

Hamiltonian domain wall fermions at strong coupling

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Abstract

We apply strong-coupling perturbation theory to gauge theories containing domain-wall fermions in Shamir's surface version. We construct the effective Hamiltonian for the color-singlet degrees of freedom that constitute the low-lying spectrum at strong coupling. We show that the effective theory is identical to that derived from naive, doubled fermions with a mass term, and hence that domain-wall fermions at strong coupling suffer both doubling and explicit breaking of chiral symmetry. Since we employ a continuous fifth dimension whose extent tends to infinity, our result applies to overlap fermions as well.

I. INTRODUCTION

For a quarter of a century, users of lattice fermions have been forced to choose between the Scylla of chiral symmetry violation [1] and the Charybdis of uncontrolled flavor doubling [2]. The constraints imposed by the Nielsen–Ninomiya theorem [3] imply that attempts to circumvent these problems must avoid its assumptions. An avenue pursued in recent years is that of employing a large number (tending to infinity) of fermion flavors, one of which survives in the continuum limit as a massless fermion supporting a chiral symmetry group. One route down this avenue is that of Kaplan [4], who introduced a fifth dimension with its concomitant excitations, one of which is kept at low energy by a defect in a background field. We here focus on Shamir’s variant [5] of domain wall fermions, sometimes known as surface fermions. The much-studied overlap fermions [6] are equivalent to this formulation in the limit that the size and lattice spacing of the fifth dimension are taken to infinity and zero, respectively [7, 8].

The domain wall formulation has been shown to yield any desired number of undoubled, massless fermions with full chiral symmetry—when the fermions are free. Shamir [5] proved that exact chiral symmetry—as evidenced by multiplicative mass renormalization—survives to all orders in perturbation theory when the fermions are coupled to gauge fields. A *non-perturbative* proof of the axial Ward identities was given by Furman and Shamir [9]. The suppression of symmetry-breaking terms requires that the length of the fifth dimension be taken to infinity. Monte Carlo calculations on finite lattices [10], however, indicate a pattern of increasing violation of chiral symmetry as the coupling is made stronger. This leads us to seek analytical knowledge of the behavior of domain wall fermions at strong coupling.

In this paper we develop a Hamiltonian formalism for domain wall fermions and analyze it in the limit of strong gauge coupling. Strong coupling analysis is natural in lattice gauge theory and, indeed, was seen immediately as one of the advantages of a non-perturbative cutoff procedure [11]. In the Hamiltonian version, strong-coupling analysis takes the form of Rayleigh-Schrödinger perturbation theory. It was first applied [12] to fermions in the newly-invented Hamiltonian lattice gauge theory in its original formulation, which employed what came to be known as staggered fermions [2]. The method was subsequently applied to Wilson fermions by Smit [13] and to naive and long-range [14] fermions by the SLAC group [15] (see also [16]).

Degenerate perturbation theory leads to an effective Hamiltonian H^{eff} that acts among the low-lying, fluxless states. The advantage of the Hamiltonian formulation in this regard is that much may be learned from the effective Hamiltonian even if one makes no attempt to diagonalize it. This is in contrast with Euclidean strong coupling methods, where some additional approximation such as a hopping parameter expansion, mean field theory, or a large- N_c limit must be applied. We note a recent study [17] of overlap fermions in the Euclidean strong-coupling theory with a hopping parameter expansion, which does not uncover the features we present below.

In all cases, H^{eff} takes the form of an antiferromagnet possessing the global symmetry of the underlying fermion formulation. For Wilson fermions, there is no symmetry in the single-flavor case; for Kogut-Susskind fermions, there is at least a discrete chiral-spin-flavor symmetry. For naive and SLAC fermions, the symmetry is continuous and contains axial generators, which makes it unfortunate that these theories must be rejected on the grounds of doubling and non-locality [18], respectively. Nevertheless, it is the analysis of these theories in [15] that will guide us through our work on domain wall fermions.

Domain wall fermions, like the Wilson fermions from which they are constructed, begin with no axial symmetry at all. When the size L of the fifth dimension is taken to infinity, its low-lying modes become topological zero-modes and turn into the chiral components of a massless Dirac fermion. Thus one may expect that H^{eff} will possess no axial symmetries until $L \rightarrow \infty$, and then the desired axial symmetry will be restored. Surprisingly, we find that in the $L \rightarrow \infty$ limit H^{eff} has *too much* symmetry. The effective theory is that of an antiferromagnet in a staggered magnetic field. In calculating the site-site couplings in the Hamiltonian, the Wilson term quickly drops out and leaves us with the naive nearest-neighbor terms of the original Hamiltonian only. Thus the site-site term in H^{eff} is identical to that stemming from a naive fermion prescription, with all its doubling and accidental symmetries [15]. The staggered magnetic field comes from coupling the zero modes to the higher modes of the fifth dimension. It is identical in form to the term generated by a mass term in the naive theory. This means that whatever the interpretation of the accidental symmetries of the nearest-neighbor theory, the symmetries that survive the magnetic field term are not axial. This is confirmed by adding Shamir's mass term [5] as a perturbation, and seeing that it gives a contribution to H^{eff} of exactly the same form. Thus our result is that in the strong-coupling limit, domain wall fermions are massive and doubled.

We note that the proof of Furman and Shamir [9] (that chiral symmetry is exact at finite gauge coupling) depends on isolating background gauge field configurations that induce zero modes in the fifth-dimension transfer matrix. This can be done when the four-dimensional Euclidean lattice is finite, but the argument could well break down in the time-continuum limit necessary for construction of the Hamiltonian.

In Sec. 2 we write out the (4+1)-dimensional Hamiltonian of domain wall fermions coupled to a (3+1)-dimensional gauge field, following the commonly used formulation of the corresponding Euclidean theory. For simplicity we work with an Abelian theory with a single flavor; we review the free theory in Appendix A. We begin a perturbation expansion in strong coupling and are thus led to solve the single-site problem, a massive fermion in (1+1) continuous dimensions on an interval $[0, L]$, in Sec. 3 and in Appendix B. The zero modes of this single-site problem, which appear only for $M > 3$, are the low-lying states that are coupled by the three-dimensional hopping Hamiltonian. Doing second-order perturbation theory in the latter gives, in Sec. 4, the effective hopping Hamiltonian H^{eff} governing the low-lying, colorless degrees of freedom. We show in Sec. 5 that the effective Hamiltonian possesses accidental internal symmetries identical to those of one derived from naive, nearest-neighbor fermions with non-zero mass. This means that domain wall fermions, as well as overlap fermions which are equivalent to them in the $L \rightarrow \infty$ limit, are massive and doubled in the strong coupling limit. We confirm in Sec. 6 our identification of the mass term by showing that Shamir's mass term generates a term of the same form in perturbation theory. Finally, we present in Sec. 7 the straightforward extension of our results to a theory with more flavors, whence we find that the mass term generated indeed breaks all axial symmetries, including those that should not be afflicted with anomalies. We close with some comments in Sec. 8, chief of which is the observation that the symmetry properties of H^{eff} are unchanged by promotion of the gauge symmetry to a non-Abelian group. Thus our result of the failure of domain wall fermions at strong coupling is quite general, and calls for careful study of their phase structure at finite gauge coupling.

II. THE HAMILTONIAN

We consider an Abelian gauge field coupled to lattice fermions defined via Shamir's domain wall prescription. Space-time consists of a three-dimensional spatial lattice with

continuous time and a continuous fifth dimension specified by the coordinate $s \in [0, L]$. (For a reprise of free fermions in this formulation, see Appendix A.) The Hamiltonian is

$$H = H_E + H_B + H_4 + H_5 + H_m, \quad (2.1)$$

with

$$H_E = \frac{1}{2}g^2 \sum_{\mathbf{x}i} E_{\mathbf{x}i}^2 \quad (2.2)$$

$$H_B = -\frac{1}{g^2} \sum_{\mathbf{x}i} \cos B_{\mathbf{x}i} \quad (2.3)$$

$$H_4 = \sum_{\mathbf{x}} \int ds \psi_{\mathbf{x}s}^\dagger \beta \sum_i \left(\frac{(1 - i\gamma_i)}{2} U_{\mathbf{x}i} \psi_{\mathbf{x}+i,s} + \frac{(1 + i\gamma_i)}{2} U_{\mathbf{x}-i,i}^\dagger \psi_{\mathbf{x}-i,s} \right) \quad (2.4)$$

$$H_5 = \sum_{\mathbf{x}} \int ds \left[\psi_{\mathbf{x}s}^\dagger (-i\alpha_5 \partial_5) \psi_{\mathbf{x}s} + (M - 3) \psi_{\mathbf{x}s}^\dagger \beta \psi_{\mathbf{x}s} \right] \quad (2.5)$$

$$H_m = m \sum_{\mathbf{x}} \left(\psi_{\mathbf{x}L}^\dagger \beta \frac{(1 + \gamma_5)}{2} \psi_{\mathbf{x}0} + \psi_{\mathbf{x}0}^\dagger \beta \frac{(1 - \gamma_5)}{2} \psi_{\mathbf{x}L} \right). \quad (2.6)$$

H_4 is the ordinary Wilson Hamiltonian for hopping on the three-dimensional spatial lattice, while H_5 is a massive Dirac Hamiltonian in the new fifth dimension. The gauge field couples equally to all fields at the same lattice site \mathbf{x} , irrespective of s , and there is no fifth component of the gauge field. H_m is Shamir's mass term which couples the layers at $s = 0$ and $s = L$. As discussed in Appendix A, we have set $r = -1$; as shown in Appendix B, we must choose $M - 3 > 0$ in order to obtain surface modes in the fifth dimension.

When $g^2 \gg 1$, we diagonalize H_E . The ground state is highly degenerate, consisting of all states with $\mathbf{E} = 0$. Gauss' Law restricts us to states $|\varphi\rangle$ satisfying

$$\left[(\nabla \cdot \mathbf{E})_{\mathbf{x}} - \int ds \psi_{\mathbf{x}s}^\dagger \psi_{\mathbf{x}s} \right] |\varphi\rangle = 0 \quad (2.7)$$

for all \mathbf{x} , so the fermion states must satisfy

$$\int ds \psi_{\mathbf{x}s}^\dagger \psi_{\mathbf{x}s} |\varphi\rangle = 0 \quad (2.8)$$

for every \mathbf{x} . The degeneracy is lifted in zeroth order by H_5 , which constitutes the single-site problem. Then H_4 , which creates and annihilates links of flux, gives a contribution in second-order perturbation theory that is of order $1/g^2$ because of the energy denominator. We take m to be small enough that H_m comes next in perturbation theory. Finally, H_B acts in second-order perturbation theory to give contributions of order $1/g^6$, which we will neglect entirely.

III. THE SINGLE-SITE PROBLEM

H_5 is a sum of terms pertaining to single lattice sites,

$$H_5 = \sum_{\mathbf{x}} \psi_{\mathbf{x}}^\dagger h \psi_{\mathbf{x}}. \quad (3.1)$$

In Appendix B, we find eigenmodes of the differential operator h satisfying

$$h u_\lambda(s) = \lambda u_\lambda(s). \quad (3.2)$$

For each eigenvector u_λ with $\lambda > 0$, we have an eigenvector $v_\lambda \equiv u_{-\lambda} = -\gamma_5 u_\lambda$. (This is because $\gamma_5 h + h \gamma_5 = 0$.) There is also a spin index $\sigma = \pm$ for each λ . Dropping the \mathbf{x} index, we write an eigenfunction expansion of ψ_s ,

$$\psi_s = \sum_{\lambda > 0, \sigma} \left(\chi_{\lambda\sigma}^+ u_{\lambda\sigma}(s) + \chi_{\lambda\sigma}^- v_{\lambda\sigma}(s) \right). \quad (3.3)$$

(The notation $\lambda > 0$ here includes $\lambda = \lambda_0$, the zero mode in the limit $L \rightarrow 0$.) The single-site Hamiltonian and the charge operator take the form

$$\psi^\dagger h \psi = \sum_{\lambda > 0, \sigma} \lambda \left((\chi_{\lambda\sigma}^+)^\dagger \chi_{\lambda\sigma}^+ - (\chi_{\lambda\sigma}^-)^\dagger \chi_{\lambda\sigma}^- \right) \quad (3.4)$$

$$Q \equiv \sum_s \psi_s^\dagger \psi_s = \sum_{\lambda > 0, \sigma} \left((\chi_{\lambda\sigma}^+)^\dagger \chi_{\lambda\sigma}^+ + (\chi_{\lambda\sigma}^-)^\dagger \chi_{\lambda\sigma}^- \right). \quad (3.5)$$

To obtain the ground state of the single-site Hamiltonian, we must fill the Dirac sea. We do this by writing

$$\chi_{\lambda\sigma}^+ = b_{\lambda\sigma} \quad (3.6)$$

$$\chi_{\lambda\sigma}^- = d_{\lambda\sigma}^\dagger. \quad (3.7)$$

Now

$$\psi^\dagger h \psi = \sum_{\lambda > 0, \sigma} \lambda \left(b_{\lambda\sigma}^\dagger b_{\lambda\sigma} + d_{\lambda\sigma}^\dagger d_{\lambda\sigma} - 1 \right) \quad (3.8)$$

$$Q = \sum_{\lambda > 0, \sigma} \left(b_{\lambda\sigma}^\dagger b_{\lambda\sigma} - d_{\lambda\sigma}^\dagger d_{\lambda\sigma} + 1 \right). \quad (3.9)$$

We can drop the 1's in these formulae: In the first case the 1 represents a constant energy shift, and in the second it is a c -number charge density that can be defined away in Gauss' Law.

The ground state at each site \mathbf{x} is $|0\rangle_{\mathbf{x}}$, the state annihilated by all $b_{\mathbf{x}\lambda\sigma}$ and $d_{\mathbf{x}\lambda\sigma}$. Low-lying states are created by $b_{\mathbf{x}0\sigma}^\dagger$ and $d_{\mathbf{x}0\sigma}^\dagger$, creation operators for the zero mode; [19] these states are degenerate with $|0\rangle_{\mathbf{x}}$ when $L \rightarrow \infty$. States with $\lambda \neq 0$ are separated by the gap $M - 3$, and lie outside the degenerate subspace. Thus there are six neutral degenerate states at each site: $|0\rangle_{\mathbf{x}}$, four single-pair states of the form $b_{\mathbf{x}0\sigma}^\dagger d_{\mathbf{x}0\sigma'}^\dagger |0\rangle_{\mathbf{x}}$, and the two-pair state $b_{\mathbf{x}0+}^\dagger d_{\mathbf{x}0+}^\dagger b_{\mathbf{x}0-}^\dagger d_{\mathbf{x}0-}^\dagger |0\rangle_{\mathbf{x}}$. We denote these states as $|\eta\rangle_{\mathbf{x}}$. A basis for the degenerate ground states of $H_E + H_5$ is then $|\{\eta\}\rangle = \prod_{\mathbf{x}} |\eta_{\mathbf{x}}\rangle_{\mathbf{x}}$.

We now define a field operator that acts only on the zero mode. It is convenient to use operators that create and annihilate particles localized near $s = 0$ or $s = L$ only, since these will be the chiral states. In view of Eq. (B2) and Eqs. (B11) and (B12), these are defined by a Bogoliubov transformation,

$$\begin{aligned} B_{\mathbf{x}\sigma} &= \frac{1}{\sqrt{2}}(b_{\mathbf{x}0\sigma} + d_{\mathbf{x}0\sigma}^\dagger) \\ D_{\mathbf{x}0\sigma}^\dagger &= \frac{1}{\sqrt{2}}(b_{\mathbf{x}0\sigma} - d_{\mathbf{x}0\sigma}^\dagger). \end{aligned} \quad (3.10)$$

Then the field operator is

$$\Psi_{\mathbf{x}} = \begin{pmatrix} B_{\mathbf{x}0+} \\ B_{\mathbf{x}0-} \\ D_{\mathbf{x}0+}^\dagger \\ D_{\mathbf{x}0-}^\dagger \end{pmatrix}. \quad (3.11)$$

IV. THE EFFECTIVE HOPPING HAMILTONIAN

We now apply second-order degenerate perturbation theory in H_4 to define an effective hopping Hamiltonian via

$$\langle \{\eta\} | H^{\text{eff}} | \{\eta'\} \rangle = \sum_{|j\rangle} \frac{\langle \{\eta\} | H_4 | j \rangle \langle j | H_4 | \{\eta'\} \rangle}{E_{\{\eta\}} - E_j}. \quad (4.1)$$

The energy denominator contains $g^2/2$ because of the electric flux, added to any energy due to continuum fermions in the intermediate state.

Let us focus on a term wherein H_4 acts on $|\{\eta'\}\rangle$ by raising the electric flux on the link $\mathbf{x}i$ by a unit. Then H_4 has to act on the intermediate state $|j\rangle$ so as to lower the flux on the same link in order to give a fluxless state of the type $|\{\eta\}\rangle$. The numerator of this term in

Eq. (4.1) is

$$\langle \{\eta\} | \int ds \psi_{\mathbf{x}+\hat{i},s}^\dagger \beta \frac{(1+i\gamma_i)}{2} \psi_{\mathbf{x}s} |j\rangle \langle j| \int ds' \psi_{\mathbf{x}s'}^\dagger \beta \frac{(1-i\gamma_i)}{2} \psi_{\mathbf{x}+\hat{i},s'} | \{\eta'\} \rangle, \quad (4.2)$$

where we have suppressed the gauge field matrix element, which is unity. We isolate the zero-mode in ψ ,

$$\begin{aligned} \psi_{\mathbf{x}s} &= \sum_{\sigma} \left(b_{\mathbf{x}0\sigma} u_{0\sigma}(s) + d_{\mathbf{x}0\sigma}^\dagger v_{0\sigma}(s) \right) + \sum_{\lambda>0,\sigma} \left(b_{\mathbf{x}\lambda\sigma} u_{\lambda\sigma}(s) + d_{\mathbf{x}\lambda\sigma}^\dagger v_{\lambda\sigma}(s) \right) \\ &\equiv \psi_{\mathbf{x}s}^{(0)} + \psi_{\mathbf{x}s}^{(1)}, \end{aligned} \quad (4.3)$$

and expand Eq. (4.2) to

$$\begin{aligned} \langle \{\eta\} | \int ds \left(\psi_{\mathbf{x}+\hat{i},s}^{(0)\dagger} + \psi_{\mathbf{x}+\hat{i},s}^{(1)\dagger} \right) \beta \frac{(1+i\gamma_i)}{2} \left(\psi_{\mathbf{x}s}^{(0)} + \psi_{\mathbf{x}s}^{(1)} \right) |j\rangle \\ \times \langle j| \int ds' \left(\psi_{\mathbf{x}s'}^{(0)\dagger} + \psi_{\mathbf{x}s'}^{(1)\dagger} \right) \beta \frac{(1-i\gamma_i)}{2} \left(\psi_{\mathbf{x}+\hat{i},s'}^{(0)} + \psi_{\mathbf{x}+\hat{i},s'}^{(1)} \right) | \{\eta'\} \rangle \end{aligned} \quad (4.4)$$

The states $|\{\eta\}\rangle$ and $|\{\eta'\}\rangle$ contain no quanta with $\lambda > 0$ at \mathbf{x} or at $\mathbf{x} + \hat{i}$. Thus the $\psi^{(1)}$ fields at either site must either be paired or be absent. This leaves us with three types of term, with which we deal in turn.

A. The trivial term

Choosing all the $\psi^{(1)}$ fields in Eq. (4.4) leaves no operators to change the zero-mode states. The term is therefore zero unless $|\{\eta\}\rangle = |\{\eta'\}\rangle$. Moreover, the matrix elements have no dependence on $\{\eta'\}$ except for possible sign factors from Fermi statistics, which cancel upon multiplying the two matrix elements. Thus this type of term gives a contribution to H^{eff} that is proportional to the unit operator, and uninteresting.

B. The site-site coupling

Consider now the result of choosing all the $\psi^{(0)}$ fields in Eq. (4.4). Then the intermediate state $|j\rangle$ differs from $|\{\eta\}\rangle$ and $|\{\eta'\}\rangle$ only in its zero-mode quanta. The energy denominator is exactly $-g^2/2$ for all accessible states $|j\rangle$, and the corresponding terms in Eq. (4.1) may be evaluated as a completeness sum. This is the site-site coupling in H^{eff} , viz.,

$$H_{\mathbf{x},\mathbf{x}+\hat{i}}^{\text{eff}} = -\frac{2}{g^2} \int ds \psi_{\mathbf{x}+\hat{i},s}^{(0)\dagger} \beta \frac{(1+i\gamma_i)}{2} \psi_{\mathbf{x}s}^{(0)} \int ds' \psi_{\mathbf{x}s'}^{(0)\dagger} \beta \frac{(1-i\gamma_i)}{2} \psi_{\mathbf{x}+\hat{i},s'}. \quad (4.5)$$

The quantity

$$\int ds \psi_{\mathbf{x}+\hat{i},s}^{(0)\dagger} \beta \psi_{\mathbf{x}s}^{(0)} \quad (4.6)$$

contains, according to Eq. (4.3), the integrals $\int ds u^\dagger \beta u$, $\int ds v^\dagger \beta v$, and $\int ds u^\dagger \beta v$. Since $\beta = \rho_1$ is off-diagonal in the chiral basis, each of these consists entirely of cross-terms of the form [see Eq. (B2)] $\int ds f(s)g(s)$. This is the overlap of functions localized at opposite ends of the interval $[0, L]$, and when L is large this integral is $\pm e^{-\kappa_0 L}$, with $\kappa_0 \simeq M - 3$. Thus the terms without γ_i in Eq. (4.5) can be dropped.

We are left with

$$H_{\mathbf{x},\mathbf{x}+\hat{i}}^{\text{eff}} = -\frac{1}{2g^2} \int ds \psi_{\mathbf{x}+\hat{i},s}^{(0)\dagger} \alpha_i \psi_{\mathbf{x}s}^{(0)} \int ds' \psi_{\mathbf{x}s'}^{(0)\dagger} \alpha_i \psi_{\mathbf{x}+\hat{i},s'}^{(0)}. \quad (4.7)$$

The first factor is

$$\int ds \psi_{\mathbf{x}+\hat{i},s}^{(0)\dagger} \alpha_i \psi_{\mathbf{x}s}^{(0)} = \int ds \sum_{\sigma\sigma'} \left(b_{\mathbf{x}+\hat{i},0\sigma}^\dagger u_{0\sigma}^\dagger(s) + d_{\mathbf{x}+\hat{i},0\sigma} v_{0\sigma}^\dagger(s) \right) \alpha_i \left(b_{\mathbf{x}0\sigma'} u_{0\sigma'}(s) + d_{\mathbf{x}0\sigma'}^\dagger v_{0\sigma'}(s) \right). \quad (4.8)$$

It is straightforward to evaluate

$$\begin{aligned} \int ds u_{0\sigma}^\dagger(s) \alpha_i u_{0\sigma'}(s) &= \int ds v_{0\sigma}^\dagger(s) \alpha_i v_{0\sigma'}(s) = 0 \\ \int ds u_{0\sigma}^\dagger(s) \alpha_i v_{0\sigma'}(s) &= -\eta_\sigma^\dagger \sigma_i \eta_{\sigma'} \\ &= \int ds v_{0\sigma}^\dagger(s) \alpha_i u_{0\sigma'}(s). \end{aligned} \quad (4.9)$$

Thus

$$\begin{aligned} \int ds \psi_{\mathbf{x}+\hat{i},s}^{(0)\dagger} \alpha_i \psi_{\mathbf{x}s}^{(0)} &= -\sum_{\sigma\sigma'} \eta_\sigma^\dagger \sigma_i \eta_{\sigma'} (b_{\mathbf{x}+\hat{i},0\sigma}^\dagger d_{\mathbf{x}0\sigma'}^\dagger + d_{\mathbf{x}+\hat{i},0\sigma} b_{\mathbf{x}0\sigma'}) \\ &= -\Psi_{\mathbf{x}+\hat{i}}^\dagger \alpha_i \Psi_{\mathbf{x}} \end{aligned} \quad (4.10)$$

and finally

$$H_{\mathbf{x},\mathbf{x}+\hat{i}}^{\text{eff}} = -\frac{1}{2g^2} \Psi_{\mathbf{x}+\hat{i}}^\dagger \alpha_i \Psi_{\mathbf{x}} \Psi_{\mathbf{x}}^\dagger \alpha_i \Psi_{\mathbf{x}+\hat{i}}. \quad (4.11)$$

The perturbation sum (4.1) also contains a term where H_4 lowers the flux on link $\mathbf{x}\hat{i}$ when creating the intermediate state $|j\rangle$. The numerator is [cf. Eq. (4.2)]

$$\langle \{\eta\} | \int ds \psi_{\mathbf{x}s}^\dagger \beta \frac{(1 - i\gamma_i)}{2} \psi_{\mathbf{x}+\hat{i},s} | j \rangle \langle j | \int ds' \psi_{\mathbf{x}+\hat{i},s'}^\dagger \beta \frac{(1 + i\gamma_i)}{2} \psi_{\mathbf{x}s'} | \{\eta'\} \rangle. \quad (4.12)$$

This gives

$$H_{\mathbf{x}+\hat{i},\mathbf{x}}^{\text{eff}} = -\frac{1}{2g^2} \Psi_{\mathbf{x}}^\dagger \alpha_i \Psi_{\mathbf{x}+\hat{i}} \Psi_{\mathbf{x}+\hat{i}}^\dagger \alpha_i \Psi_{\mathbf{x}}. \quad (4.13)$$

Dropping the β term (4.6) means that the Wilson term in H_4 does not contribute to H^{eff} . It is hence not surprising that the result shown in Eqs. (4.11) and (4.13) is no different from that in the naive nearest-neighbor theory [15].

C. Fierz transformation

Before completing our analysis of Eq. (4.4), we execute a Fierz transformation [15] on H^{eff} in order to express its terms as products of single-site operators. We express

$$(\alpha_i)_{ij}(\alpha_i)_{kl} = \frac{1}{4} \sum_a s_{ia} \Gamma_{il}^a \Gamma_{kj}^a, \quad (4.14)$$

choosing all the Γ^a to be hermitian (e.g., $i\gamma_i$) and satisfying $\text{Tr} \Gamma^a \Gamma^b = 4\delta_{ab}$. The coefficients are

$$\begin{aligned} s_{ia} &= \frac{1}{4} \text{Tr} (\alpha_i) \Gamma^a (\alpha_i) \Gamma^a \\ &= \pm 1, \end{aligned} \quad (4.15)$$

according to whether Γ^a commutes or anticommutes with α_i ,

$$\Gamma^a (\alpha_i) = s_{ia} (\alpha_i) \Gamma^a. \quad (4.16)$$

Thus the Fierz transformation takes the form

$$\begin{aligned} \Psi_i^\dagger (\alpha_i)_{ij} \Psi'_j \Psi_k^\dagger (\alpha_i)_{kl} \Psi_l &= \sum_a s_{ia} \Psi_i^\dagger \Gamma_{il}^a \Psi_l (-\Psi_k^\dagger \Psi'_j + \delta_{jk}) \Gamma_{kj}^a \\ &= -\sum_a s_{ia} \Psi_i^\dagger \Gamma^a \Psi \Psi'^\dagger \Gamma^a \Psi' + 4\Psi_i^\dagger \Psi \end{aligned} \quad (4.17)$$

and the site-site part of the effective Hamiltonian is

$$H_{s-s}^{\text{eff}} = \frac{1}{g^2} \sum_{\mathbf{x}\hat{i}} \sum_a s_{ia} \Psi_{\mathbf{x}}^\dagger \Gamma^a \Psi_{\mathbf{x}} \Psi_{\mathbf{x}+\hat{i}}^\dagger \Gamma^a \Psi_{\mathbf{x}+\hat{i}}. \quad (4.18)$$

We have combined $H_{\mathbf{x},\mathbf{x}+\hat{i}}^{\text{eff}}$ with $H_{\mathbf{x}+\hat{i},\mathbf{x}}^{\text{eff}}$ and dropped the $\Psi^\dagger \Psi$ which is a constant.

D. The single-site term

We return to Eq. (4.4) and examine the term containing $\psi_{\mathbf{x}s}^{(0)}$ and $\psi_{\mathbf{x}+\hat{i},s}^{(1)}$. This term allows $|\{\eta\}\rangle$ to differ from $|\{\eta'\}\rangle$ only at \mathbf{x} . The energy denominator will depend on the energy of

the intermediate $\lambda > 0$ state, and there is no completeness sum. The contribution to H^{eff} is

$$\begin{aligned} \langle \{\eta\} | H_{\mathbf{x}i}^{\text{eff}(1)} | \{\eta'\} \rangle &= \sum_{|j\rangle} \frac{1}{E_{\{\eta\}} - E_j} \langle \{\eta\} | \int ds \psi_{\mathbf{x}+i,s}^{(1)\dagger} \beta \frac{(1+i\gamma_i)}{2} \psi_{\mathbf{x}s}^{(0)} | j \rangle \\ &\quad \times \langle j | \int ds' \psi_{\mathbf{x}s'}^{(0)\dagger} \beta \frac{(1-i\gamma_i)}{2} \psi_{\mathbf{x}+i,s'}^{(1)} | \{\eta'\} \rangle. \end{aligned} \quad (4.19)$$

The intermediate state has a different zero-mode state $|\hat{\eta}\rangle$ at site \mathbf{x} and an additional continuum quantum in state $|(\lambda\sigma)\rangle$ at site $\mathbf{x} + \hat{i}$, and we write it as the product

$$|j\rangle = |\hat{\eta}\rangle_{\mathbf{x}} |\eta'_{\mathbf{x}+i}(\lambda\sigma)\rangle_{\mathbf{x}+\hat{i}}. \quad (4.20)$$

The energy denominator is

$$E_{\{\eta\}} - E_j = -\frac{g^2}{2} - \lambda, \quad (4.21)$$

and the rightmost matrix element is

$$\int ds \sum_{\sigma'} \left(\langle \hat{\eta} | b_{\mathbf{x}0\sigma'}^\dagger | \eta'_{\mathbf{x}} \rangle u_{0\sigma'}^\dagger(s) + \langle \hat{\eta} | d_{\mathbf{x}0\sigma'} | \eta'_{\mathbf{x}} \rangle v_{0\sigma'}^\dagger(s) \right) \beta \frac{(1-i\gamma_i)}{2} \langle (\lambda\sigma) | d_{\mathbf{x}+i,\lambda\sigma}^\dagger | 0 \rangle v_{\lambda\sigma}(s). \quad (4.22)$$

Using orthogonality relations derived in Appendix B, it is easy to show that

$$\int ds u_{0\sigma}^\dagger(s) \frac{1}{2} \alpha_i v_{\lambda\sigma'}(s) = \int ds v_{0\sigma}^\dagger(s) \frac{1}{2} \alpha_i v_{\lambda\sigma'}(s) = 0. \quad (4.23)$$

Moreover,

$$\int ds u_{0\sigma}^\dagger \frac{1}{2} \beta v_{\lambda\sigma'} = 0, \quad (4.24)$$

and the only non-zero matrix element is

$$\int ds v_{0\sigma}^\dagger \frac{1}{2} \beta v_{\lambda\sigma'} = -\delta_{\sigma\sigma'} \mathcal{I}(\lambda), \quad (4.25)$$

where we define

$$\mathcal{I}(\lambda) = \int_0^L ds f_0(s) g_\lambda(s). \quad (4.26)$$

The matrix element (4.22) is thus that of the operator $-\mathcal{I}(\lambda) d_{\mathbf{x}0\sigma}$; treating the other matrix element in Eq. (4.19) in the same way, we get the adjoint, and hence

$$H_{\mathbf{x}i}^{\text{eff}(1)} = \sum_{\lambda\sigma} \frac{\mathcal{I}(\lambda)^2 d_{\mathbf{x}0\sigma}^\dagger d_{\mathbf{x}0\sigma}}{-g^2/2 - \lambda}. \quad (4.27)$$

There is a corresponding term from the flux-lowering matrix element (4.12). The counterpart to Eq. (4.19) is

$$\begin{aligned} \langle \{\eta\} | H_{\mathbf{x}i}^{\text{eff}(2)} | \{\eta'\} \rangle &= \sum_{|j\rangle} \frac{1}{E_{\{\eta\}} - E_j} \langle \{\eta\} | \int ds \psi_{\mathbf{x}s}^{(0)\dagger} \beta \frac{(1-i\gamma_i)}{2} \psi_{\mathbf{x}+i,s}^{(1)} | j \rangle \\ &\quad \times \langle j | \int ds' \psi_{\mathbf{x}+i,s'}^{(1)\dagger} \beta \frac{(1+i\gamma_i)}{2} \psi_{\mathbf{x}s'}^{(0)} | \{\eta'\} \rangle. \end{aligned} \quad (4.28)$$

The rightmost matrix element is now

$$\int ds \sum_{\sigma'} \langle (\lambda\sigma) | b_{\mathbf{x}+\hat{i},\lambda\sigma}^\dagger | 0 \rangle u_{\lambda\sigma}^\dagger(s) \beta \frac{(1+i\gamma_i)}{2} \left(\langle \hat{\eta}_{\mathbf{x}} | b_{\mathbf{x}0\sigma'} | \eta'_{\mathbf{x}} \rangle u_{0\sigma'}(s) + \langle \hat{\eta}_{\mathbf{x}} | d_{\mathbf{x}0\sigma'}^\dagger | \eta'_{\mathbf{x}} \rangle v_{0\sigma'}(s) \right). \quad (4.29)$$

Proceeding as before, we find that the only non-zero integral is

$$\int ds u_{\lambda\sigma}^\dagger \frac{1}{2} \beta u_{0\sigma'} = \delta_{\sigma\sigma'} \mathcal{I}(\lambda), \quad (4.30)$$

which gives the matrix element of the operator $\mathcal{I}(\lambda)b_{\mathbf{x}0\sigma}$. The other matrix element gives the adjoint, and the result is

$$H_{\mathbf{x}i}^{\text{eff}(2)} = \sum_{\lambda\sigma} \frac{\mathcal{I}(\lambda)^2 b_{\mathbf{x}0\sigma}^\dagger b_{\mathbf{x}0\sigma}}{-g^2/2 - \lambda}. \quad (4.31)$$

Since there is no dependence on the link direction i , we obtain a factor of 3; the backward link ($\mathbf{x}-\hat{i}, \mathbf{x}$) gives the same result, giving another factor of 2 in the final result for the single-site term. We note that

$$\sum_{\sigma} b_{\mathbf{x}0\sigma}^\dagger b_{\mathbf{x}0\sigma} + d_{\mathbf{x}0\sigma}^\dagger d_{\mathbf{x}0\sigma} = \Psi_{\mathbf{x}}^\dagger \beta \Psi_{\mathbf{x}} - 1, \quad (4.32)$$

and so the single-site effective Hamiltonian is

$$H_{\text{site}}^{\text{eff}} = -G \sum_{\mathbf{x}} \Psi_{\mathbf{x}}^\dagger \beta \Psi_{\mathbf{x}}, \quad (4.33)$$

with

$$G = 6 \sum_{\lambda} \frac{\mathcal{I}(\lambda)^2}{\lambda + g^2/2}. \quad (4.34)$$

We show in Appendix C that the summation in Eq. (4.34) is convergent.

V. SYMMETRIES OF H^{eff}

The fermion bilinears appearing in H^{eff} form a $U(4)$ algebra that acts on the Dirac indices of Ψ . Identifying the generators of this algebra will permit analysis of the symmetries of the effective theory [15].

We define local charges by

$$\tilde{Q}_{\mathbf{x}}^a = \Psi_{\mathbf{x}}^\dagger \Gamma^a \Psi_{\mathbf{x}}. \quad (5.1)$$

Since the 16 matrices Γ^a are a hermitian basis for the $U(4)$ algebra, so are the $\tilde{Q}_{\mathbf{x}}^a$,

$$[\tilde{Q}_{\mathbf{x}}^a, \tilde{Q}_{\mathbf{x}}^b] = i f^{abc} \tilde{Q}_{\mathbf{x}}^c. \quad (5.2)$$

The effective Hamiltonian can be expressed as

$$H^{\text{eff}} = \frac{1}{g^2} \sum_{\mathbf{x}i} \sum_a s_{ia} \tilde{Q}_{\mathbf{x}}^a \tilde{Q}_{\mathbf{x}+\hat{i}}^a - G \sum_{\mathbf{x}} \tilde{Q}_{\mathbf{x}}^m. \quad (5.3)$$

We have defined $\tilde{Q}_{\mathbf{x}}^m = \Psi_{\mathbf{x}}^\dagger \gamma_0 \Psi_{\mathbf{x}}$, the generator appearing in the single-site term in H^{eff} , soon to be identified as a mass term.

In order to get rid of the s_{ia} sign factors in Eq. (5.3), we absorb them into new local charges $Q_{\mathbf{x}}^a$ according to

$$Q_{\mathbf{x}}^a Q_{\mathbf{x}+\hat{i}}^a = s_{ia} \tilde{Q}_{\mathbf{x}}^a \tilde{Q}_{\mathbf{x}+\hat{i}}^a \quad (5.4)$$

which has the solution

$$Q_{\mathbf{x}}^a = \prod_i (s_{ia})^{x_i} \tilde{Q}_{\mathbf{x}}^a. \quad (5.5)$$

The commutation relations of the new charges are the same as those of the old ones, since

$$\begin{aligned} [Q_{\mathbf{x}}^a, Q_{\mathbf{x}}^b] &= \prod_i (s_{ia} s_{ib})^{x_i} [\tilde{Q}_{\mathbf{x}}^a, \tilde{Q}_{\mathbf{x}}^b] \\ &= \prod_i (s_{ia} s_{ib})^{x_i} i f^{abc} \tilde{Q}_{\mathbf{x}}^c \\ &= \prod_i (s_{ic})^{x_i} i f^{abc} \tilde{Q}_{\mathbf{x}}^c \end{aligned} \quad (5.6)$$

$$= i f^{abc} Q_{\mathbf{x}}^c. \quad (5.7)$$

Equation (5.6) is true because given a and b , the structure factor f^{abc} is nonzero for at most one value of c (i.e., the commutator of Γ^a and Γ^b is just a particular Γ^c), and for this combination of a, b, c we have $s_{ia} s_{ib} = s_{ic}$. Now we can write

$$H^{\text{eff}} = \frac{1}{g^2} \sum_{\mathbf{x}i} \sum_a Q_{\mathbf{x}}^a Q_{\mathbf{x}+\hat{i}}^a - G \sum_{\mathbf{x}} (-1)^{x+y+z} Q_{\mathbf{x}}^m. \quad (5.8)$$

Note that $s_{im} = -1$ for all i .

Defining global charges according to

$$Q^a = \sum_{\mathbf{x}} Q_{\mathbf{x}}^a, \quad (5.9)$$

it is clear that

$$[Q^a, Q^b] = i f^{abc} Q^c \quad (5.10)$$

and

$$[Q^a, H_{s-s}^{\text{eff}}] = 0, \quad (5.11)$$

but $H_{\text{site}}^{\text{eff}} = -G \sum_{\mathbf{x}} (-1)^{x+y+z} Q_{\mathbf{x}}^m$ breaks all charges except for those that commute with Q^m . This breaks the symmetry group from $U(4)$ to $SU(2) \times SU(2) \times U(1)_B \times U(1)_m$, where the new $U(1)_m$ is generated by Q^m and the $SU(2)$'s are generated by the Q^a corresponding to the axial vectors $\sigma \pm \gamma_5 \gamma$.

The neutral states $|\eta\rangle$ at each site can be classified as follows. Begin with the drained state $|\text{Dr}\rangle \equiv d_{\mathbf{x}0+}^\dagger d_{\mathbf{x}0-}^\dagger |0\rangle_{\mathbf{x}}$, which is the unique state with charge $Q = -2$. Since charge is the $U(1)_B$ factor of the $U(4)$, all the generators of $U(4)$ commute with it, and hence $|\text{Dr}\rangle$ is a singlet. From this we form the neutral states $|ij\rangle = \Psi_i^\dagger \Psi_j^\dagger |\text{Dr}\rangle$, where i, j are Dirac indices. Since Ψ_i transforms as a fundamental **4** of $SU(4)$, the states $|ij\rangle$, which are the $|\eta\rangle$ states, transform as the antisymmetric product **6** of two quartets. Under the $SU(2) \times SU(2) \times U(1)_m$ subgroup left unbroken by $H_{\text{site}}^{\text{eff}}$, the **6** is composed of the representations $(0, 0)^{\pm 1}, (\frac{1}{2}, \frac{1}{2})^0$. This is exactly the structure of the naive, nearest-neighbor theory as discussed in [15], with an important exception: There an explicit Dirac mass term furnishes the single-site $Q_{\mathbf{x}}^m$ term in H^{eff} , while here this term is inherent in the domain wall formulation, even without a fermionic mass term.

VI. THE SHAMIR MASS TERM

The final perturbation we consider is the mass term H_m . Acting among the low-energy states $|\{\eta\}\rangle$, it takes the form

$$H_m^{\text{eff}} = m \sum_{\mathbf{x}} \left(\psi_{\mathbf{x}L}^{(0)\dagger} \beta \frac{(1 + \gamma_5)}{2} \psi_{\mathbf{x}0}^{(0)} + \psi_{\mathbf{x}0}^{(0)\dagger} \beta \frac{(1 - \gamma_5)}{2} \psi_{\mathbf{x}L}^{(0)} \right). \quad (6.1)$$

The boundary conditions (A11) allow the replacement $\frac{1}{2}(1 \pm \gamma_5) \rightarrow 1$. Expanding $\psi^{(0)}$, we have

$$H_m^{\text{eff}} = m \sum_{\mathbf{x}} \sum_{\sigma\sigma'} \left[\left(b_{\mathbf{x}0\sigma}^\dagger u_{0\sigma}^\dagger(L) + d_{\mathbf{x}0\sigma}^\dagger v_{0\sigma}^\dagger(L) \right) \beta \left(b_{\mathbf{x}0\sigma'} u_{0\sigma'}(0) + d_{\mathbf{x}0\sigma'}^\dagger v_{0\sigma'}(0) \right) \right. \\ \left. + \left(b_{\mathbf{x}0\sigma}^\dagger u_{0\sigma}^\dagger(0) + d_{\mathbf{x}0\sigma}^\dagger v_{0\sigma}^\dagger(0) \right) \beta \left(b_{\mathbf{x}0\sigma'} u_{0\sigma'}(L) + d_{\mathbf{x}0\sigma'}^\dagger v_{0\sigma'}(L) \right) \right]. \quad (6.2)$$

Now we note that (see Appendix B for notation)

$$u_{0\sigma}(0) = -v_{0\sigma}(0) = \begin{pmatrix} f_0(0) \\ 0 \end{pmatrix} \eta_\sigma \quad (6.3)$$

$$u_{0\sigma}(L) = v_{0\sigma}(L) = \begin{pmatrix} 0 \\ g_0(L) \end{pmatrix} \eta_\sigma, \quad (6.4)$$

and $g_0(L) = -f_0(0) = A_0 \sinh \kappa_0 L \simeq \sqrt{\kappa_0} \simeq \sqrt{M-3}$. Hence

$$\begin{aligned} H_m^{\text{eff}} &= -2\kappa_0 m \sum_{\mathbf{x}} \sum_{\sigma} \left(b_{\mathbf{x}0\sigma}^\dagger b_{\mathbf{x}0\sigma} - d_{\mathbf{x}0\sigma} d_{\mathbf{x}0\sigma}^\dagger \right) \\ &= -2\kappa_0 m \sum_{\mathbf{x}} \Psi_{\mathbf{x}}^\dagger \beta \Psi_{\mathbf{x}}. \end{aligned} \quad (6.5)$$

Shamir's mass term leads to a term in H^{eff} that is of the same form as $H_{\text{site}}^{\text{eff}}$. It thus breaks no hitherto unbroken symmetry of the theory. This confirms our conclusion that the domain wall formulation possesses no axial symmetry in the strong coupling limit. The symmetries that do survive breaking are accidental symmetries of the nearest-neighbor theory.

VII. MORE FLAVORS

One might wonder at this point whether the explicit breaking of chiral symmetry is due in some mysterious way to the anomaly. To settle this point we extend the above analysis to $N_f > 1$. This extension is straightforward [15]. It is easiest to regard Ψ as a spinor with $4N_f$ components and to combine transformations of the Dirac and flavor indices into a $U(4N_f)$ group. Retracing our derivations, one may verify that the effective Hamiltonian is unchanged in form; it is only necessary to substitute $M^{ar} \equiv \Gamma^a \otimes \lambda^r$ for Γ^a in the definition of the charges \tilde{Q} and Q , where λ^r are the generators of the flavor $U(N_f)$. Thus

$$H^{\text{eff}} = \frac{1}{g^2} \sum_{\mathbf{x}i} \sum_{ar} Q_{\mathbf{x}}^{ar} Q_{\mathbf{x}+i}^{ar} - G \sum_{\mathbf{x}} (-1)^{x+y+z} Q_{\mathbf{x}}^m. \quad (7.1)$$

The single-site term is diagonal in flavor and contains $\tilde{Q}_{\mathbf{x}}^m = \Psi_{\mathbf{x}}^\dagger \beta \Psi_{\mathbf{x}}$ as before. The same goes for H_m^{eff} . The single-site and mass terms therefore break $SU(4N_f)$ to $SU(2N_f) \times SU(2N_f) \times U(1)_m$. Again, there is too much symmetry, an indication of doubling; moreover, whatever symmetry is broken by the mass term is already broken by the single site term. Each of these terms is flavor-diagonal and breaks all generators of the form $\gamma_5 \otimes \lambda^a$, including the off-diagonal generators that should be unaffected by anomalies.

For completeness, we present the construction of the state space in which H^{eff} acts. The development follows closely that for the single-flavor case. Start with the drained state $|\text{Dr}\rangle = \prod_f d_{\mathbf{x}0+}^{f\dagger} d_{\mathbf{x}0-}^{f\dagger} |0\rangle_{\mathbf{x}}$. This state has the $U(1)_B$ charge $Q = -2N_f$ and is unique, and hence a singlet under $SU(4N_f)$. To get back to the neutral states, we raise with Ψ_i^\dagger according to $|ij\dots\rangle = (\Psi_i^\dagger \Psi_j^\dagger \dots) |\text{Dr}\rangle$, where i, j are Dirac-flavor indices and there are $2N_f$ of them. Ψ_i transforms in the fundamental representation of $SU(4N_f)$, and hence the neutral states

transform in the antisymmetric representation whose Young diagram has one column of $2N_f$ boxes.

VIII. DISCUSSION

We have shown that an Abelian gauge theory with domain wall fermions breaks chiral symmetry explicitly in strong coupling, and moreover becomes invariant under the same symmetry group as is a theory with naive, nearest-neighbor fermions. Our conclusions are derived from construction of the effective Hamiltonian in strong-coupling perturbation theory. A symmetry-breaking single-site term appears that echoes that produced by a Dirac mass term in the naive theory or by a Shamir mass term in the domain wall theory. We conclude with some speculations and directions for further work.

A. Domain wall fermions

The interpretation of our result is not difficult. The gauge interaction offers a mechanism for communication between the would-be chiral modes that reside on opposite surfaces of the five-dimensional space. As seen above, a chiral mode on one site, living at $s = 0$, communicates with a continuum mode on an adjoining site via the gauge field, and the continuum mode carries the interaction to $s = L$ whence it is passed to the other chiral mode, on the original site, by the gauge field. This is the origin of the single-site term in the effective Hamiltonian. Shamir's mass term, then, merely adds contact between the surfaces to what is already there.

At present, the arguments in favor of exact chiral symmetry in the domain wall theory are perturbative [5] or limited to a finite Euclidean lattice [9]. Numerical results indicate that it is quite possible that non-perturbative physics breaks the symmetry at moderate couplings, with the result depending on the ratio of the fifth dimension to the four-dimensional volume. Our results make more urgent the question of whether chiral symmetry is thus broken even in weak coupling. If it is, then one must demonstrate a way to take a continuum (and infinite-volume) limit at weak coupling in such a way that the symmetry breaking phenomena may be neglected. If it is not, then there must be a phase transition at intermediate coupling. This phase transition would be an interesting topic of research in its own right.

B. The spectrum

Despite the dire consequences for chiral symmetry, it is interesting to pursue the present model a bit further. One may ask about the spectrum of this effective Hamiltonian, following [15]. In the absence of the single site term and of a mass perturbation, H^{eff} represents a Heisenberg antiferromagnet with symmetry group $SU(4N_f)$ and spins in the antisymmetric representation indicated above. The analysis of [15] shows that the ground state of this theory breaks the symmetry spontaneously to $SU(2N_f) \times SU(2N_f) \times U(1)$, where the direction of the $U(1)$ factor is arbitrary, and thus there are $8N_f^2$ Goldstone bosons. A mass perturbation will select the direction of the $U(1)$ by fixing its generator to be Q^m . The broken generators, then, will be all those whose Dirac structure anticommutes with β , including γ_5 . The Goldstone bosons will develop small masses, proportional to m^2 .

In the domain wall model, however, the symmetry-breaking mass-like term is not small. It breaks the $SU(4N_f)$ symmetry immediately, and may invalidate the above picture, which is perturbative in H_m^{eff} . One must start with an antiferromagnet with $SU(2N_f) \times SU(2N_f) \times U(1)$ symmetry, with spins in a reducible representation, and calculate the pattern of spontaneous symmetry breaking. As indicated above, the unbroken symmetry of the Hamiltonian does not include chiral symmetry. In the context of a nearest-neighbor theory perturbed by next-nearest-neighbor (or long-ranged) interactions, these strange symmetries may be interpreted [15] as generalizations of the approximate, non-relativistic $SU(6)$ symmetry (for $N_f = 3$). The value of doing this here is unclear.

C. Non-Abelian gauge fields

In one sense, the introduction of non-Abelian gauge fields changes almost nothing. One simple modification is that the energy denominator contains, instead of $g^2/2$, the energy of a fundamental link of flux given by $\frac{g^2}{2}C_F$, where C_F is the Casimir of the fundamental representation of the $SU(N_c)$ group. Apart from this, the effective Hamiltonian is unchanged from that given above. Thus the explicit breaking of chiral symmetry is still there.

The actual solution of the Hamiltonian, however, is considerably more complicated than in the Abelian case. The reason, of course, is the existence of color singlet baryons. Consider constructing the space of single-site states. The drained state is $|\text{Dr}\rangle = \prod_{c,f} d_{\mathbf{x}0+}^{cf\dagger} d_{\mathbf{x}0-}^{cf\dagger} |0\rangle_{\mathbf{x}}$,

where c is the color index. Again, this is the unique state with $U(1)$ charge (baryon number) equal to $-2N_f N_c$, and hence is a singlet under both Dirac-flavor and color. To return to the neutral sector, apply $\Psi_i^{c\dagger}$ to this state $2N_f N_c$ times. In order to make this state a color singlet, the color indices on the Ψ^\dagger 's must be contracted to singlets. But this is like making $2N_f$ baryons out of $2N_f N_c$ quarks (e.g., four nucleons out of 12 quarks), and there are many color-singlet contraction schemes, each giving a different flavor representation. The representation of the Q^{ar} charges on each site is therefore highly reducible, even before considering the breaking of $SU(4N_f)$.

An additional feature, related to this difficulty, is the possibility of considering states with non-zero baryon density. In view of our results, domain wall fermions offer no advantage over naive fermions for this problem, especially as the latter may be made massless.

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APPENDIX A: REPRISÉ OF FREE DOMAIN WALL FERMIONS

We review here the spectrum of free fermions on a 3-dimensional spatial lattice with continuous time and a continuous fifth dimension with coordinate $0 \leq s \leq L$. The Hamiltonian is

$$H = H_4 + H_5, \tag{A1}$$

where H_4 is the ordinary (massless) Wilson Hamiltonian in $3 + 1$ dimensions,

$$H_4 = \sum \left[\psi^\dagger (-i\boldsymbol{\alpha} \cdot \nabla) \psi - \frac{r}{2} \psi^\dagger \beta \nabla^2 \psi \right], \tag{A2}$$

and H_5 is the continuum kinetic term in the 5th dimension with a mass term,

$$H_5 = \sum \left[\psi^\dagger (-i\alpha_5 \partial_5) \psi + M \psi^\dagger \beta \psi \right]. \tag{A3}$$

The lattice derivatives are

$$\nabla_i \psi = \frac{1}{2}(\psi_{\mathbf{x}+i} - \psi_{\mathbf{x}-i}) \quad (\text{A4})$$

$$\nabla^2 \psi = \sum_i (\psi_{\mathbf{x}+i} + \psi_{\mathbf{x}-i} - 2\psi_{\mathbf{x}}). \quad (\text{A5})$$

Under Fourier transformation,

$$\nabla_i \psi \rightarrow i \sin p_i \psi \equiv i s_i \psi \quad (\text{A6})$$

$$\nabla^2 \psi \rightarrow -2 \sum_i (1 - \cos p_i). \quad (\text{A7})$$

We use the following Dirac matrices:

$$\begin{aligned} \gamma_0 &= \beta = \rho_1 & \gamma_5 &= \rho_3 \\ \boldsymbol{\alpha} &= \rho_3 \boldsymbol{\sigma} & \boldsymbol{\gamma} &= \beta \boldsymbol{\alpha} = -i \rho_2 \boldsymbol{\sigma} \\ \alpha_5 &= i \beta \gamma_5 = \rho_2. \end{aligned} \quad (\text{A8})$$

H_4 takes the explicit form

$$H_4 = \sum_{\mathbf{x}} \int ds \left[-\psi_{\mathbf{x}s}^\dagger \beta \sum_i \left(\frac{(r + i\gamma_i)}{2} \psi_{\mathbf{x}+i,s} + \frac{(r - i\gamma_i)}{2} \psi_{\mathbf{x}-i,s} \right) + 3r \psi_{\mathbf{x}s}^\dagger \beta \psi_{\mathbf{x}s} \right]. \quad (\text{A9})$$

Transforming to (3-d) momentum space, we have

$$H = \int d^3 p ds \psi^\dagger \left(\boldsymbol{\alpha} \cdot \mathbf{s} + r \beta \sum_i (1 - \cos p_i) - i \alpha_5 \partial_5 + M \beta \right) \psi. \quad (\text{A10})$$

The boundary conditions in s are

$$(1 + \gamma_5) \psi(L) = (1 - \gamma_5) \psi(0) = 0. \quad (\text{A11})$$

Writing

$$\psi = \begin{pmatrix} \chi \\ \eta \end{pmatrix}, \quad (\text{A12})$$

where χ and η are 2-spinors, the equations diagonalizing the quadratic form are

$$(\mathbf{s} \cdot \boldsymbol{\sigma}) \chi + \mu \eta - \partial_5 \eta = E \chi \quad (\text{A13})$$

$$-(\mathbf{s} \cdot \boldsymbol{\sigma}) \eta + \mu \chi + \partial_5 \chi = E \eta, \quad (\text{A14})$$

where $\mu = r \sum (1 - \cos p_i) + M$, with the boundary conditions

$$\chi(L) = \eta(0) = 0. \quad (\text{A15})$$

We can work with helicity eigenstates according to

$$(\mathbf{s} \cdot \boldsymbol{\sigma})\chi = h\chi, \quad (\mathbf{s} \cdot \boldsymbol{\sigma})\eta = h\eta. \quad (\text{A16})$$

Now we solve for χ ,

$$\chi = \frac{1}{E - h}(\mu - \partial_5)\eta, \quad (\text{A17})$$

leaving us with

$$(\partial_5^2 - \lambda^2)\eta = 0, \quad (\text{A18})$$

where $\lambda^2 = \mu^2 + h^2 - E^2$. Assuming $\lambda^2 > 0$, the boundary condition on η leaves us the solution

$$\eta(s) = A \sinh \lambda s \eta_h \quad (\text{A19})$$

where η_h is the (normalized) helicity eigenspinor. Then

$$\chi(s) = \frac{A}{E - h}(\mu \sinh \lambda s - \lambda \cosh \lambda s) \eta_h. \quad (\text{A20})$$

The boundary condition $\chi(L) = 0$ gives us the quantization condition

$$\lambda = \mu \tanh \lambda L. \quad (\text{A21})$$

The quantization condition has a non-trivial solution for λ only if $\mu > 0$. We choose r and M so that this solution will exist only near $\mathbf{p} = 0$ and not near $p_i = \pm\pi$. At the points $p_i = 0, \pm\pi$ we have $\mu = 2nr + M$, where n is the number of nonzero components of \mathbf{p} . Thus if we choose $r = -1$ and $0 < M < 2$, the discrete mode at $\mathbf{p} = 0$ has no counterparts (doublers) at the zone faces.

There are also continuum solutions with $\lambda^2 < 0$, which are similar to the single-site continuum modes discussed in the next appendix.

APPENDIX B: THE SINGLE-SITE PROBLEM

The single-site eigenvalue equation is

$$hu \equiv -i\alpha_5 \partial_5 u + \mu\beta u = \lambda u, \quad (\text{B1})$$

where we have defined $\mu = M - 3$. There is no spin dependence, so we write

$$u = \begin{pmatrix} \hat{f}(s)\eta \\ \hat{g}(s)\eta \end{pmatrix} \quad (\text{B2})$$

where η is any 2-spinor. Then we have the equations

$$-\partial\hat{g} + \mu\hat{g} = \lambda\hat{f} \quad (\text{B3})$$

$$\partial\hat{f} + \mu\hat{f} = \lambda\hat{g}. \quad (\text{B4})$$

The boundary conditions

$$(1 - \gamma_5)u(0) = 0 \quad (\text{B5})$$

$$(1 + \gamma_5)u(L) = 0 \quad (\text{B6})$$

translate into

$$\hat{g}(0) = 0 \quad (\text{B7})$$

$$\hat{f}(L) = 0. \quad (\text{B8})$$

We solve Eq. (B3) for \hat{f} ,

$$\hat{f} = \frac{1}{\lambda}(-\partial + \mu)\hat{g}, \quad (\text{B9})$$

and substitute into Eq. (B4), giving

$$\partial^2\hat{g} - \kappa^2\hat{g} = 0 \quad (\text{B10})$$

where $\kappa^2 = \mu^2 - \lambda^2$.

1. $\kappa^2 > 0$

The solutions satisfying the boundary conditions are

$$\hat{f} = \pm A \sinh \kappa(s - L) \equiv \pm f_0(s) \quad (\text{B11})$$

$$\hat{g} = A \sinh \kappa s \equiv g_0(s), \quad (\text{B12})$$

where the eigenvalue condition is satisfied by $\kappa = \kappa_0$,

$$\kappa_0 = \mu \tanh \kappa_0 L, \quad (\text{B13})$$

and thus

$$\lambda = \pm\lambda_0 = \pm \frac{\mu}{\cosh \kappa_0 L}. \quad (\text{B14})$$

As $L \rightarrow \infty$, we have $\kappa_0 \rightarrow \mu$ and $\lambda_0 \rightarrow 0$. Note that we require $\mu = M - 3 > 0$ in order for this solution to exist. The normalization constant when L is large is

$$A_0 \simeq 2\sqrt{\kappa}e^{-\kappa L}. \quad (\text{B15})$$

2. $\kappa^2 < 0$

Defining $\kappa = ik$, we have

$$\partial^2 g + k^2 g = 0 \quad (\text{B16})$$

where $k^2 = \lambda^2 - \mu^2$. The solutions are

$$\hat{f} = \pm A \sin k(s - L) \equiv \pm f_\lambda(s) \quad (\text{B17})$$

$$\hat{g} = A \sin ks \equiv g_\lambda(s), \quad (\text{B18})$$

with

$$k = \mu \tan kL \quad (\text{B19})$$

and

$$\lambda = \pm \frac{\mu}{\cos kL}. \quad (\text{B20})$$

There is a gap to this pseudo-continuum, $|\lambda| > \mu = M - 3$. When $k \gg \mu$ the solutions of Eq. (B19) approach $(2n + 1)\pi/L$, so λ grows without bound as well. The normalization constant is

$$A_\lambda = \left(L - \frac{1}{2k} \sin 2kL \right)^{-1/2}. \quad (\text{B21})$$

The eigenvectors of h are orthogonal, and in particular,

$$\int ds u_{0\sigma}^\dagger v_{\lambda\sigma'} = \int ds v_{0\sigma}^\dagger v_{\lambda\sigma'} = 0. \quad (\text{B22})$$

Moreover, since $\{\gamma_5, h\} = 0$,

$$\int ds u_{0\sigma}^\dagger \gamma_5 v_{\lambda\sigma'} = \int ds v_{0\sigma}^\dagger \gamma_5 v_{\lambda\sigma'} = 0, \quad (\text{B23})$$

and thus

$$\int ds u_{0\sigma}^\dagger (1 \pm \gamma_5) v_{\lambda\sigma'} = \int ds v_{0\sigma}^\dagger (1 \pm \gamma_5) v_{\lambda\sigma'} = 0. \quad (\text{B24})$$

This means, according to Eq. (B2),

$$\int ds f_0 f_\lambda = \int ds g_0 g_\lambda = 0. \quad (\text{B25})$$

APPENDIX C: THE COEFFICIENT OF THE SINGLE-SITE TERM

First we evaluate

$$\begin{aligned}\mathcal{I}(\lambda) &= \int_0^L ds f_0(s)g_\lambda(s) \\ &= A_0A_\lambda \int_0^L ds \sinh \kappa(s-L) \sin ks.\end{aligned}\tag{C1}$$

A straightforward calculation gives

$$\begin{aligned}\mathcal{I}(\lambda) &= A_0A_\lambda\mu^{-1} \frac{\sin 2kL \sinh 2\kappa L}{4(\cosh \kappa L + \cos kL)} \\ &\simeq \frac{\sin 2kL}{2\sqrt{\mu L}}\end{aligned}\tag{C2}$$

as $L \rightarrow \infty$ (with k not too small, i.e., $k \gg 1/L$). The coefficient is thus

$$G = 6 \int dk \rho(k) \frac{1}{4\mu L} \frac{\sin^2 2kL}{\sqrt{k^2 + \mu^2} + g^2/2},\tag{C3}$$

where $\rho(k)$ is the density of solutions of the quantization condition

$$k = \mu \tan kL.\tag{C4}$$

To find ρ , we define $u = k/\mu$ and write Eq. (C4) as

$$\tan^{-1} u = \mu Lu - n\pi\tag{C5}$$

and we differentiate, keeping always to the same branch of the arctan,

$$\left(\frac{1}{1+u^2} - \mu L \right) du = -\pi dn,\tag{C6}$$

so

$$\rho(k) = \frac{dn}{dk} = \frac{L}{\pi} - \frac{1}{\pi} \frac{\mu}{k^2 + \mu^2}.\tag{C7}$$

The other factor in the integral is

$$\sin^2 2kL = \frac{4\mu^2 k^2}{(\mu^2 + k^2)^2}\tag{C8}$$

so the integral is convergent. In the large g limit, $G = 3/g^2$.

[1] K. G. Wilson, in *New Phenomena in Subnuclear Physics*, Erice 1975, edited by A. Zichichi (Plenum, New York, 1977), Part A p. 69.

- [2] L. Susskind, Phys. Rev. D **16**, 3031 (1977).
- [3] H. B. Nielsen and M. Ninomiya, Nucl. Phys. **B185**, 20 (1981); **B193**, 173 (1981).
- [4] D. B. Kaplan, Phys. Lett. B **288**, 342 (1992).
- [5] Y. Shamir, Nucl. Phys. **B406**, 90 (1993).
- [6] R. Narayanan and H. Neuberger, Phys. Lett. B **302**, 62 (1993); Nucl. Phys. **B412**, 574 (1994); **B443**, 305 (1995).
- [7] H. Neuberger, Phys. Rev. D **57**, 5417 (1998).
- [8] Y. Kikukawa and T. Noguchi, hep-lat/9902022.
- [9] V. Furman and Y. Shamir, Nucl. Phys. **B439**, 54 (1995).
- [10] R. G. Edwards, U. M. Heller, and R. Narayanan, Phys. Rev. D **60**, 034502 (1999); A. Ali Khan *et al.* (CP-PACS collaboration), hep-lat/9909049; S. Aoki, T. Izubuchi, Y. Kuramashi, and Y. Taniguchi, hep-lat/9909154; P. Vranas, hep-lat/9911002.
- [11] K. G. Wilson, Phys. Rev. D **10**, 2445 (1974).
- [12] T. Banks, L. Susskind and J. Kogut, Phys. Rev. D **13**, 1043 (1976).
- [13] J. Smit, Nucl. Phys. **B175**, 307 (1980).
- [14] S.D. Drell, M. Weinstein and S. Yankielowicz, Phys. Rev. D **14**, 1627 (1976).
- [15] B. Svetitsky, S. D. Drell, H. R. Quinn and M. Weinstein, Phys. Rev. D **22**, 490 (1980); M. Weinstein, S. D. Drell, H. R. Quinn and B. Svetitsky, *ibid.* **22**, 1190 (1980).
- [16] J. Greensite and J. Primack, Nucl. Phys. **B180**, 170 (1981).
- [17] I. Ichinose and K. Nagao, hep-lat/9910031.
- [18] L. H. Karsten and J. Smit, Nucl. Phys. **B144**, 536 (1978); **B183**, 103 (1981).
- [19] Henceforth we use the notation $\lambda = 0$ for the zero mode, and $\lambda > 0$ for the continuum above $M - 3$.