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Results from lattice simulations of $\mathcal{N} = 4$ supersymmetric Yang–Mills

Simon Catterall^{* a}, Joel Giedt^{* b}, David Schaich^{* a}, Poul H. Damgaard^c and Thomas DeGrand^d

- ^a Department of Physics, Syracuse University, Syracuse, NY 13244, USA
- ^b Department of Physics, Applied Physics and Astronomy, Rensselaer Polytechnic Institute, Troy NY 12065, USA
- ^c Niels Bohr International Academy and Discovery Center, Niels Bohr Institute, University of Copenhagen, Copenhagen, Denmark
- ^d Department of Physics, University of Colorado, Boulder, CO 80309, USA

We report recent results and developments from our ongoing lattice studies of $\mathcal{N} = 4$ supersymmetric Yang–Mills theory. These include a proof that only a single fine-tuning needs to be performed, so long as the moduli space is not lifted by nonperturbative effects. We extend our investigations of supersymmetry restoration in the continuum limit by initiating Monte Carlo renormalization group studies. We present additional numerical evidence that the lattice theory does not suffer from a sign problem. Finally we study the static potential, which we find to be Coulombic at both weak and strong coupling. We compare the static potential Coulomb coefficients to perturbation theory, including initial results for N = 3 colors in addition to N = 2.

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*Speaker.

 $\mathcal{N} = 4$ supersymmetric Yang–Mills (SYM) is a fascinating quantum field theory which plays an important role in many areas of theoretical physics including holographic approaches to quantum gravity, understanding the structure of scattering amplitudes in $\mathcal{N} = 8$ supergravity, and the conformal bootstrap program. It is the only known example of a four-dimensional theory admitting a supersymmetric lattice discretization, and the only non-trivial example of a lattice gauge theory with a β function that vanishes at least at one loop. (The β function of the continuum theory vanishes to all orders.) Many people have contributed to the lattice formulation that we employ in our studies; see the review [1] and references therein for more information. Alternate approaches to numerically studying $\mathcal{N} = 4$ SYM include Refs. [2, 3, 4, 5, 6, 7, 8].

In this proceedings we report recent results and developments from our ongoing lattice studies of $\mathcal{N} = 4$ SYM. Some of these results are new and preliminary; others have appeared in Refs. [9, 10, 11, 12, 13]. In Sections 1–3 we summarize the main elements of the lattice formulation. In particular, in Section 2 we show that so long as the moduli space is not lifted by nonperturbative effects only a single marginal operator must be tuned to recover the $\mathcal{N} = 4$ supersymmetric continuum limit. Sections 3 and 4 discuss observables that we measure to monitor supersymmetry restoration in the continuum limit, including initial results from new Monte Carlo renormalization group (MCRG) studies. We briefly update our measurements of the Pfaffian phase in Section 5, exploring more lattice volumes for larger gauge groups. The Pfaffian remains nearly real and positive in all cases, further evidence that the lattice theory does not suffer from a sign problem. We focus on the static potential in Section 6, presenting initial results for N = 3 colors. We conclude with some discussion of the next steps planned for our wide-ranging investigations.

To encourage independent work on lattice $\mathcal{N} = 4$ SYM, we have developed a publicly available parallel software package suitable for use on clusters and supercomputers.¹ This software evolved from the MILC code for lattice QCD, generalized to handle fermions in the adjoint representation for an arbitrary number of colors. Its central feature is rational hybrid Monte Carlo importance sampling. Many additional measurements are provided, including parallel computation of the complex Pfaffian of the fermion operator. A detailed presentation recently appeared in Ref. [14].

1. Discretizing the twisted continuum action

The lattice theory we employ results from discretizing a topologically twisted form of continuum $\mathcal{N} = 4$ SYM with gauge group SU(N). There are three independent topological twists of $\mathcal{N} = 4$ SYM, of which we employ the Marcus or Geometric-Langlands twist [15, 16]. This twisted theory is most conveniently written as a dimensional reduction of a five-dimensional action

$$S = \frac{N}{2\lambda} \mathcal{Q} \int d^5 x \left(\chi_{ab} \mathcal{F}_{ab} + \eta \left[\overline{\mathcal{D}}_a, \mathcal{D}_a \right] + \frac{1}{2} \eta d \right) - \frac{N}{8\lambda} \int d^5 x \, \epsilon_{abcde} \, \chi_{ab} \overline{\mathcal{D}}_c \chi_{de}. \tag{1.1}$$

 $\lambda = g^2 N$ is the 't Hooft coupling. The gauge fields \mathcal{A}_a (and their associated covariant derivatives \mathcal{D}_a and $\overline{\mathcal{D}}_a$) have real parts which are just the usual Yang–Mills gauge fields, while their imaginary parts arise from the five scalars that appear after reducing $\mathcal{N} = 1$ SYM in ten dimensions down

¹http://github.com/daschaich/susy

to five dimensions. The appearance of the scalar fields as vectors is odd at first sight, but is at the heart of the twisting process.

Twisting decomposes the fields under the diagonal subgroup of the Euclidean Lorentz symmetry and the R symmetry of the theory. Since the scalars transform as vectors under the R symmetry they remain vectors under the twisted symmetry, and naturally combine with the R-singlet gauge fields. Similar arguments show that twisting transforms the original four Majorana fermions into a set of antisymmetric fields with integer spins under the twisted symmetry, appearing as $(\eta, \psi_a, \chi_{ab})$ in the above action. Likewise the $\mathcal{N} = 4$ fermionic supercharges $Q_{\alpha i}$, $Q_{\alpha i}^{\dagger}$ become integer-spin $(\mathcal{Q}, \mathcal{Q}_a, \mathcal{Q}_{ab})$. In flat space twisting does not change the physical content of the theory. It is merely an exotic change of variables, to a spinor-free form better suited to discretization.

The first part of the action in Eq. 1.1 is the Q variation of a function. This scalar supercharge Q is the supersymmetry that can be preserved under discretization. It acts on the fields as follows:

$$Q \mathcal{A}_{a} = \psi_{a} \qquad \qquad \mathcal{Q} \psi_{a} = 0$$

$$Q \chi_{ab} = -\overline{\mathcal{F}}_{ab} \qquad \qquad \mathcal{Q} \overline{\mathcal{A}}_{a} = 0 \qquad (1.2)$$

$$Q \eta = d \qquad \qquad \mathcal{Q} \mathcal{A} = 0.$$

Its nilpotent character guarantees that the first piece of the action is trivially Q invariant. The second term can be shown to be invariant using the Bianchi identity, so that QS = 0.

We recover the standard Marcus twist of four-dimensional $\mathcal{N} = 4$ SYM from naive dimensional reduction, by simply setting to zero the momentum in the fifth dimension, $\partial_4 = 0$, and identifying the corresponding component of the gauge field as the sixth scalar field, $\mathcal{A}_4 \to \phi$. A more symmetric choice is to project the theory onto the four-dimensional hyperplane with normal vector $\hat{\mathbf{n}} = \frac{1}{\sqrt{5}}(1, 1, 1, 1, 1)$. This corresponds to setting $\sum_{a=0}^4 \partial_a = 0$ and $\sum_{a=0}^4 \mathcal{A}_a \to \phi$, preserving an S_5 permutation symmetry.

Using this formalism, the transition to the lattice is quite straightforward. The lattice structure follows from a similar dimensional reduction. Starting with the five-dimensional hypercubic lattice in momentum space, the constraint $\sum_{a=0}^{4} \partial_a = 0 \rightarrow \sum_{a=0}^{4} n_a = 0$ produces the A_4 lattice. Transforming back to real space leaves us with the dual A_4^* lattice, which retains an S_5 point group symmetry manifest as five (linearly dependent) basis vectors symmetrically spanning four spacetime dimensions. This S_5 symmetry provides a set of irreducible representations that match those of the continuum twisted SO(4) symmetry [17]. It can be useful to think of A_4^* as the four-dimensional analog of the triangular lattice A_2^* in two dimensions, or as the weight lattice of SU(5).

The complex gauge fields $\mathcal{A}_a(x)$ become complexified $\mathcal{U}_a(n)$ living on the five links of the A_4^* lattice. The form of the nilpotent scalar supersymmetry \mathcal{Q} remains the same as in the continuum (Eq. 1.2), and requires that these $\mathcal{U}_a(n)$ be elements of the algebra $\mathfrak{gl}(N, \mathbb{C})$, with consequences we will discuss in Section 3. Supersymmetry also forces us to assign the ψ_a fermions to the same links, while the η fermion and bosonic auxiliary field d are placed on sites. The χ_{ab} fermions are associated with the field strengths

$$\overline{\mathcal{F}}_{ab}(n) = \overline{\mathcal{U}}_a(n + \widehat{\mu}_b)\overline{\mathcal{U}}_b(n) - \overline{\mathcal{U}}_b(n + \widehat{\mu}_a)\overline{\mathcal{U}}_a(n),$$

and connect lattice sites $n + \hat{\mu}_a + \hat{\mu}_b$ and n so that their contractions with

$$\mathcal{D}_{a}^{(+)}\psi_{b} = \mathcal{U}_{a}(n)\psi_{b}(n+\widehat{\mu}_{a}) - \psi_{b}(n)\mathcal{U}_{a}(n+\widehat{\mu}_{b})$$

will produce gauge invariant loops. After integrating out the auxiliary field d, on the lattice the two terms in Eq. 1.1 become

$$S_{exact} = \frac{N}{2\lambda_{\text{lat}}} \sum_{n} a^{4} \operatorname{Tr} \left[-\overline{\mathcal{F}}_{ab} \mathcal{F}_{ab}(n) + \frac{1}{2} \left(\overline{\mathcal{D}}_{a}^{(-)} \mathcal{U}_{a}(n) \right)^{2} - \chi_{ab} \mathcal{D}_{[a}^{(+)} \psi_{b]}(n) - \eta \overline{\mathcal{D}}_{a}^{(-)} \psi_{a}(n) \right]$$
$$S_{closed} = -\frac{N}{8\lambda_{\text{lat}}} \sum_{n} a^{4} \operatorname{Tr} \left[\epsilon_{abcde} \chi_{de}(n + \widehat{\mu}_{a} + \widehat{\mu}_{b} + \widehat{\mu}_{c}) \overline{\mathcal{D}}_{c}^{(-)} \chi_{ab}(n) \right], \tag{1.3}$$

where the specific form of the forward and backward difference operators can be found in Ref. [12].

This exactly supersymmetric lattice action was studied analytically in Refs. [9, 11, 13]. These studies have found that the single exact supersymmetry suffices to establish the following results:

- The moduli space survives to all orders of lattice perturbation theory, which prohibits any scalar potential being induced through radiative corrections.
- The β function vanishes at least at one loop in lattice perturbation theory.
- Certain quantities, such as the partition function, may be computed exactly in the semiclassical limit.
- Lattice symmetries imply that only a single tuning of a marginal operator is necessary to restore R symmetry and all supersymmetries in the continuum limit.

We derive the last result in the next section.

2. Renormalization and the long-distance effective theory

Since $\mathcal{N} = 4$ SYM possesses a line of conformal fixed points, we reach the continuum limit of the lattice theory by taking $1/L \rightarrow 0$ for any fixed 't Hooft coupling λ , where L is the linear length of the lattice volume. We must ensure that the long-distance effective action S_{eff} of the lattice theory appropriately recovers $\mathcal{N} = 4$ SYM in the continuum limit. Simply defining S_{eff} requires the existence of a real-space renormalization group (RG) blocking transformation that preserves both the symmetries of the system and the geometric interpretation of the fields. In this section we will present an explicit example of such a blocking scheme, derive the long-distance effective action, and show that only a single tuning is required if the moduli space is not lifted by nonperturbative effects. These results were recently derived in Ref. [13].

Writing the original A_4^* lattice as $\Lambda = \left\{ a \sum_{i=0}^3 n_i \hat{\mu}_i \mid n \in \mathbb{Z}^4 \right\}$, with $\hat{\mu}_i$ the first four of the five (degenerate) basis vectors, the blocked lattice $\Lambda' = \left\{ 2a \sum_{i=0}^3 n_i \hat{\mu}_i \mid n \in \mathbb{Z}^4 \right\}$ is merely doubled in every direction. The blocked fields must begin and end on sites of Λ' . We will denote the blocked fields by primes, and work in lattice units with a = 1. The following blocking transforma-

tion preserves both the Q algebra and the geometric interpretation of the fields:

$$\begin{aligned} \mathcal{U}_{a}'(n) &= \xi \mathcal{U}_{a}(n) \mathcal{U}_{a}(n+\widehat{\mu}_{a}) & \overline{\mathcal{U}}_{a}'(n) = \xi \overline{\mathcal{U}}_{a}(n+\widehat{\mu}_{a}) \overline{\mathcal{U}}_{a}(n) \\ d'(n) &= d(n) & \eta'(n) = \eta(n) \\ \psi_{a}'(n) &= \xi \left[\psi_{a}(n) \mathcal{U}_{a}(n+\widehat{\mu}_{a}) + \mathcal{U}_{a}(n) \psi_{a}(n+\widehat{\mu}_{a}) \right] & (2.1) \\ \chi_{ab}'(n) &= \frac{\xi^{2}}{2} \left[\overline{\mathcal{U}}_{a}(n+\widehat{\mu}_{a}+2\widehat{\mu}_{b}) \overline{\mathcal{U}}_{b}(n+\widehat{\mu}_{a}+\widehat{\mu}_{b}) + \overline{\mathcal{U}}_{b}(n+2\widehat{\mu}_{a}+\widehat{\mu}_{b}) \overline{\mathcal{U}}_{a}(n+\widehat{\mu}_{a}+\widehat{\mu}_{b}) \right] \chi_{ab}(n) \\ &+ \xi^{2} \left[\overline{\mathcal{U}}_{a}(n+\widehat{\mu}_{a}+2\widehat{\mu}_{b}) \chi_{ab}(n+\widehat{\mu}_{b}) \overline{\mathcal{U}}_{b}(n) + \overline{\mathcal{U}}_{b}(n+2\widehat{\mu}_{a}+\widehat{\mu}_{b}) \chi_{ab}(n+\widehat{\mu}_{a}) \overline{\mathcal{U}}_{a}(n) \right] \\ &+ \frac{\xi^{2}}{2} \chi_{ab}(n+\widehat{\mu}_{a}+\widehat{\mu}_{b}) \left[\overline{\mathcal{U}}_{a}(n+\widehat{\mu}_{b}) \overline{\mathcal{U}}_{b}(n) + \overline{\mathcal{U}}_{b}(n+\widehat{\mu}_{a}) \overline{\mathcal{U}}_{a}(n) \right]. \end{aligned}$$

Because the link variables are non-compact elements of $\mathfrak{gl}(N,\mathbb{C})$, we allow for the possibility that they are rescaled by a factor ξ under the transformation. For the site variables η and d we simply use decimation. From Eq. 2.1 it is easy to see that \mathcal{U}'_a and ψ'_a connect lattice sites n and $n + 2\hat{\mu}_a$, while χ'_{ab} connects $n + 2\hat{\mu}_a + 2\hat{\mu}_b$ and n, as desired. As a consequence, the properties of the system under the S_5 point group symmetry are preserved. Any S_5 invariant of the original fields, such as $\sum_a \mathcal{U}_a \overline{\mathcal{U}}_a$, remains invariant when expressed in terms of the blocked fields.

It is obvious that $\mathcal{Q}\overline{\mathcal{U}}_{a}' = 0$, $\mathcal{Q}\eta' = d'$ and $\mathcal{Q}d' = 0$, just as in Eq. 1.2. $\mathcal{Q}\mathcal{U}_{a}' = \psi_{a}'$ also follows from the original algebra, while $\mathcal{Q}\psi_{a}' = 0$ due to the negative sign from anticommuting \mathcal{Q} past ψ_{a} :

$$\mathcal{Q}\psi_a'(n) = \xi \left[-\psi_a(n)\psi_a(n+\widehat{\mu}_a) + \psi_a(n)\psi_a(n+\widehat{\mu}_a)\right] = 0$$

Finally, defining the blocked field strength as expected,

$$\overline{\mathcal{F}}'_{ab}(n) = \overline{\mathcal{U}}'_a(n+\widehat{\mu}_b)\overline{\mathcal{U}}'_b(n) - \overline{\mathcal{U}}'_b(n+\widehat{\mu}_a)\overline{\mathcal{U}}'_a(n),$$

just a few lines of algebra are required to check $Q\chi'_{ab} = \mathcal{F}'_{ab}$. So we see that the blocking scheme in Eq. 2.1 preserves the supersymmetry $Q^2 = 0$ along with the other symmetries of the system.

We can now consider what lattice operators could possibly be generated under RG flow based on this blocking scheme. Those that give relevant or marginal operators in the continuum limit must be included in the most general long-distance effective action S_{eff} . If two lattice operators produce the same relevant or marginal continuum operator, and only differ by irrelevant operators in the continuum limit, then we only need to include one of them in S_{eff} . We must ensure that every operator in the effective action acquires its canonical coefficient, either by fine-tuning the corresponding coefficients in the UV theory, Eq. 1.3, or by adding new counterterms to the lattice action. As we will describe below, some operators can be given their canonical coefficients simply by rescaling the fields.

In both the continuum and lattice theories, there is only one renormalizable Q-closed operator: the second line of Eq. 1.3 gives the second term in Eq. 1.1. The only renormalizable Q-exact terms must take the form QTr $[\Psi f(U, \overline{U}, d)]$ or $Q \{ \text{Tr} [\eta] \text{Tr} [f(U, \overline{U}, d)] \}$ where Ψ stands for one of the fermion fields. Q must act on a fermionic quantity so that S_{eff} is bosonic, while cubic or higher powers of fermions would be nonrenormalizable. Only η can be used in double-trace operators, since traces must be evaluated at sites to be gauge invariant. Finally, the effective action must be invariant under two further global symmetries of the system: the shift symmetry

$$\eta \to \eta + c \mathbb{I}_N, \tag{2.2}$$

with c a constant Grassmann parameter, and a U(1) "ghost number" symmetry [1] corresponding to the untwisted part of the SO(6) R symmetry.

For $\Psi = \eta$ these requirements leave only three possible terms,

$$\mathcal{Q}\mathrm{Tr}\left[\eta\overline{\mathcal{D}}_{a}^{(-)}\mathcal{U}_{a}\right] \qquad \mathcal{Q}\mathrm{Tr}\left[\eta d\right] \qquad \mathcal{Q}\mathrm{Tr}\left[\eta\mathcal{U}_{a}\overline{\mathcal{U}}_{a}\right] - \frac{1}{N}\mathcal{Q}\left\{\mathrm{Tr}\left[\eta\right]\mathrm{Tr}\left[\mathcal{U}_{a}\overline{\mathcal{U}}_{a}\right]\right\}, \qquad (2.3)$$

the first two of which are already present in Eq. 1.3. The only nontrivial operator with $\Psi = \psi_a$ is forbidden by the U(1) ghost number symmetry [13]. Finally, for $\Psi = \chi_{ab}$ the antisymmetry of χ_{ab} requires that $Q \operatorname{Tr} [\chi_{ab} \mathcal{U}_a \mathcal{U}_b]$ and $Q \operatorname{Tr} [\chi_{ab} \mathcal{U}_b \mathcal{U}_a]$ be combined with a negative sign, leaving

$$Q\mathrm{Tr}\left[\chi_{ab}\mathcal{F}_{ab}\right] \tag{2.4}$$

which is also already in Eq. 1.3. These arguments only hold because the blocking scheme in Eq. 2.1 preserves the geometric interpretation of the blocked fields on Λ' , so that the same terms appear in both the original action and the long-distance effective theory. It is also important that the blocking preserves the S_5 symmetry, without which many other operators would have been generated.

Thus the most general long-distance effective action is

$$S_{eff} = \mathcal{Q} \operatorname{Tr} \left[\alpha_1 \chi_{ab} \mathcal{F}_{ab} + \alpha_2 \eta \overline{\mathcal{D}}_a^{(-)} \mathcal{U}_a - \frac{\alpha_3}{2} \eta d \right] - \frac{\alpha_4}{4} \epsilon_{abcde} \operatorname{Tr} \left[\chi_{de} \overline{\mathcal{D}}_c^{(-)} \chi_{ab} \right] + \beta \mathcal{Q} \left\{ \operatorname{Tr} \left[\eta \mathcal{U}_a \overline{\mathcal{U}}_a \right] - \frac{1}{N} \operatorname{Tr} \left[\eta \right] \operatorname{Tr} \left[\mathcal{U}_a \overline{\mathcal{U}}_a \right] \right\},$$
(2.5)

suppressing the overall $\frac{N}{2\lambda_{\text{lat}}} \sum_n a^4$ for brevity. Acting with \mathcal{Q} , followed by the rescaling

$$\eta \to \lambda_\eta \eta \qquad \psi_a \to \lambda_\psi \psi_a \qquad \chi_{ab} \to \lambda_\chi \chi_{ab} \qquad d \to \lambda_d d,$$
 (2.6)

$$S_{eff} = \operatorname{Tr} \left[-\alpha_1 \overline{\mathcal{F}}_{ab} \mathcal{F}_{ab} - \alpha_1 \lambda_{\chi} \lambda_{\psi} \chi_{ab} \mathcal{D}_{[a}^{(+)} \psi_{b]} + \alpha_2 \lambda_d d \overline{\mathcal{D}}_a^{(-)} \mathcal{U}_a - \alpha_2 \lambda_\eta \lambda_{\psi} \eta \overline{\mathcal{D}}_a^{(-)} \psi_a - \frac{\alpha_3}{2} \lambda_d^2 d^2 \right] - \frac{\alpha_4}{4} \lambda_{\chi}^2 \epsilon_{abcde} \operatorname{Tr} \left[\chi_{de} \overline{\mathcal{D}}_c^{(-)} \chi_{ab} \right]$$

$$(2.7)$$

$$+\beta \left\{ \lambda_{d} \operatorname{Tr} \left[d\mathcal{U}_{a} \overline{\mathcal{U}}_{a} \right] - \lambda_{\eta} \lambda_{\psi} \operatorname{Tr} \left[\eta \psi_{a} \overline{\mathcal{U}}_{a} \right] - \frac{\lambda_{d}}{N} \operatorname{Tr} \left[d \right] \operatorname{Tr} \left[\mathcal{U}_{a} \overline{\mathcal{U}}_{a} \right] + \frac{\lambda_{\eta} \lambda_{\psi}}{N} \operatorname{Tr} \left[\eta \right] \operatorname{Tr} \left[\psi_{a} \overline{\mathcal{U}}_{a} \right] \right\}.$$

We are now free to set many of the coefficients to α_1 by imposing four constraints,

$$\alpha_1 \lambda_\chi \lambda_\psi = \alpha_1 \qquad \alpha_2 \lambda_d = \alpha_1 \qquad \alpha_2 \lambda_\eta \lambda_\psi = \alpha_1 \qquad \alpha_4 \lambda_\chi^2 = \alpha_1.$$
 (2.8)

Solving this system of equations produces the rescaling

$$\lambda_{\eta} = \sqrt{\frac{\alpha_1^3}{\alpha_4 \alpha_2^2}} \qquad \qquad \lambda_{\chi} = \frac{1}{\lambda_{\psi}} = \sqrt{\frac{\alpha_1}{\alpha_4}} \qquad \qquad \lambda_d = \frac{\alpha_1}{\alpha_2}. \tag{2.9}$$

We can then absorb an overall factor of α_1 into the renormalized gauge coupling, which does not need to be tuned since the continuum theory is conformal for any λ . This leaves

$$S_{eff} = \operatorname{Tr} \left[-\overline{\mathcal{F}}_{ab} \mathcal{F}_{ab} - \chi_{ab} \mathcal{D}_{[a}^{(+)} \psi_{b]} + d\overline{\mathcal{D}}_{a}^{(-)} \mathcal{U}_{a} - \eta \overline{\mathcal{D}}_{a}^{(-)} \psi_{a} - \frac{1}{2} \frac{\alpha_{1} \alpha_{3}}{\alpha_{2}^{2}} d^{2} \right] - \frac{1}{4} \epsilon_{abcde} \operatorname{Tr} \left[\chi_{de} \overline{\mathcal{D}}_{c}^{(-)} \chi_{ab} \right]$$

$$+ \frac{\beta}{\alpha_{2}} \left\{ \operatorname{Tr} \left[d\mathcal{U}_{a} \overline{\mathcal{U}}_{a} \right] - \operatorname{Tr} \left[\eta \psi_{a} \overline{\mathcal{U}}_{a} \right] - \frac{1}{N} \operatorname{Tr} \left[d \right] \operatorname{Tr} \left[\mathcal{U}_{a} \overline{\mathcal{U}}_{a} \right] + \frac{1}{N} \operatorname{Tr} \left[\eta \right] \operatorname{Tr} \left[\psi_{a} \overline{\mathcal{U}}_{a} \right] \right\}.$$

$$(2.10)$$

At most two fine-tunings are required to obtain continuum $\mathcal{N} = 4$ SYM: $\alpha_3 \rightarrow \alpha_2^2/\alpha_1$ and $\beta \rightarrow 0$. In particular, it is remarkable that the relation $\lambda_d = \lambda_\eta \lambda_\psi$ keeps the β term from bifurcating into multiple operators with different coefficients. We can already appreciate the benefits of maintaining an exact supersymmetry at nonzero lattice spacing by comparing these two tunings against the eight that would be necessary with naive Wilson fermions [13]. In addition, as mentioned at the end of Section 1, this exact supersymmetry ensures that the moduli space survives to all orders of lattice perturbation theory [9]. Ref. [13] demonstrates that the β term in Eq. 2.10 lifts the moduli space. Therefore if nonperturbative effects such as instantons also preserve the moduli space, then the β term is forbidden and $\beta = 0$ is preserved by the RG flow.

In the absence of the β term, we can easily integrate out the auxiliary field d in Eq. 2.10. The resulting long-distance effective action reproduces Eq. 1.3 up to one change:

$$\frac{1}{2} \left(\overline{\mathcal{D}}_a^{(-)} \mathcal{U}_a(n) \right)^2 \longrightarrow \frac{c_2}{2} \left(\overline{\mathcal{D}}_a^{(-)} \mathcal{U}_a(n) \right)^2 \tag{2.11}$$

in S_{exact} , with $c_2 = 2 - \alpha_2^2/(\alpha_1 \alpha_3)$. That is, a single fine-tuning of this marginal c_2 should suffice to recover $\mathcal{N} = 4$ SYM in the continuum limit defined by $1/L \to 0$ with fixed λ_{lat} . We have begun to explore this tuning in numerical calculations. Before we describe our initial results in Section 4, we must discuss certain sources of soft \mathcal{Q} supersymmetry breaking that are required to stabilize numerical simulations of lattice $\mathcal{N} = 4$ SYM.

3. Stabilizing numerical computations

For the lattice theory to recover $\mathcal{N} = 4$ SYM in the continuum limit, we must also require that the lattice gauge fields have the expansion

$$\mathcal{U}_a(n) = \frac{1}{a} \mathbb{I}_N + \mathcal{A}_a(x) + \mathcal{O}(a)$$
(3.1)

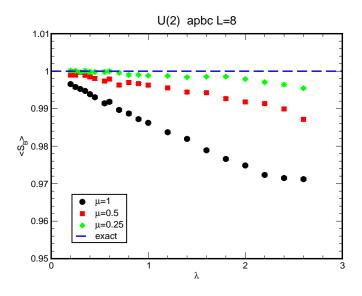


Figure 1: Relative deviations of the bosonic action from its exact supersymmetric value, $\langle s_B \rangle / 18$, plotted vs. the 't Hooft coupling λ_{lat} for 8^4 lattices with several nonzero values of the Q-breaking coefficient μ in Eq. 3.2. As $\mu \to 0$, $\langle s_B \rangle / 18 \to 1$ for all λ_{lat} , indicating the restoration of supersymmetry. From Ref. [10].

in some appropriate gauge. Without the unit matrix the lattice \mathcal{F}_{ab} would lack the derivative (kinetic) terms required to correspond to the continuum field strength. If the link variables were elements of the gauge group this unit matrix would arise automatically from expanding the exponential of terms in the algebra. Eq. 3.1 is nontrivial since our links are themselves elements of the algebra $\mathfrak{gl}(N,\mathbb{C})$, which also implies that the lattice gauge group is U(N) as opposed to the target SU(N) [18, 19, 20]. In the continuum limit the U(1) degrees of freedom decouple as $U(N) = SU(N) \otimes U(1)$, and below we will discuss how to suppress lattice artifacts associated with the U(1) sector. In addition, the integration measure for the lattice gauge fields is not the usual Haar measure but a flat measure. This flat measure is gauge invariant, since the fields are complex and the measure contains both $D\mathcal{U}$ and $D\overline{\mathcal{U}}$. The Jacobian resulting from a gauge transformation on $D\mathcal{U}$ cancels against the corresponding quantity for $D\overline{\mathcal{U}}$.

The appearance of the unit matrix in Eq. 3.1 corresponds to the U(1) component of the scalar fields taking on a vacuum expectation value. While this is indeed a classical vacuum state it is not unique. There are infinitely many flat directions in the theory and a priori it is not clear that the vacuum needed to generate the kinetic terms is picked out. To stabilize this vacuum we include in the lattice action (with no sum over repeated indices)

$$\delta S_1 = \frac{N}{2\lambda} \mu^2 \sum_{n, c} a^4 \left(\frac{1}{N} \operatorname{Tr} \left[\mathcal{U}_c(n) \overline{\mathcal{U}}_c(n) \right] - 1 \right)^2.$$
(3.2)

This term also lifts the SU(N) flat directions, biasing the system towards the continuum superconformal phase in which the scalars do not pick up any non-trivial expectation values.

Although δS_1 breaks supersymmetry, it does so softly: the exact Q supersymmetry at $\mu = 0$ guarantees that any Q-breaking counterterms must possess coefficients that vanish as $\mu \to 0$. This can be confirmed by considering the bosonic action per lattice site, $\langle s_B \rangle$. This quantity is related to

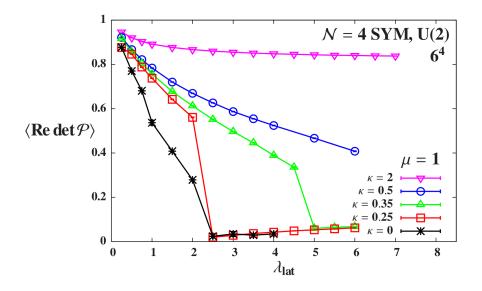


Figure 2: The real part of the plaquette determinant vs. the 't Hooft coupling λ_{lat} on 6⁴ lattices with a variety of κ in Eq. 3.3 and fixed $\mu = 1$. As κ increases the confinement transition associated with the U(1) sector moves to larger λ_{lat} , and disappears entirely for $\kappa \ge 0.5$. Lines connect points with the same κ to guide the eye. From Ref. [12].

(3.4)

a Q Ward identity, which for unbroken Q supersymmetry predicts the exact λ -independent value $s_B = 9N^2/2$ for gauge group U(N). In Fig. 1 we plot $\langle s_B \rangle / 18$ (for N = 2) vs. the 't Hooft coupling for several nonzero values of μ . Clearly as $\mu \to 0$ we approach the supersymmetric value for all 't Hooft couplings.

We mentioned above the issue of lattice artifacts linked to the undesired U(1) sector. The determinant of the plaquette is a gauge-invariant quantity associated with this sector. For sufficiently large $\lambda_{\text{lat}} \approx 2.5$, in Ref. [12] we observed (Re det \mathcal{P}) falling to zero as shown by the black bursts in Fig. 2. At the same λ_{lat} the system confined, with the Polyakov loop also vanishing while the fermion operator developed a large number of near-zero eigenvalues. While Fig. 2 considers N = 2, we have also observed all the same effects with gauge group U(1).

This behavior suggests that we are observing confinement in the compact lattice U(1) sector, which has a well-known dual description in terms of monopole world lines, one-dimensional objects that form closed loops of monopole flux. Indeed, if we measure the density of monopole world lines we find it becomes nonzero at precisely the same $\lambda_{\text{lat}} \approx 2.5$ [12]. To suppress this lattice artifact we penalize small plaquette determinant values by adding to the lattice action

$$\delta S_2 = \kappa \sum_{n, a < b} a^4 |\det \mathcal{P}_{ab}(n) - 1|^2,$$
(3.3)

where \mathcal{P}_{ab} is the plaquette in the *a*-*b* plane. To leading order in the lattice spacing, δS_2 just generates a U(1) field strength term

$$\delta S_2 = 2\kappa \sum_{n, a < b} a^4 \left[1 - \cos \mathcal{F}^0_{ab}(n) \right] + \mathcal{O}(a^5).$$

In agreement with simulations of compact U(1) gauge theory [21], we observe that for $\kappa \ge 0.5$ we suppress monopoles and avoid the confinement transition for arbitrarily large 't Hooft coupling, as shown in Fig. 2.

The lattice action we use in numerical computations includes all of Eqs. 1.3, 3.2 and 3.3: $S = S_{exact} + S_{closed} + \delta S_1 + \delta S_2.$ (6)

$$\frac{(s_{R})-18}{18} = 0.1$$

$$\frac{(s_{R})-18}{0.1} = 0.2$$

$$\frac{(s_{R})-18}{0.1} = 0.2$$

$$\frac{(\mu, \kappa) = (0.2, 0.6)$$

Figure 3: Relative deviations of the bosonic action from its exact supersymmetric value. Left (from Ref. [12]): We fix $\lambda_{\text{lat}} = 1$ and plot vs. κ for 4⁴ lattices with several nonzero μ . The lines are cubic $\kappa \to 0$ extrapolations that confirm the restoration of supersymmetry in the limit (μ, κ) \rightarrow (0,0). Right: We fix (μ, κ) = (0.2, 0.6) and plot vs. λ_{lat} for 8³ × 24 lattices with gauge groups U(2) and U(3), finding ~10% deviations that decrease $\propto 1/N^2$. The lines connect points to guide the eye.

Both δS_1 and δS_2 softly break the Q supersymmetry. As in Fig. 1, the left panel of Fig. 3 shows that the bosonic action s_B acquires its exact supersymmetric value in the limit $(\mu, \kappa) \rightarrow (0, 0)$, indicating the recovery of supersymmetry. This plot also makes it clear that nonzero κ leads to much more severe supersymmetry breaking than does nonzero μ . However, this effect is confined to the U(1) sector that decouples in the continuum limit. The right panel of Fig. 3 illustrates the ~10% level of Q supersymmetry breaking on representative $8^3 \times 24$ lattice ensembles that we will discuss further below. Here we compare N = 2 and 3, finding that violations of supersymmetry decrease $\propto 1/N^2$. In Ref. [12] we discuss the breaking and restoration of Q in more detail.

4. Restoration of the other 15 supersymmetries Q_a and Q_{ab}

Although the lattice formulation discussed above exactly preserves the Q supersymmetry for $(\mu, \kappa) \rightarrow (0, 0)$, the other fifteen Q_a and Q_{ab} are broken, and must be recovered in the continuum limit. In Ref. [11] we showed how restoration of the full symmetries of $\mathcal{N} = 4$ SYM follows from preservation of both Q and any one of a set of discrete R symmetries $\{R_a, R_{ab}\}$, subgroups of the continuum SO(6)_R symmetry that fix the correct coefficients in the long-distance effective action, Eq. 2.5. On the lattice, the R_a acts on the links as [12]

$$R_a \mathcal{U}_a = \mathcal{U}_a \qquad R_a \overline{\mathcal{U}}_a = \overline{\mathcal{U}}_a \qquad R_a \mathcal{U}_b = \overline{\mathcal{U}}_b^{-1} \qquad R_a \overline{\mathcal{U}}_b = \mathcal{U}_b^{-1}, \qquad (4.1)$$

for all $b \neq a$. This transformation commutes with lattice gauge invariance, allowing us to measure violations of the R_a symmetry by considering its action on $m \times n$ Wilson loops in the a-b plane,

$$\mathcal{W}_{ab} = \operatorname{Tr}\left[\prod_{m} \mathcal{U}_{a}(x) \prod_{n} \mathcal{U}_{b}(x+m\widehat{\mu}_{a}) \prod_{m} \overline{\mathcal{U}}_{a}(x+n\widehat{\mu}_{b}) \prod_{n} \overline{\mathcal{U}}_{b}(x)\right]$$

$$\implies R_{a}\mathcal{W}_{ab} = \operatorname{Tr}\left[\prod_{m} \mathcal{U}_{a}(x) \prod_{n} \overline{\mathcal{U}}_{b}^{-1}(x+m\widehat{\mu}_{a}) \prod_{m} \overline{\mathcal{U}}_{a}(x+n\widehat{\mu}_{b}) \prod_{n} \mathcal{U}_{b}^{-1}(x)\right] \equiv \widetilde{\mathcal{W}}_{ab}.$$

$$(4.2)$$

Since our links are non-unitary, $\widetilde{\mathcal{W}}_{ab} \neq \mathcal{W}_{ab}$ even though they follow the same path in the lattice. By computing the relative difference $(\widetilde{\mathcal{W}} - \mathcal{W})/\frac{1}{2}(\widetilde{\mathcal{W}} + \mathcal{W})$ we can assess how badly R_a is broken on the lattice and monitor its restoration as we approach the continuum limit.

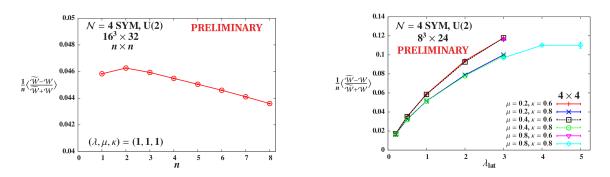


Figure 4: Violations of R_a symmetry, with lines connecting points to guide the eye. Left: For $n \times n$ loops on $16^3 \times 32$ lattices with $(\lambda_{\text{lat}}, \mu, \kappa) = (1, 1, 1)$, the violations decrease (slightly) at larger distance scales, as expected. Right: For 4×4 loops on $8^3 \times 24$ lattices with various (μ, κ) , we find 5–10% violations that vanish in the $\lambda_{\text{lat}} \rightarrow 0$ limit as desired.

We present some results for these R_a symmetry violations in Fig. 4. A complication we encountered is that inverting the non-unitary links changes their gauge-invariant magnitudes,

$$\operatorname{Tr}\left[\overline{\mathcal{U}}_{a}^{-1}\mathcal{U}_{a}^{-1}\right] > \operatorname{Tr}\left[\mathcal{U}_{a}\overline{\mathcal{U}}_{a}\right].$$
(4.3)

This causes the relative differences to be unreasonably sensitive to the number (2n) of inverted links in the $m \times n$ modified Wilson loops \widetilde{W} . The question of how best to handle this artifact is not yet settled. We suspect that projecting W and \widetilde{W} from U(N) to SU(N), by dividing out the determinant of each loop, will prove beneficial. Such a determinant-divided analysis is not yet complete, so for the time being we empirically note that simply normalizing the relative differences by $\frac{1}{2n}$ stabilizes the results, leaving the expected decrease in violations as we approach the $1/L \to 0$ continuum limit. This is shown in the left panel of Fig. 4, where we can see that the resulting violations based on square $n \times n$ loops decrease only gradually on larger distance scales, falling by just ~5% as nincreases from 2 to 8 on $16^3 \times 32$ lattices. This $\frac{1}{2n}$ normalization is not well motivated, and is likely to be replaced by a more robust prescription in the future.

The right panel of Fig. 4 considers R_a symmetry breaking based on 4×4 loops for our main $8^3 \times 24$ U(2) lattice ensembles, as functions of the 't Hooft coupling. These systems exhibit 5–10% violations that vanish in the $\lambda_{\text{lat}} \rightarrow 0$ limit, as desired. There is no visible sensitivity to μ , but larger values of the U(1)-suppressing κ parameter noticeably reduce R_a symmetry breaking, which motivates the determinant projection to SU(N) discussed above. These results are similar to those based on the 1×1 plaquette that we presented in Ref. [12].

Given this observable sensitive to the restoration of Q_a and Q_{ab} , we can revisit the fine-tuning discussed in Section 2. Our goal is to choose the parameter c_2 in Eq. 2.11 in order to minimize R_a symmetry breaking and thereby approach the continuum theory as rapidly and smoothly as possible. Fig. 5 presents an initial study of this tuning on small $4^3 \times 12$ and $6^3 \times 18$ lattices with

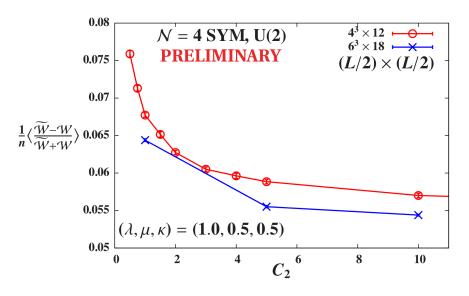


Figure 5: Study of tuning c_2 in Eq. 2.11. We consider R_a symmetry breaking in 2×2 loops on $4^3 \times 12$ lattices and 3×3 loops on $6^3 \times 18$ lattices, all with $(\lambda_{\text{lat}}, \mu, \kappa) = (1, 0.5, 0.5)$. We observe gradual improvement as c_2 increases above its classical value $c_2 = 1$, while violations increase steeply for smaller $c_2 < 1$. Lines connect points with the same volume to guide the eye.

fixed $(\lambda_{\text{lat}}, \mu, \kappa) = (1, 0.5, 0.5)$. As desired, violations decrease for larger 3×3 Wilson loops on the larger volume, though again the approach to the $1/L \rightarrow 0$ continuum limit is gradual. In addition we see that values of c_2 larger than the classical $c_2 = 1$ lead to ~10% smaller violations of R_a symmetry, though on these volumes there is no strongly preferred value; the results flatten out as c_2 increases. For $c_2 < 1$, in contrast, R_a symmetry breaking increases steeply. Conveniently, we also find that Q breaking behaves similarly (not shown), improving gradually for larger c_2 . This is important because restoration of the full symmetries of $\mathcal{N} = 4$ SYM requires both Q and R_a . For the time being we continue to use the classical $c_2 = 1$ in our larger-volume studies. Fig. 5 suggests that larger c_2 would only result in modest improvement of R_a symmetry breaking.

In Fig. 6 we present an initial exploration of MCRG blocking based on Eq. 2.1. We block 12^4 lattices down to 6^4 and measure observables on the blocked lattices as well as on independent 6^4 ensembles with fixed $0.5 \le \lambda_{\text{lat}} \le 1.5$ and $(\mu, \kappa) = (0.5, 0.5)$. As in Ref. [13], we fix $\xi^4 \approx 1.2$ by requiring that the blocked plaquette matches the "bare" value on the 6^4 ensembles. The left panel of Fig. 6 shows that all larger blocked Wilson loops then approximately match the corresponding bare loops, extending the results in Ref. [13]. This evidence that the couplings don't flow noticeably under RG blocking is consistent with our observation of a Coulombic static potential in Section 6. However, the blocked R_a symmetry violations in the right panel of Fig. 6 end up significantly smaller than the bare 6^4 results, by ~25% for these 3×3 loops. This is the behavior we expect as we move towards the $1/L \rightarrow 0$ continuum conformal theory.

At present we have investigated the effects of the R_a transformation only for the gauge links, through Eqs. 4.1 and 4.2. Additional tests of the discrete R symmetries involving fermions should also be carried out in the future, to ensure that symmetry restoration is consistent across all sectors of the theory. Unfortunately, the rescalings that we applied to the fermion fields in Eq. 2.6 interfere with such analyses, and further exploration is required to determine the best approach.

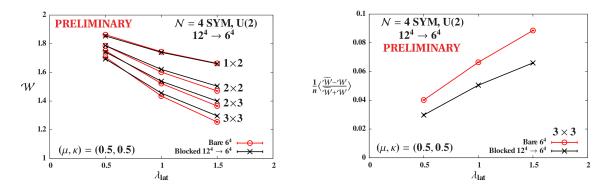


Figure 6: Wilson loops (left) and violations of R_a symmetry (right) after one step of $12^4 \rightarrow 6^4$ MCRG blocking. Lines connect points for loops of the same size to guide the eye. The approximate agreement between the blocked Wilson loops and "bare" measurements on independent 6^4 ensembles with the same couplings indicates that those couplings don't flow significantly under RG blocking. The reduced R_a symmetry breaking is consistent with the blocked measurements probing larger distance scales closer to the $1/L \rightarrow 0$ continuum limit.

5. Measuring the complex Pfaffian to assess a potential sign problem

A worrisome aspect of our numerical calculations is that they must be phase-quenched to take advantage of efficient importance sampling algorithms [14]. Gaussian integration over the fermion fields $(\eta, \psi_a, \chi_{ab})$ in the path integral produces the Pfaffian of the fermion operator \mathcal{D} defined by Eq. 1.3. This Pfaffian is not manifestly real for any given gauge configuration, pf $\mathcal{D} = |pf\mathcal{D}|e^{i\alpha}$, and we retain only its magnitude in the path integral, omitting its phase $e^{i\alpha}$. So long as $\langle e^{i\alpha} \rangle$ is nonzero, bona fide operator expectation values may be determined through phase reweighting. If α fluctuates so much that $\langle e^{i\alpha} \rangle$ is consistent with zero, then simulations would suffer from a sign problem and numerical results could not be trusted.

In Ref. [12] we initiated an ongoing study of the complex Pfaffian of lattice $\mathcal{N} = 4$ SYM. Focusing on gauge group U(2) we found that $e^{i\alpha}$ was approximately real and positive on every investigated system, close enough to unity that phase reweighting is not necessary to obtain reliable expectation values. Fluctuations in the phase did not grow with the system size on the largest accessible volumes up to $4^3 \times 6$, and on $2^3 \times 4$ lattices $\langle \cos \alpha \rangle$ showed little dependence on the number of colors N = 2, 3 or 4. These results are included in Fig. 7, which also considers additional U(3) and U(4) lattice volumes, finding similar behavior for all gauge groups. All points in this figure are for ($\lambda_{\text{lat}}, \mu, \kappa$) = (1,1,1). In initial investigations we have observed even better behavior for smaller λ_{lat} and no significant dependence on κ , so long as κ is large enough to suppress the U(1) confinement transition discussed in Section 3.

Evaluating the Pfaffian is a notoriously hard problem, and these results are only obtainable thanks to new parallel software described in detail by Ref. [14]. Despite this new code, we remain limited to lattice volumes around $4^3 \times 6$ (and smaller for larger N), where several days are already required for each measurement. However, now that we have designed and implemented the RG blocking transformation in Eq. 2.1, it will be interesting to measure the Pfaffian on lattices blocked

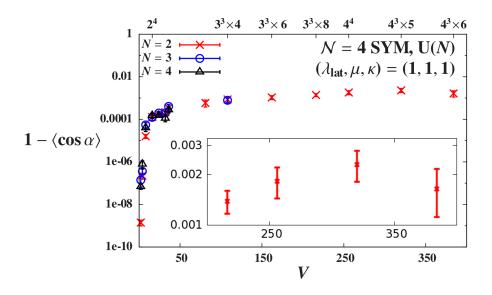


Figure 7: Semi-log plot of $1 - \langle \cos \alpha \rangle$ vs. lattice volume, where α is the phase of the pfaffian. All points are for $(\lambda_{\text{lat}}, \mu, \kappa) = (1, 1, 1)$, for gauge groups U(2), U(3) and U(4). The inset zooms in on the four largest-volume U(2) results with $V \ge 3^3 \times 8$, where the small phase does not grow with volume.

down to a small number of sites. In addition to searching for information about the phase on our larger-volume lattice ensembles, by comparing systems after different numbers of RG blocking steps we may gain insight into the behavior of the Pfaffian in the continuum limit.

6. Static potentials and Coulomb coefficients

In previous sections we focused on validating our lattice system and phase-quenched calculations, to ensure that we simulate $\mathcal{N} = 4$ SYM to a good approximation and can recover the appropriate theory in the continuum limit. This work is necessary to establish that numerical results from lattice $\mathcal{N} = 4$ SYM can be trusted. Let us now consider a more physical quantity, the static potential. In Ref. [12] we presented results for the U(2) potential, finding Coulombic behavior V(r) = A - C/r at both weak and strong 't Hooft coupling, with the Coulomb coefficients Cin good agreement with leading-order perturbation theory. Here we supplement these results with a first look at the U(3) static potential.

We extract the static potential from Wilson loops $W(\vec{r}, t)$ measured on $8^3 \times 24$ lattices. In Section 4 we considered $m \times n$ Wilson loops oriented along the principal axes of the lattice. For the static potential we want to be more general and consider all possible spatial separations \vec{r} , which we do by gauge fixing to Coulomb gauge and computing

$$W(\vec{r},t) = \text{Tr}\left[P(\vec{x},t,t_0)P^{\dagger}(\vec{x}+\vec{r},t,t_0)\right].$$
(6.1)

Here $P(\vec{x}, t, t_0)$ is a product of temporal links U_t at spatial location \vec{x} , extending from timeslice t_0 to timeslice $t_0 + t$. We extract V(r) by fitting $W(r, t) = w \exp(-V(r)t)$, combining all \vec{r} with the same magnitude $r \equiv |\vec{r}|$ and comparing fit ranges $t_{min} \leq t < N_t/2 = 12$ to find the $t_{min} = 5-6$ above which the results form stable plateaus.

To directly compare our data and results with continuum expectations, we must take care to account for the physical structure of the A_4^* lattice. As discussed in Ref. [12] we use a convenient representation of the A_4^* lattice that adds the fifth link $\hat{\mu}_4 = (-1, -1, -1, -1)$ to the usual four hypercubic basis vectors. This allows us to use familiar integer displacement vectors in the code,

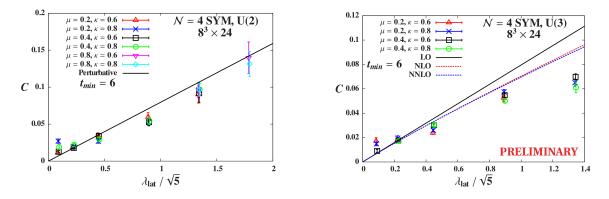


Figure 8: Static potential Coulomb coefficients C for the usual Wilson loops from $8^3 \times 24$ lattices with gauge groups U(2) (left, from Ref. [12]) and U(3) (right). The N = 2 results are consistent with the leading-order perturbative prediction. The N = 3 results fall below even the NNLO perturbative prediction from Refs. [22, 23].

which must be converted to physical distances on the true A_4^* lattice. Doing so, we find that displacements corresponding to any r > 2.6 can wrap around the L = 8 spatial lattice, making them unusable in the gauge-fixed correlator Eq. 6.1. Limited investigations of L = 12 and 16 suggest that our signals disappear into the background noise for $r \gtrsim 3$ in any case, so that we don't lose much due to the restriction $\sqrt{3/4} \le r \le 2.6$ on our $8^3 \times 24$ lattices.

An additional consequence of the A_4^* lattice structure is the normalization factor relating the continuum and lattice couplings, $\lambda = \lambda_{\text{lat}}/\sqrt{5}$. This is simply the Jacobian of the transformation between the lattice and continuum space-time coordinates, which is non-trivial because the A_4^* links are not orthogonal. We provide a more intuitive derivation of this rescaling in Ref. [12].

Fitting V(r) to the Coulomb form V(r) = A - C/r, using all accessible $\sqrt{3/4} \le r \le 2.6$, produces the Coulomb coefficient results shown in Fig. 8. For each gauge group U(2) and U(3) we consider several values of (μ, κ) , observing no significant dependence on these parameters. We plot the Coulomb coefficients as functions of the continuum 't Hooft coupling $\lambda = \lambda_{\text{lat}}/\sqrt{5}$, and compare our numerical results with perturbative predictions. For N = 2 in the left panel of Fig. 8 we find values of C in good agreement with leading-order (LO) perturbation theory,

$$C_{LO} = \frac{\lambda_{\text{lat}}/\sqrt{5}}{4\pi} = \frac{\lambda}{4\pi}.$$
(6.2)

For N = 3 in the right panel, however, our results begin deviating significantly from the LO line at stronger couplings. Although the next-to-leading-order and next-to-next-to-leading-order corrections from Refs. [22, 23] move the perturbative prediction closer to our results, a clear difference remains. These NLO and NNLO contributions are

$$\frac{C_{NLO}}{C_{LO}} = \frac{\lambda}{2\pi^2} \left(\log\left[\frac{\lambda}{2\pi}\right] + \gamma_E - 1 \right)$$

$$\frac{C_{NNLO}}{C_{LO}} = \frac{\lambda^2}{8\pi^4} \left\{ \log\left[\frac{\lambda}{2\pi}\right] + \gamma_E + \left(1 + \frac{\pi^2}{3}\right) \left(\log\left[\frac{\lambda}{2\pi}\right] + \gamma_E\right) - \frac{\pi^2}{12} - \frac{7}{2} + \frac{9}{4}\zeta(3) \right\}$$
(6.3)

where γ_E is Euler's constant, $\zeta(3) \approx 1.202$ is Apéry's constant, and we have set $C_A \alpha = 2C_F \alpha = \frac{\lambda}{4\pi}$ for gauge group U(N).

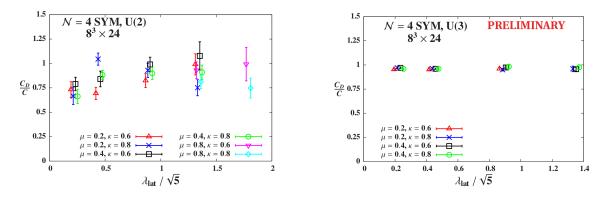


Figure 9: Ratios C_D/C of static potential Coulomb coefficients for the determinant-divided Wilson loops relative to those for the usual loops, from $8^3 \times 24$ lattices with gauge groups U(2) (left, from Ref. [12]) and U(3) (right). Both plots show the same vertical range to illustrate that the new N = 3 results are much less noisy than N = 2, though the former are consistently larger than the expected ratio $1 - \frac{1}{N^2} = \frac{8}{9}$.

While it is tempting to interpret the right panel of Fig. 8 as a suggestion that we are seeing C depart from perturbation theory to approach the famous large-N prediction $C \propto \sqrt{\lambda}$ at strong coupling [24, 25], we suspect that both N = 3 and $\lambda \leq 3/\sqrt{5}$ are still too small for this to happen. We can see that the NLO and (especially) NNLO corrections are still small compared to C_{LO} , suggesting that our results should remain within the perturbative regime. Although the contrast between C for U(2) and U(3) in the two plots of Fig. 8 is striking, its cause is not yet clear. It will be very interesting to add U(4) results to this picture, to clarify any systematic trends. This work is underway, but is challenging due to the steeply increasing computation costs, which grow $\propto N^5$ [14].

Another interesting difference between the U(2) and U(3) static potentials is visible when we divide out the determinant of the Wilson loops from which V(r) is extracted. As discussed in Section 4, this procedure should approximately project the observables to SU(N), removing the U(1) sector that decouples in the continuum limit. We expect that the Coulomb coefficients C_D resulting from determinant-divided Wilson loops will be reduced by a factor $(N^2 - 1)/N^2$ compared to the usual C in Fig. 8. (This holds in perturbation theory through NLO, and the departure from this simple scaling at NNLO [23] appears negligible.)

In Fig. 9 we test our expectation by plotting the ratios C_D/C , again as functions of the continuum $\lambda = \lambda_{\text{lat}}/\sqrt{5}$ for several values of (μ, κ) . As before, there is no clear dependence on (μ, κ) , and the ratios are also insensitive to the coupling. Our U(2) results are fairly noisy, but appear consistent with the expected ratio of 3/4. In contrast, our results for N = 3 are much more stable, which we emphasize by showing the same vertical range in both plots of Fig. 9. Numerically we find $C_D/C \approx 0.96$ with uncertainties small enough to clearly differ from the expected ratio of 8/9. Again, ongoing investigations of U(4) will be important to clarify the interpretation of these initial U(3) results.

Finally, we can also consider the $\mathcal{N} = 4$ static potential with the contribution of the scalar fields removed [22, 23]. We implement this by building Wilson loops from unitary matrices u defined through the polar decomposition of each link variable, $\mathcal{U} = uH$ where H is Hermitian and positive definite. Decoupling the scalar fields simplifies the perturbative predictions for the

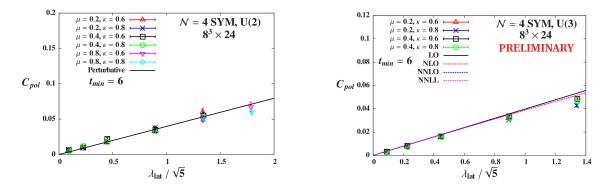


Figure 10: Static potential Coulomb coefficients C_{pol} for the polar-projected Wilson loops that decouple the scalar fields, from $8^3 \times 24$ lattices with gauge groups U(2) (left, from Ref. [12]) and U(3) (right). The N = 2 results are consistent with the leading-order perturbative prediction, and the N = 3 results also remain close to perturbation theory.

Coulomb coefficient in Refs. [22, 23], with Eqs. 6.2 and 6.3 becoming

$$C_{pol} = \frac{\lambda}{8\pi} \left\{ 1 - \frac{\lambda}{4\pi^2} + A\left(\frac{\lambda}{4\pi^2}\right)^2 + B\left(\frac{\lambda}{4\pi^2}\right)^3 \log\left[\frac{\lambda}{4\pi}\right] + \mathcal{O}\left(\lambda^3\right) \right\}$$

$$A = \frac{5}{2} + \frac{\pi^2}{4} - \frac{\pi^4}{64} \qquad B = \frac{1}{12\pi}.$$
(6.4)

At leading order the polar-projected C_{pol} is just half the usual Coulomb coefficient, $C_{pol}/C = 1/2$.

In Fig. 10 we plot C_{pol} vs. $\lambda_{lat}/\sqrt{5}$ for N = 2 and 3, using the same axes and considering the same values of (μ, κ) as in Fig. 8. The U(2) results for C_{pol} in the left panel of Fig. 10 are in good agreement with leading-order perturbation theory, and appear less noisy than the usual Coulomb coefficients in Fig. 8. The U(3) results in the right panel of Fig. 10 are significantly closer to leading-order perturbation theory than the corresponding C in Fig. 8. At the same time, the higher-order perturbative corrections have barely visible effects in this range of λ_{lat} . The $N = 3 C_{pol}$ may still begin to drop below the perturbative predictions at the strongest coupling we have investigated so far, but this deviation is not significant and may be cured by extrapolating $(\mu, \kappa) \rightarrow (0, 0)$.

7. Recapitulation of status, prospects and next steps

Our investigations of lattice $\mathcal{N} = 4$ SYM are making progress addressing a number of interesting questions, regarding both the validity of our lattice calculations as well as the resulting predictions that may be compared to continuum perturbation theory and gauge–gravity duality. In this proceedings, we presented several new developments that go beyond the results appearing in Refs. [12, 13]. These include an initial study of fine-tuning to recover the full symmetries of $\mathcal{N} = 4$ SYM in the continuum limit, which makes use of the recent derivation that only a single parameter needs to be tuned so long as the moduli space is not lifted by nonperturbative effects. We also extended our measurements of the complex Pfaffian, considering more lattice volumes for larger N = 3 and 4, which also indicate that the lattice theory does not suffer from a sign problem. These results provide additional evidence that our phase-quenched lattice calculations simulate $\mathcal{N} = 4$ SYM to a good approximation, and can recover the appropriate theory in the continuum limit.

We then focused on the static potential, extending our previous study of gauge group U(2) with initial results for N = 3. In both cases, we find Coulombic behavior at both weak and strong coupling, and compare the Coulomb coefficients to perturbation theory. Our initial U(3) results appear significantly less noisy than those for N = 2, but fall away from perturbative predictions at the strongest couplings we have investigated. When we consider Wilson loops approximately projected to SU(3), we find Coulomb coefficients C_D closer to the usual U(3) C than we expected from naive $(N^2 - 1)/N^2$ scaling. It is not yet clear how we should interpret the current contrast between N = 2 and 3. We are working on extending these investigations to N = 4, which we hope will clarify any systematic trends.

Additional future work will include further studies of the Pfaffian phase, with the aim of improving our qualitative understanding of why we see no sign problem for lattice $\mathcal{N} = 4$ SYM. Although it is extremely expensive to measure the Pfaffian on larger lattice volumes or for larger values of N, we may be able to gain insight by applying our new RG blocking scheme to obtain blocked systems with fewer degrees of freedom, which are more practical to analyze. MCRG anal-

yses may also prove useful to predict the anomalous dimensions of single-trace operators like the Konishi, a particularly exciting goal. Such predictions would be complementary to those obtainable in perturbation theory [26] or via the conformal bootstrap program [27].

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