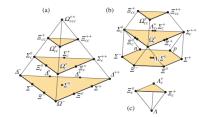


 $L\sim 3$ fm and $a\sim 0.1$ fm; $r_0\sim 0.5$ fm

SU(4) representations:

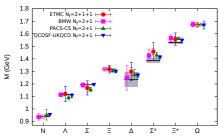
$$4 \otimes 4 \otimes 4 = 20 \oplus 20 \oplus 20 \oplus \overline{4}$$

$$\square \otimes \square \otimes \square = \square \square \oplus \square \oplus \square \oplus \square \oplus \square$$



Spectrum using $N_f = 2 + 1 + 1$ for a range of pion masses from about 450 MeV to 210 MeV

- Continuum extrapolation using three lattice spacings, a = 0.094 fm, 0.082 fm and 0.065 fm
- Volume dependence no observable effects within our statistics
- Chiral extrapolation biggest systematic error

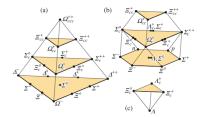


C.A., V. Drach, K. Jansen, Ch. Kallidonis, G. Koutsou, arXiv:1406.4310

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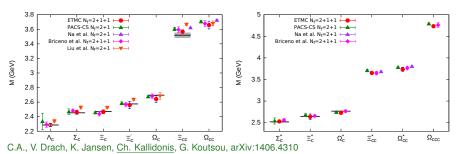
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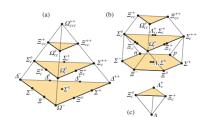
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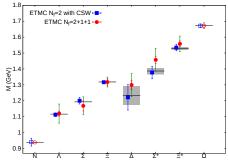


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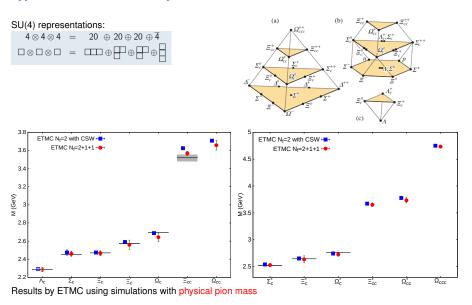
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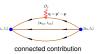
Results by ETMC using simulations with physical pion mass



Nucleon form factors

Evaluation of three-point functions:

$$G^{\mu\nu}(\Gamma, \vec{q}, t_s, t_{ins}) = \sum_{\vec{x}_s, \vec{x}_{ins}} e^{i\vec{x}_{ins} \cdot \vec{q}} \Gamma_{\beta\alpha} \langle J_{\alpha}(\vec{x}_s, t_s) \mathcal{O}^{\mu\nu}(\vec{x}_{ins}, t_{ins}) \overline{J}_{\beta}(\vec{x}_0, t_0) \rangle^{-\alpha}$$



Form ratio by dividing the three-point correlator by an appropriate combination of two-point functions:

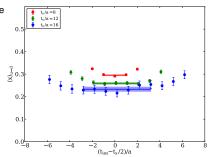
$$R(t_{s}, t_{\text{ins}}, t_{0}) \xrightarrow[(t_{s} - t_{\text{inc}})\Delta \gg 1]{(t_{s} - t_{\text{inc}})\Delta \gg 1}} \mathcal{M}[1 + \dots e^{-\Delta(\mathbf{p})(t_{\text{ins}} - t_{0})} + \dots e^{-\Delta(\mathbf{p}')(t_{s} - t_{\text{ins}})}]$$

- M the desired matrix element
- t_s, t_{ins}, t₀ the sink, insertion and source time-slice
- lacktriangle $\Delta(\mathbf{p})$ the energy gap with the first excited state

Connect lattice results to measurements:

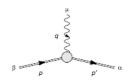
$$\mathcal{O}_{\overline{\mathrm{MS}}}(\mu) = Z(\mu, a)\mathcal{O}_{\mathrm{latt}}(a)$$

 \implies evaluate $Z(\mu, a)$ non-perturbatively



Electromagnetic form factors

$$\langle \textit{N}(\textit{p}',\textit{s}')|\textit{j}^{\mu}(0)|\textit{N}(\textit{p},\textit{s})\rangle = \bar{\textit{u}}_\textit{N}(\textit{p}',\textit{s}')\left[\gamma^{\mu}\textit{F}_{1}(\textit{q}^{2}) + \frac{\textit{i}\sigma^{\mu\nu}\textit{q}_{\nu}}{2m}\textit{F}_{2}(\textit{q}^{2})\right]\textit{u}_\textit{N}(\textit{p},\textit{s})$$





- lack Proton radius extracted from muonic hydrogen is 7.7 σ different from the one extracted from electron scattering, R. Pohl et al., Nature 466 (2010) 213
- Muonic measurement is ten times more accurate

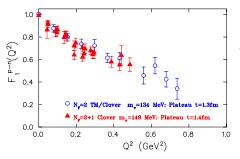
Electromagnetic form factors

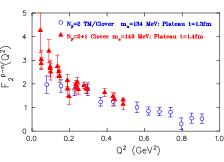
$$\langle N(p',s')|j^{\mu}(0)|N(p,s)\rangle = \bar{u}_N(p',s')\left[\gamma^{\mu}F_1(q^2) + \frac{i\sigma^{\mu\nu}q_{\nu}}{2m}F_2(q^2)\right]u_N(q^2)$$
The good news

The good news

Two studies at near physical pion mass:

- ETMC: $N_f = 2$ twisted mass with clover, a = 0.091 fm, $m_{\pi} = 134$ MeV, 1020 statistics
- MIT: $N_f = 2 + 1$ clover produced by the BMW collaboration, a = 0.116 MeV, $m_{\pi} = 149$ MeV, \sim 7750 statistics, J.M. Green et al. 1404,4029



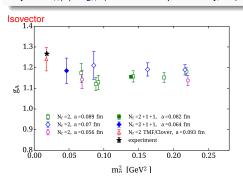


Agreement even before taking the continuum limit

Axial charge g_A

The good news:

Axial-vector FFs:
$$A_{\mu}^3 = \bar{\psi} \gamma_{\mu} \gamma_5 \frac{\tau^3}{2} \psi(x) \Longrightarrow \frac{1}{2} \bar{u}_N(\vec{p}') \left[\gamma_{\mu} \gamma_5 \frac{G_0(q^2)}{2m} \frac{q^{\mu} \gamma_5}{2m} \frac{G_p(q^2)}{2} \right] u_N(\vec{p})|_{q^2=0}$$
 \rightarrow yields $G_A(0) \equiv g_A$: i) well known experimentally, & ii) no quark loop contributions



- q_A at the physical point mass indicates agreement with the physical value → important to reduce error
- many results from other collaborations, e.g.
 - N_f = 2 + 1 Clover, J. R. Green et al., arXiv:1209.1687
 - N_f = 2 Clover, R.Hosley et al., arXiv:1302.2233
 - N_f = 2 Clover, S. Capitani et al. arXiv:1205.0180
 - N_f = 2 + 1 Clover, B. J. Owen et al., arXiv:1212.4668
 - N_f = 2 + 1 + 1 Mixed action (HISQ/Clover), T. Bhattacharya et
 - al., arXiv:1306.5435

Conclusions

- Lattice QCD is a successful framework for non-perturbative calculations → simulations of the full theory
 at the physical point are now possible
- After getting agreement with experiment it enables predictions of new quantities
- New ideas and techniques are needed for the study of e.g. resonances, multi-quark states, decays, etc
- Can be used to study models beyond the SM

As simulations at the physical pion mass and more computer resources are becoming available we expect many physical results on key hadron observables that will impact both experiments and phenomenology



New ideas and approaches are needed and welcome

Please join and contribute

Acknowledgments

European Twisted Mass Collaboration (ETMC)





Cyprus (Univ. of Cyprus, Cyprus Inst.), France (Orsay, Grenoble), Germany (Berlin/Zeuthen, Bonn, Frankfurt, Hamburg, Münster), Italy (Rome I, II, III, Trento), Netherlands (Groningen), Poland (Poznan), Spain (Valencia), Switzerland (Bern), UK (Liverpool)