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Lattice baryons in the $1/N$ expansion

Thomas DeGrand

Department of Physics, University of Colorado, Boulder, CO 80309, USA

Abstract

 $$\rm{F}^{\rm{2}}$$ respectively. The excitation levels of a rigid rotor. ons use the quenched approximation. Lattice spacings are matched using the static potential.
 \sqrt{N} , The spectra show independence on N and vacuum-to-hadron matrix elements scale as \sqrt{N} . The
 \sqrt{N} in spectrum s Results are presented for hadron spectroscopy with gauge groups $SU(N)$ with $N = 3, 5, 7$. Calculations use the quenched approximation. Lattice spacings are matched using the static potential. Meson spectra show independence on N and vacuum-to-hadron matrix elements scale as \sqrt{N} . The baryon spectrum shows the excitation levels of a rigid rotor.

I. INTRODUCTION

Replacing the "3" of color $SU(3)$ by "N"" and then taking N to infinity has a long history in the (continuum) phenomenology of the strong interactions, dating back to Refs. [1, 2].

There is also a literature of lattice simulations applied to gauge theories with the group $SU(N)$, for moderately large N. Most of it [3-10] is directed at the properties of pure gauge theory. I know of two papers on meson spectroscopy: Refs. [11, 12]. They are done at smaller values of "large N ," $N = 2, 3, 4, 6$. They are pretty standard lattice QCD spectroscopy calculations and reveal the N-independence of meson masses. Narayanan, Neuberger, and collaborators [13, 14] have measured masses and the pseudoscalar decay constant from simulations at much larger N (exploiting reduction, to simulate on smaller lattice sizes).

But only Ref. [15] discusses large N expectations for baryons, and it only makes comparisons to actual lattice data for $N = 3$. So it seemed like an appropriate time to look at a lattice calculation of baryon spectroscopy at several values of N.

Baryons in large N seem to be fascinating objects, either viewed as many-quark states [16] or as topological objects in effective theories of mesons[17, 18]. There is an enormous (continuum) literature about spectroscopy and matrix elements for large N baryons. Assorted early references include [19–23], summarized in a review, Ref. [24]. Perhaps a lattice study might reveal something interesting?

In fact, it does: general arguments [18, 19] state that the mass of an N color baryon of angular momentum J should show a rotor spectrum:

$$
M(N,J) = NA + \frac{J(J+1)}{N}B.
$$
\n⁽¹⁾

The formula applies to baryons made of an $SU(2)$ isospin doublet of equal mass quarks. The parameters A and B depend on the quark mass and both should be some "typical hadron size," a few hundred MeV. The observation of Eq. 1 in simulation data is the main new result of this paper. Common expectations are that Eq. 1 is only true for small values of J, because then the terms have meaning as a good expansion in $1/N$. However, I observe that it holds both for the bottom and the top of the multiplets where I did measurements.

The first response of the lattice simulator to a proposal to look at baryons at large N is undoubtedly negative: The cost of an N-color simulation scales roughly like N^3 , just from

the cost of multiplying $SU(N)$ matrices. Baryons are large objects so big volumes will be needed. Ordinary $(N = 3)$ baryon correlators are noisy and because baryon masses scale as N, baryons in higher N are probably noisier. And most of the literature involves some combination of small $1/N$ effects and expectation values of operators $(\xi \sim \langle B|O|B\rangle)$, which are already hard enough for $N = 3$ [25]. Fortunately, large N tests seem (mostly) not to require extrapolations to zero quark mass.

Only for odd N are baryons fermions, of course, so I am performing simulations of $SU(N)$ gauge theories with $N = 3, 5, 7$.

I used the quenched approximation for all N. The quenched approximation contains many of the same ingredients as the $N \to \infty$ limit of QCD: hadrons are bound states of the minimum number of fermions, and decay widths and other effects effects of virtual $\bar{q}q$ loops die away as inverse powers of N . At nonzero N all these effects are present. That means that, truly, I am not simulating QCD for any value of N. However, a first study does not quite demand the same level of quality as a later one might. I am interested in comparing simple observables in systems which differ only in N , but otherwise have the same UV cutoff (lattice spacing), lattice action, and physical volume. This is easy to set up in quenched approximation. Quenched simulations for $N = 3$ are obsolete, but the effects of dynamical fermions on spectroscopy are really not all that large: away from the deep chiral limit they are small, order ten per cent for simple observables (compare Fig. 1 in Ref. [26]).

A practical difficulty with using the quenched approximation comes when one wishes to convert a lattice number to MeV. The spectrum of quenched QCD is simply different from the spectrum of "real" QCD. I will compare dimensionless ratios of lattice quantities in my tests of scaling with N. Any conversions to MeV I make are only qualitative statements.

One could take the position that one should test everything about large- N QCD at once, by doing simulations with dynamical fermions at several values of N. This just postpones the matching problem: the spectrum of the different N QCD's will be different. But perhaps that is not the right point of view. The real question one asks when comparing several values of N is whether dimensionless ratios of masses show some smooth behavior with N . And for that, one can begin simply, make observations, and then ask how they change as one does ever more realistic simulations.

For all N's, I use lattices of size $16³ \times 32$ points. I match the bare couplings so that the lattice spacings, as measured by various observables from the heavy quark potential, are the same, and then sweep in quark mass over similar ranges. This insures that finite volume and nonzero lattice spacing artifacts will be similar across the board. I set the common scale using the shorter version of the Sommer $[27]$ parameter, r_1 , defined in terms of the force $F(r)$ so that $r^2F(r) = -1.0$ at $r = r_1$. The real-world value is $r_1 = 0.31$ fm [28], and with it my lattice spacings would be about 0.08 fm.

Before going on, it's useful to set some definitions. The 't Hooft coupling is

$$
g^2 N = \lambda \tag{2}
$$

and the usual gauge coupling is thus

$$
\beta = \frac{2N}{g^2} = \frac{2N^2}{\lambda}.\tag{3}
$$

The combination g^2C_F , which appears in all perturbative calculations of renormalization factors, is equal to $\lambda(1-1/N^2)/2$. For comparisons at fixed 't Hooft coupling, like the ones I present, this suggests that differences which might be perturbative (lattice–to–continuum matching factors, for example) will scale up to corrections of order $1/N^2$.

The outline of the paper is as follows: The next section describes some technical problems I faced: using fat links, gauge fixing, and the construction of baryon operators. Only the third topic might be of interest to continuum physicists. Next, I present lattice results for the potential and for meson masses and simple matrix elements. The potential measurements show the extent to which lattice spacings are matched. The mesonic observables illustrate the N – independence of masses and the \sqrt{N} scaling of quantities like the pseudoscalar decay constant. Finally, in Section IV, I show some results for baryon spectra. The major one is the presence of a rotor spectrum of excitations for two-flavor baryons.

II. TECHNICAL DETAILS

A. Simulation techniques

My simulations use a version of the publicly available package of the MILC collaboration [29], modified to generate gauge configurations and quark propagators at arbitrary N. Prior to this project, it had been extensively used in studies of $N = 2, 3, 4$ [30–32].

The gauge action is the usual Wilson action. Quenched simulations are performed using the standard mix of Brown - Woch microcanonical over-relaxation [33] and Cabibbo - Marinari heat bath updates [34], performed on all $N(N-1)/2$ $SU(2)$ subgroups of the $SU(N)$ link variables. It is known that simulations performed on $SU(2)$ subgroups suffer critical slowing down at larger N , but the largest N is only 7 and this problem did not appear. Lattices are spaced 100 sweeps of the lattice apart, for later analysis.

The spectroscopy was intended to be a typical $SU(3)$ lattice QCD calculation writ large. This involved an improved fermion action and extended sources for hadronic correlation functions. To achieve these goals a number of small technical problems had to be solved. Many of them have been encountered in other large- N studies, but perhaps my solutions are a bit different than what is found there, and might be worth reporting.

B. $SU(N)$ fat links; gauge fixing

My lattice fermions are clover fermions with normalized hypercubic (nHYP) smeared links as their gauge connections[35]. The clover coefficient is fixed at its tree level value, $c_{SW} = 1$. This particular discretization is known to work well, with small scaling violations, in both ordinary QCD phenomenology and in beyond-Standard Model studies in $SU(2)$, $SU(3)$, $SU(4)$. So the first technical problem is the construction of the nHYP link for arbitrary N . To describe the nHYP link in words, it is a local average of gauge connections over the hypercubes surrounding the link, which smears out the gauge field for the fermion. The specific problem to be solved is that the fat link is the average of a set of paths which produces a sum of $SU(N)$ matrices, call it Σ. The fat link V_μ is defined as

$$
V_{\mu} = \frac{\Sigma}{\sqrt{\Sigma^{\dagger}\Sigma}}.\tag{4}
$$

This is the matrix which maximizes Re Tr $V_\mu \Sigma^{\dagger}$. The quantity $Q^{-1/2} = 1/\sqrt{\Sigma^{\dagger}\Sigma}$ is computed using the Cayley - Hamilton theorem as described in Appendix B of Ref. [32]. This construction involves finding the eigenvalues of Q, which is done using a Jacobi algorithm.

As written, V_{μ} is an element of $U(N)$, not $SU(N)$. This is not an issue when it is used for the fermions since all that we care about is that our action should be gauge invariant, and under a gauge transformation both the thin and fat links transform the same way.

State of the art spectroscopic calculations use extended sources as interpolating fields for hadrons. In this work configuration of link variables is gauge fixed to lattice Coulomb gauge and the source for the quark propagator is some spatially extended function. So, we need to gauge fix our lattices to lattice Coulomb gauge, rotating the links to maximize $\sum_{x}\sum_{i}$ Re Tr $U_i(x)$, *i* labeling the spatial directions. This is done by finding a matrix V at each site of the lattice x which maximizes Re Tr $V\Sigma^{\dagger}$ where Σ is the sum of forward going $U_i(x)$'s emanating from x minus the sum of U_i^{\dagger} $\hat{i}^{\dagger}(x - \hat{i})$'s terminating on site x.

In the author's $SU(3)$ code, gauge fixing is done iteratively by sweeping through the lattice and, site by site, determining an optimal V by maximizing $\text{Re Tr } V\Sigma^{\dagger}$. As in the case of configuration generation, the relaxation is done using the $SU(2)$ subgroups of V. Unfortunately, for $N > 4$ relaxing on the subgroups suffers critical slowing down and it becomes impossible to carry out gauge fixing without performing an enormous number of iterations.

This is a variation of a problem which has been previously observed (and solved) by several groups simulating large N gauge theories, with $N > 10$ [36, 37]. There the problem is in the update step. Updating on $SU(2)$ subgroups produces a simulation algorithm with a long autocorrelation time, which becomes longer with increasing N . The solution is to perform over-relaxation on the full $SU(N)$ group, implementing an old idea of Creutz [38]. I do this a little differently than Refs. [36, 37]. The maximizing V is given by Eq, 4, which I have already dealt with while generating the fat links for the fermion action. The same application of the Cayley-Hamilton theorem gives V . As I already remarked, this V is an element of $U(N)$, not $SU(N)$. Now one needs $SU(N)$ elements, so I must compute the determinant of V, extracting its phase ϕ and performing an additional multiplication by the diagonal matrix $\exp(-i\phi/N)$. Doing this makes gauge fixing no more expensive for $SU(7)$ than for $SU(3)$.

C. Baryon operators

Mesonic states are constructed in the usual way, by sandwiching fermion propagators with Dirac matrices. Baryons are a bit more complicated. In $SU(3)$ it is common to use relativistic sources $(\epsilon_{abc}[u^a C \gamma_5 d^b]u^c$, for example, for the proton). I don't know a nice way to generalize this to $N > 3$, so I built baryon states using non-relativistic quark model states. So far, I have only considered $N_f = 2$ flavors of quarks. In what follows, I will usually label the flavors as u and d, although one can give them different masses – in $SU(3)$ the same operator can give the p or Ξ , for example.

In the MILC code, quark propagators are constructed in the Weyl $(\gamma_5$ diagonal) basis, and are rotated to γ_0 basis. I keep the two "large" components of the propagator's source and sink spin indices to be the fields which contract the quark model states. There are two choices for projection (states which are eigenvectors of γ_0 with eigenvalues ± 1 allow forward going in time and backward going two-component quark propagators). One projection has the lightest positive parity state in a channel as the forward-going state, and the lightest negative-parity state as the backward-going one. The situation is reversed for the other γ_0 projector. In order to reduce noise, these two propagators are summed (actually subtracted), configuration by configuration. This produces a single correlator extending halfway across the lattice, which is a candidate for fitting to a single decaying exponential.

A generic N−color baryon operator can be written as

$$
|B\rangle = \epsilon^{abc...N} \sum_{\{s_j\}} C_{\{s_j\}} u_{s_1}^a u_{s_2}^b \dots d_{s_N}^N |0\rangle.
$$
 (5)

The C's are an appropriate set of Clebsch-Gordan coefficients. The baryon propagator is a multiple sum over source and sink colors and spin coefficients of products of N quark propagators. This creates the possibility a large number of terms. Fortunately, many are redundant.

I condense states as follows. (I suspect that this is quite inefficient and that more efficiency is almost certainly possible.) I begin with states of definite u and d content. These states are eigenstates of I_3 , J , and J_3 . I then anticommute fermion fields into a "standard order," shown in Eq. 5: moving from the left, spin-up u quarks, spin-down u quarks, spin-up d quarks, and spin-down d quarks. The epsilon symbol absorbs the resulting minus signs.

Then, two examples are the $I = J = J_3 = N/2$ state

$$
|B, I = J = N/2\rangle = \epsilon^{abc...N} u_{\uparrow}^a \dots u_{\uparrow}^N |0\rangle
$$
 (6)

and the $SU(3)$ proton operator

$$
|p,\uparrow\rangle = \epsilon^{abc} \sqrt{\frac{2}{3}} (u_{\uparrow}^a u_{\uparrow}^b d_{\downarrow}^c - u_{\uparrow}^a u_{\downarrow}^b d_{\uparrow}^c) |0\rangle. \tag{7}
$$

Wick's theorem says that the n -quark propagator itself is also a determinant:

$$
\langle q_{i_1}(x)\bar{q}_{j_1}(y) q_{i_2}(x) \bar{q}_{j_2}(y) \dots q_{i_n}(x)\bar{q}_{j_n}(y) \rangle
$$

\n
$$
= (-)^n \sum_{P(1,2...n)} sign(P)(D_{i_1,j_{P_1}}^{-1}(x-y)D_{i_2,j_{P_2}}^{-1}(x-y) \dots D_{i_n,j_{P_n}}^{-1}(x-y)).
$$
\n(8)

 $(P \text{ is a permutation of indices.})$ The baryon propagator, then is a product of an up-quark determinant times a down quark determinant, summed over all the color combinations of the individual up and down quarks. Many terms are redundant in this product, as can be seen in an example, the $SU(3)$ proton: The contributions from the $\epsilon^{abc}u_{\uparrow}^a u_{\uparrow}^b$ source terms are antisymmetric in both Eqs. 5 and 8. This means that it is only necessary to keep one color ordering $(b > a$, for example) for each pair of colors (a, b) in the baryon propagator, and each term can be reweighted by a factor of 2!. This generalizes straightforwardly so that a a term in any wave function with N^u_\uparrow spin-up u quarks, N^u_\downarrow spin-down u quarks, N^d_\uparrow spin-up d quarks, N_{\downarrow}^{d} spin-down d quarks, picks up a restricted color sum, a single color ordering for each individual spin-flavor label and a multiplicity $N^u_\uparrow! N^u_\downarrow! N^d_\uparrow! N^d_\downarrow!$.

An extreme example of this pruning procedure is the the propagator for the $I = N/2, J =$ $N/2$ state of Eq. 6. It is

$$
\Delta(x, y) = (N!)^2 \det M(x, y)
$$
\n(9)

where M is the $N \times N$ matrix of u_{\uparrow} propagators from a source color to a sink one.

The lower the J, the more complicated are the wave functions. By $SU(7)$, the $J = 1/2$ states involve several hundred color combinations per spin configuration. This is getting quite unwieldy. (Since the nonrelativistic quark propagator is itself only a $(2N) \times (2N)$ matrix, there must be more redundancy.) But for now, I go on, naively.

To summarize, flavor $SU(2)$ wave functions and propagators for the various states are

- $I = J = N/2$: as already described, there is one determinant of an $N \times N$ matrix, for a cost N^3 .
- Analogs of Σ^* and Ξ^* states are

$$
|B, I = N/2 - 1, J = N/2\rangle = \sqrt{(N-1)!} (u_{\uparrow}^a \dots u_{\uparrow}^{N-1} s_{\uparrow}^N |0\rangle).
$$
 (10)

The baryon correlator is built of N^2 terms, one for each s color in the source and sink. Each is a determinant of an $(N-1) \times (N-1)$ matrix – the propagator of the up quarks.

• $I = J = N/2 - 1$: We have an $(N - 1) \times (N - 1)$ matrix of u-quark propagators and $N + N(N-1)$ color terms in the interpolating field for a cost of roughly $N^{4+3} = N^7$:

$$
|B, I = J = N/2 - 1\rangle = \sqrt{\frac{N-1}{N}} (u_1^a \dots u_1^{N-1} d_\downarrow^N - u_1^a \dots u_1^{N-2} u_\downarrow^{N-1} d_\uparrow^N) |0\rangle \tag{11}
$$

• $I = J = N/2 - 2$: An $(N - 2) \times (N - 2)$ matrix of propagators. The last term in the interpolating field has about N^3 color possibilities, for a cost of N^9 .

$$
|B, I = J = N/2 - 2\rangle \propto (u_{\uparrow}^{a} ... u_{\uparrow}^{N-2} d_{\downarrow}^{N-1} d_{\downarrow}^{N}
$$

$$
-2u_{\uparrow}^{a} ... u_{\uparrow}^{N-3} u_{\downarrow}^{N-2} d_{\uparrow}^{N-1} d_{\downarrow}^{N}
$$

$$
+u_{\uparrow}^{a} ... u_{\uparrow}^{N-4} u_{\downarrow}^{N-3} u_{\downarrow}^{N-2} d_{\uparrow}^{N-1} d_{\uparrow}^{N}) |0\rangle
$$

(12)

• $I = J = N/2 - 3$: our lowest state if $N = 7$; there are N^{11} terms to evaluate.

$$
|B, I = J = N/2 - 3\rangle \propto (u_{\uparrow}^{a} ... u_{\uparrow}^{N-3} d_{\downarrow}^{N-2} ... d_{\downarrow}^{N}
$$

\n
$$
-3u_{\uparrow}^{a} ... u_{\uparrow}^{N-4} u_{\downarrow}^{N-3} d_{\uparrow}^{N-2} d_{\downarrow}^{N-1} d_{\downarrow}^{N}
$$

\n
$$
+3u_{\uparrow}^{a} ... u_{\uparrow}^{N-5} u_{\downarrow}^{N-4} u_{\downarrow}^{N-3} d_{\uparrow}^{N-2} d_{\uparrow}^{N-1} d_{\downarrow}^{N}
$$

\n
$$
-u_{\uparrow}^{a} ... u_{\uparrow}^{N-6} u_{\downarrow}^{N-5} ... u_{\downarrow}^{N-3} d_{\uparrow}^{N-2} ... d_{\uparrow}^{N}) |0\rangle
$$

\n(13)

Still to be constructed are three flavor (u, d, s) states.

An obvious solution to the problem of increasing multiplicity is to prune states by doing an incomplete color sum. This will collide with another big problem: $SU(N)$ baryon signals degrade as N increases. For now, the only way to fight this is to collect larger data sets. Along the way, however, one can try to improve our signals by tactics such as averaging over the propagators with the same J and different m_J 's. (In practice, I combine the m_J and $-m_J$ propagators into a single correlator.) Another approach involves forcing a fit of several correlators which couple to the same states to a common mass. Ultimately, a variational calculation along the lines of what is done for excited state baryon spectroscopy might be necessary.

III. RESULTS FOR PURE GAUGE AND MESONIC OBSERVABLES

The simulations use the Wilson gauge action. Pure gauge systems with the Wilson action and $N = 2$ and 3 have a rapid crossover from strong to weak coupling, which becomes increasingly strong as N rises, converting to a real first order bulk transition by $N = 6$ or so. The location of the bulk transition has been tabulated in Ref. [3]. It is a lattice artifact.

The weak coupling side of the transition is the physical one from the point of view of making comparisons to continuum physics: only there is physics in the ultraviolet governed by asymptotic freedom so that the lattice spacing can be taken to zero by tuning the bare g^2 to zero. It is necessary to perform simulations above this transition, and I checked that my simulations were above it, by using the table and by doing my own simulations on smaller lattices to find the transition.

I present quenched results from three N's, $N = 3, 5, 7$. I have data sets of 80, 120, and 160 propagators, respectively. The bare gauge coupling β is tuned to match potentials through the Sommer parameters r_0 or r_1 . (Recall $r_0^2 F(r_0) = -1.65$. r_1 is the shorter-distance version of the Sommer parameter, $r_1^2 F(r_1) = -1.00$. In the real world, $r_1 \sim 0.31$ fm.)

In the present simulations, $\beta(N=3) = 6.0175$ (to match to previous work by Ref. [12]), and then the $N = 5$ and 7 couplings were tuned to match r_1 . Couplings and derived quantities are recorded in Table I. The bare 't Hooft couplings λ turn out to be quite similar. Potentials for the three values of N are shown in Fig. 1. They seem satisfactorily matched.

The combination $r_0\sqrt{\sigma}$ or $r_1\sqrt{\sigma}$ (σ is the string tension) gives a dimensionless combination of the Sommer radius and the string tension σ . The table shows that this quantity scales well.

My spectroscopy is based on smeared Gaussian sources and $\vec{p} = 0$ point sinks. At each N I collected sets for several different values of the width of R_0 for the source. These correlation functions are not variational since the source and sink are different. Thus, as R_0 is varied, the effective mass $(m_{eff}$ is defined through fitting correlators at a single distance to a single exponential; with open boundary conditions for correlator $C(t)$, $m_{eff} = \log C(t)/C(t+1)$ can approach its asymptotic value from above or below. I observe that typically, as R_0 rises, the mixed Gaussian - point correlators make their approach from above for smaller R_0 , and from below for bigger R_0 . An example of this behavior is shown in Fig. 2. (At large t, the signal deteriorates – a characteristic feature quite familiar from many $SU(3)$ studies. The noise is enhanced by the small data sets -40 lattices $-$ used to make the figure.) Then, rather than re-running the propagator code with yet another source, I can combine pairs of sources to produce a flat m_{eff} distribution using say

$$
C(t) = C(R_0 = 6, t) + fC(R_0 = 8, t). \tag{14}
$$

FIG. 1: Static potentials from $16^3 \times 32$ volumes: (a) $SU(3)$ at $\beta = 6.0175$; (b) $SU(5)$ at $\beta = 17.5$; (c) $SU(7)$ at $\beta = 34.9$. Effective mass fits for several values of t are overlaid.

For $SU(3)$, the optimal source is a linear combination of $R_0 = 4$ and 6 sources, favoring larger R_0 as the quark mass falls. For $SU(5)$ I mix sources with $R_0 = 6$ and 8. For $SU(7)$ the $R_0 = 8$ source produced flat effective mass plots across my mass range, and I did not do any source mixing.

Tables II-IV contain the resulting spectroscopy. The values of the masses are highly correlated because they come from the same underlying configurations. Mass differences, which will be described below, are taken from jackknife averages of the data.

FIG. 2: Effective masses for the highest-spin baryon for different size sources, labeled by octagons for $R_0 = 4$, diamonds for $R_0 = 6$, and squares, $R_0 = 8$. (a) $SU(3)$, $\kappa = 0.1265$; (b) $SU(5)$ 12³ volume, $\kappa = 0.1275$; (c) $SU(7)$, 12^3 volume, $\kappa = 0.129$.

Now I turn to results for mesonic observables. The critical hopping parameter κ_c is determined through the vanishing point for the Axial Ward Identity (AWI) quark mass, defined as

$$
\partial_t \sum_{\mathbf{x}} \langle A_0^a(\mathbf{x}, t) \mathcal{O}^a \rangle = 2m_q \sum_{\mathbf{x}} \langle P^a(\mathbf{x}, t) \mathcal{O}^a \rangle. \tag{15}
$$

where the axial current $A^a_\mu = \bar{\psi} \gamma_\mu \gamma_5 (\tau^a/2) \psi$, the pseudoscalar density $P^a = \bar{\psi} \gamma_5 (\tau^a/2) \psi$, and \mathcal{O}^a could be any source. Here it is my Gaussian shell model source.

FIG. 3: Additive mass shift versus $1/N^2$. My data are shown as octagons. The diamonds are the pure Wilson fermion data of Ref. [11], just for comparison over a wider N range.

The mass is shifted from its free value in the usual way. To compare to usual expectations, recall the relation between g^2C_F and λ : we expect the mass shift

$$
\delta m = \frac{1}{2\kappa_c} - 4\tag{16}
$$

to show $1/N^2$ variation. This I roughly see; compare Fig. 3. Just for comparison over a wider N range, I show the older results of Ref. [11], done with unimproved Wilson fermions.

Next, we turn to spectroscopy, shown for mesons in Fig. 4 and baryons in Fig. 5. Data are plotted in terms of the AWI quark mass, in units of r_1 , to make the x-axis the same for all N. The near independence of meson masses on N (when expressed in terms of a common variable) is apparent. This is not so, for baryons.

One is tempted to compare the chiral limit of the vector meson mass as a function of N. A simple linear fit to the data of Fig. 4 gives $r_1m_V = 1.58(2)$, 1.44(1) and 1.52(1) for $N = 3$, 5, 7. With a nominal $1/r_1 = 635$ MeV, this is about 900 MeV to 1 GeV versus 770 MeV for the physical rho meson. One always has to be careful with quenched lattice results from a single volume and lattice spacing, but the number is completely sensible. It is close to the results of Refs. [11, 12].

FIG. 4: Meson spectroscopy in units of r_1 from $N = 3$, 5 and 7, plotted as squares, diamonds and octagons, respectively. Panel (a) shows the squared pseudoscalar mass. Panel(b) shows the vector meson mass.

(Using completely different methodology plus a different choice of physical input to set the lattice scale, the authors of Ref. [14] have a mass which is from thirty per cent to a factor of two higher than this. The difference between their result and the low-N ones will probably only be resolved by removing the many differences in methodology one at at time.)

In what follows, I will replace the quark mass as the independent variable by the square of the ratio of the pseudoscalar to vector meson mass. The quark mass is scheme dependent; the ratio is not. Of course, nothing I say is affected by this choice.

Vacuum-to-meson matrix elements are expected to scale as \sqrt{N} . I have looked at the pseudoscalar, vector meson, and axial vector meson decay constants. To be explicit, they are defined as follows:

$$
\langle 0|\bar{u}\gamma_0\gamma_5 d|\pi\rangle = m_{\pi}f_{\pi} \tag{17}
$$

(so $f_{\pi} \sim 132$ MeV) while the vector meson decay constant of state V is defined as

$$
\langle 0|\bar{u}\gamma_i d|V\rangle = m_V^2 f_V \epsilon_i \tag{18}
$$

and the axial vector meson decay constant of state A is

$$
\langle 0|\bar{u}\gamma_i\gamma_5 d|A\rangle = m_A^2 f_A \epsilon_i. \tag{19}
$$

FIG. 5: Baryon spectroscopy in units of r_1 from $N = 3$, 5 and 7, plotted as squares, diamonds and octagons, respectively. For all N, higher J lies higher in mass: for $N = 3$ and 5 I have all states from $J = 1/2$ to $J = N/2$. For $N = 7$ only the $J = 7/2$, $5/2$, and $3/2$ states are shown.

 ϵ_i is a unit polarization vector. The lattice quantities f^L are converted to continuum convention by $f = f^L(1 - (3\kappa)/(4\kappa_c))$. I have left out the lattice-to-continuum Z-factor. With nHYP clover fermions, it is a few percent away from unity. (Ref. [39] has a table of one-loop Z-factors for various actions and HYP (hypercubic) links, which are a predecessor of nHYP links with the same one-loop Z-factor: For a current of type i, $Z_i = 1 + g^2 C_f/(16\pi^2)z_i$. The pseudoscalar factor is $z_P = 0.04$. The corresponding vector and axial vector factors are $z_V = -1.38$ and $z_A = -1.30$.)

The pseudoscalar decay constant is shown in Fig. 6. I show the dimensionless combination $r_1 f_{PS}/\sqrt{N}$. It's nice to see the \sqrt{N} scaling. Naive linear extrapolations give $r_1 f_\pi/\sqrt{N}$ 0.154(2), 0.151(1) and 0.154(2) for $N = 3, 5, 7$. The real world value is $1/\sqrt{3} \times 0.31$ fm $\times 132$ $MeV = 0.12$, so the quenched decay constant at this lattice spacing is coming in about 15 per cent high. Experts know that this kind of extrapolation is far too naive, to say nothing about comparing quenched QCD to the real world. Nevertheless, the answer is not absurd.

FIG. 6: Pseudoscalar decay constant, matched from the different simulations using the r_1 parameter from the potential and rescaled by $1/\sqrt{N}$. Squares, $N_c = 3$; diamonds, $N_c = 5$, octagons, $N_c = 7$.

Ref. [26] shows a figure with the ratio of quenched f_{π} to its experimental result about ten per cent high.

 \sqrt{N} scaling for the pseudoscalar decay constant was first observed by the authors of Ref. [13]. Their numerical result for the decay constant, translated to $SU(3)$, is also high, but by 40 per cent. Detecting the origin of this difference will probably again require detailed numerical work.

The same comparison is shown for f_V , and f_{a_1} in Fig. 7. The a_1 decay constant is quite noisy at small quark mass and I omit those results as untrustworthy. The \sqrt{N} scaling rule works well here, too.

IV. RESULTS FOR BARYONS – THE ROTOR SPECTRUM

My baryon data in Fig. 5 shows masses which increase roughly linearly in N . All data shows that for the baryons, higher J does lie higher in energy. This is no surprise, so let's look deeper. Fig. 8 shows the splitting between the various members of each multiplet. It

FIG. 7: Vector meson and axial vector meson decay constants, rescaled by $1/\sqrt{N}$. Squares, $N_c = 3$; diamonds, $N_c = 5$, octagons, $N_c = 7$.

is extracted using a jackknife average of differences of the two baryon masses.

I now demonstrate that the masses in Fig. 8 form a rotor spectrum. First, we can test the numerator of the rotor term of Eq. 1 N by N . This is done by looking at the ratio of differences

$$
\Delta(J_1, J_2, J_3) = \frac{M(N, J_2) - M(N, J_3)}{M(N, J_1) - M(N, J_3)},\tag{20}
$$

for which the constants (A, B) cancel. The result is shown in Fig. 9. I just plot one mass difference as a function of the other one and compare the data to a straight line of zero intercept whose slope is given by Eq. 20. The rotor spectrum is confirmed for all the members of the $N = 5$ and 7 multiplets I observe.

Second, one can look at the $J=3/2-1/2$ splitting, and check the N dependence for the bottom of the multiplets:

$$
M(N,3/2) = M(N,1/2) = \frac{3B}{N}.
$$
\n(21)

Given the states I have recorded, this can only be done for $N = 3$ and 5. Rescaling the mass difference by $N/3$ exposes B. The result is shown in Fig. 10(a). This is quite promising: the $N = 3$ and 5 data coincide.

Next, we can look at the top of the multiplet. Eq. 1 gives us a rescaled Landé interval rule: it says that the splitting between the $J = N/2$ and $J = N/2 - 1$ states is a constant,

FIG. 8: (a) Delta-proton $(J = 3/2 - 1/2)$ mass splitting versus quark mass, in quenched $SU(3)$. (b) $SU(5)$ mass splittings: octagons for $J = 5/2 - 1/2$, squares for $J = 5/2 - 3/2$, and crosses, for $J = 3/2 - 1/2$. (c) $SU(7)$ mass splittings: octagons for $J = 7/2 - 3/2$, squares for $J = 7/2 - 5/2$, and crosses, for $J = 5/2 - 3/2$. Eq. 1 says that the $SU(5)$ $J = 5/2 - 3/2$ splitting is supposed to be equal to the $SU(3)$ $J = 3/2 - 1/2$ mass splitting, and also to the $SU(7)$ $J = 7/2 - 5/2$ splitting.

B, independent of J. Fig. 10(b) shows this difference. It and panel (a) share the common $N = 3$ points, but the other points are different. The envelope of the curve is $B(m_q)$.

By design, these differences ignore the A term in Eq. 1. To get it, we can look at the

FIG. 9: Mass differences in the $SU(5)$ and $SU(7)$ multiplets, panels (a) and (b) respectively. For $SU(5)$ the line has a slope which is equal to the ratio $\Delta(5/2, 3/2, 1/2) = 3/8$ (recall Eq. 20). For $SU(7)$ the slope of the line is $\Delta(7/2, 5/2, 3/2) = 5/12$.

FIG. 10: Exposing the B term of Eq. 1: (a) $N/3$ times the $J = 3/2 - 1/2$ mass differences in the $SU(3)$ and $SU(5)$ multiplets, shown respectively as squares and diamonds. (b) The $J = N/2$ vs $J = N/2 - 1$ mass difference in the $SU(3)$, $SU(5)$, and $SU(7)$ multiplets, shown respectively as squares, diamonds, and octagons.

FIG. 11: Exposing the A term of Eq. 1: (a) Eq. 22 for the $SU(3)$ and $SU(5)$ multiplets, shown respectively as squares and diamonds. (b) The $J = N/2$ vs $J = N/2-1$ mass difference of Eq. 23 in the $SU(3)$, $SU(5)$, and $SU(7)$ multiplets, shown respectively as squares, diamonds, and octagons.

bottom of the multiplets,

$$
A = \frac{5}{4N}M(N, J = 1/2) - \frac{1}{4N}M(N, J = 3/2),\tag{22}
$$

or the top of the multiplets,

$$
A = \frac{N+2}{4N}M(N, J = N/2 - 1) - \frac{N-2}{4N}M(N, J = N/2).
$$
 (23)

Fig. 11 shows these two mass formulas. Again, they behave consistently.

Figs. 10 and 11 show that the A and B coefficients in Eq. 1 have typical hadronic sizes. Inserting a nominal lattice spacing, $1/a \sim 2100$ MeV, we see that $A = 400$ MeV at small quark mass and is an increasing function of quark mass. In quark model language, A is the constituent quark mass, and its value and dependence on quark mass are both quite reasonable. $B = 300$ MeV and falls with energy. Again this is a typical hadronic scale.

These values address an old conundrum of large-N phenomenology: the mass difference of the nucleon and Δ is supposed to be order Λ_{QCD}/N . It is measured to be about 300 MeV, which is itself order Λ_{QCD} . The plots of mass differences show that the nucleon- Δ mass splitting is indeed well-described by large-N QCD. Having data at several N's as well

as at several quark masses makes this result appear immediately. In particular, the data disfavor the large- N scenario of Ref. [40].

V. CONCLUSIONS

The qualitative features of large - N QCD phenomenology are easy to observe. The shape of the potential (characterized by the scaling combination $r_1\sqrt{\sigma}$) is N - independent. Meson masses show little N dependence – $m_Hr₁$ is N-independent. Vacuum - to -hadron matrix elements scale as \sqrt{N} . Two-flavor baryons show a prominent rotor spectrum. This seems to be true for both the bottom of the spectrum (low J) and the top $(J = N \text{ states})$.

Naively, one might expect that the wave function of the baryon would be N - independent. Lattice simulations do not generally directly reveal wave function information, but one might expect that the same interpolating fields might behave the same for different N 's. That does not seem to be the case; larger N seems to prefer larger Gaussian trial wave functions. But perhaps even $N = 7$ is not such large N and one should push farther.

Readers who do lattice simulations can undoubtedly list many obvious extensions of this project: smaller lattice spacing, to do an honest extrapolation to the continuum limit, bigger volumes to check that the answers are trustworthy, smaller quark masses, and bigger N 's are obvious technical improvements. Giving the two fermions different masses, or better yet, considering flavor $SU(3)$, would allow more tests of large-N spectroscopy. One might really want to test whether dynamical fermions become less important at large N. Writing the code to simulate $SU(N)$ fundamental fermions is straightforward; running it might be costly, and seeing the expected small differences shrink as N rises would be even more costly.

The continuum literature of large - N baryons is more than thirty years old, and I clearly have only scratched its surface in this paper. Several simple ingredients were useful: having data at several N 's, at several J 's for each N , and having data at many quark masses. Most of the continuum literature I have read restricts itself to statements about $J = 1/2$ and $3/2$ – presumably because that is all that exists in experimental data. Predictions for any J can challenge the lattice and would be candidates for future studies.

	SU(3)	SU(5)	SU(7)
β	6.0175	17.5	34.9
$\lambda = q^2 N$	2.99	2.86	2.81
r_0/a	5.36(6)	5.20(5)	5.41(3)
r_1/a	3.90(3)	3.77(3)	3.91(2)
$r_0\sqrt{\sigma}$		$1.175(3)$ $1.172(3)$ $1.167(2)$	
$r_1\sqrt{\sigma}$		$0.856(5)$ $0.850(4)$ $0.845(2)$	

TABLE I: Bare parameters and observables from potentials.

κ	am_q	am_{PS}	am_V	$am_B(J=3/2)$ $am_B(J=1/2)$	
		0.1230 0.119 $0.525(2)$ $0.617(3)$		0.996(6)	0.942(5)
		0.1240 0.089 $0.449(2)$ $0.561(3)$		0.915(7)	0.846(5)
		0.1250 0.055 $0.362(2)$ $0.505(4)$		0.835(8)	0.744(6)
		0.1253 0.046 $0.333(2)$ $0.491(5)$		0.804(9)	0.700(6)
		0.1260 0.029 $0.255(3)$ $0.457(6)$		0.747(12)	0.613(7)
			0.1265 0.014 $0.184(4)$ $0.427(10)$	0.703(13)	0.516(12)

TABLE II: Spectra from $SU(3)$ simulations.

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κ	am_a	am_{PS}	am_V		$am_B(J=5/2)$ $am_B(J=3/2)$ $am_B(J=1/2)$	
		0.1240 0.127 $0.565(1)$ $0.655(1)$		1.864(6)	1.814(5)	1.786(5)
		0.1250 0.099 $0.488(1)$ $0.593(1)$		1.708(6)	1.650(6)	1.616(5)
		0.1260 0.070 $0.403(1)$ $0.532(2)$		1.566(7)	1.488(6)	1.446(6)
		0.1265 0.055 $0.356(1)$ $0.500(2)$		1.486(7)	1.404(6)	1.356(6)
		0.1270 0.041 $0.302(2)$ $0.469(3)$		1.419(8)	1.325(7)	1.270(7)
		0.1275 0.026 $0.240(2)$ $0.440(4)$		1.357(10)	1.243(9)	1.177(8)

TABLE III: Spectra from $SU(5)$ simulations.

κ	am_a	am_{PS}	am_V		$am_B(J = 7/2)$ $am_B(J = 5/2)$ $am_B(J = 3/2)$	
		0.1260 0.115 $0.565(1)$ $0.663(1)$		2.681(11)	2.649(8)	2.612(8)
		0.1270 0.088 $0.488(1)$ $0.603(1)$		2.483(9)	2.420(10)	2.386(9)
		0.1280 0.062 $0.401(1)$ $0.537(2)$		2.279(11)	2.215(11)	2.174(10)
		0.1290 0.036 $0.299(1)$ $0.471(3)$		2.082(15)	1.994(12)	1.929(11)

TABLE IV: Spectra from $SU(7)$ simulations.

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