Quenched chiral logarithms

Stephen R. Sharpe*

Continuous Electron Beam Accelerator Facility, Newport News, Virginia 23606

(Received 4 May 1992)

I develop a diagrammatic method for calculating chiral logarithms in the quenched approximation. While not rigorous, the method is based on physically reasonable assumptions, which can be tested by numerical simulations. The main results are that, at leading order in the chiral expansion, (a) there are no chiral logarithms in quenched f_{π} , for $m_u = m_d$, (b) the chiral logarithms in B_K and related kaon B parameters are, for $m_d = m_s$, the same in the quenched approximation as in the full theory, and (c) for m_{π} and the condensate there are extra chiral logarithms due to loops containing the η' , which lead to a peculiar nonanalytic dependence of these quantities on the bare quark mass. Following the work of Gasser and Leutwyler, I discuss how there is a predictable finite volume dependence associated with each chiral logarithm. I compare the resulting predictions with numerical results: for most quantities the expected volume dependence is smaller than the errors, but for B_V and B_A there is an observed dependence which is consistent with the predictions.

PACS number(s): 12.38.Gc, 11.10.Ef, 14.40.Aq

I. INTRODUCTION

The approximate chiral symmetry of QCD provides strong constraints on the form of amplitudes involving pseudo Goldstone bosons. A classic example is that the pion scattering amplitudes at threshold are related to the pion decay constant [1]

$$T_{I}^{\pi\pi}(s=4m_{\pi}^{2},t=u=0)=c_{I}\frac{m_{\pi}^{2}}{f_{\pi}^{2}}+O(m_{\pi}^{4}).$$
(1)

The c_I are known numerical constants that depend on the isospin I of the pion pair. The corrections of $O(m_{\pi}^4)$ fall into two classes:¹ those that are analytic in m_{π}^2 , coming from nonleading terms in the chiral Lagrangian [2]; and the nonanalytic "chiral logarithms"² proportional to $m_{\pi}^4 \ln(m_{\pi})$. These logarithms come from infrared divergences in pion loops [3,4]. They have the important property that they are proportional to the lowest-order term in the chiral expansion. For example, in the limit $m_{\mu} = m_d = 0$, f_{π} has the expansion [4,2]

$$f_{\pi} = f[1 - \frac{1}{2}L(m_K) + cm_K^2], \qquad (2)$$

where the chiral logarithm is

$$L(m) = \left[\frac{m}{4\pi f}\right]^2 \ln\left[\frac{m^2}{\Lambda^2}\right].$$
 (3)

*Permanent address: Physics Department, FM-15, University of Washington, Seattle, WA 98195.

¹Here and in the following I often use "pion" to refer generically to all of the pseudo Goldstone bosons, i.e., the pions, kaons, and η 's. In the present example, $O(m_{\pi}^4)$ includes the possibility of $O(m_{\pi}^2 m_K^2)$.

²For baryon properties some of the nonanalytic terms are not logarithms, but are proportional to $(m_{\pi}^2)^{3/2}$ [4]. I do not consider baryon properties here.

(I use the normalization such that $f_{\pi} = 93$ MeV.) Chiral symmetry alone does not determine f, Λ , or c (in fact c can be absorbed into Λ), but does fix the coefficient of the chiral logarithm L.

In the continuum, chiral logarithms are useful as a guide to the size of the corrections to the leading-order chiral behavior. If the logarithms are large when one takes a reasonable scale $\Lambda \sim m_{\rho}$, then there is reason to worry about the reliability of leading-order chiral predictions. This approach has been taken for example in Refs. [4] and [5].

On the lattice, chiral logarithms are potentially of greater importance, because they allow one to test whether the contribution of pion loops is being correctly described. These loops give rise not only to chiral logarithms, but also to the pion cloud surrounding hadrons. This cloud is an important part of the structure of hadrons, affecting charge radii and other form factors. The obvious way to search for the chiral logarithm, by studying the dependence of amplitudes on m_{π}^2 or m_K^2 , is not very useful, for it is difficult to disentangle a logarithm from a power series. It is better to study the volume dependence of the amplitude. As explained by Gasser and Leutwyler [6], associated with each chiral log is a known volume-dependent correction. For the particular case of the masses, the correction can also be obtained from a general formula given by Lüscher [7].

Chiral logarithmic corrections have been calculated for a number of quantities in QCD. Those that are of interest here are the pseudo-Goldstone-boson masses and decay constants, the condensate, and B_K . I have extended these calculations to B_V and B_A . These results can be converted into predictions of finite volume dependence for lattice simulations of full QCD, and I give some examples below. At present, however, most lattice calculations of these quantities use the so-called "quenched" approximation, in which internal quark loops are not included. The lack of internal quark loops means that many, though not all, of the pion loops present in full

<u>46</u> 3146

QCD are absent in the quenched approximation. The main focus of this article is to understand how to adapt the calculation of chiral logarithms in full QCD to the quenched approximation. This is done here for the quantities just mentioned, but the method can be extended to many other quantities. These results allow one to predict the finite volume dependence of quenched matrix elements.

In addition to their diagnostic power, chiral logarithms provide an estimate of the size of the error introduced by the quenched approximation. In most quantities the chiral logarithms have a different coefficient in quenched and full theories. For example, it is shown below that there are no leading-order chiral logarithms in f_{π} in the quenched approximation,³ in contrast with the result for full QCD, Eq. (2). Taking $\Lambda = m_{\rho}$, the chiral logarithm in full QCD gives a 6% contribution to f_{π} . This suggests that the absence of chiral logarithms in the quenched approximation should have an effect on f_{π} of this size. This is a suggestion, and not a calculation, because the quantities f and c need not be the same in full and quenched QCD. Nevertheless, even a rough guide is useful since we have so few tools for studying the effect of quenching.

The idea of calculating chiral logarithms in the quenched approximation was first suggested by Morel [8], who was inspired by numerical evidence of nonanalytic terms in m_{π}^2/m_q [9]. In order to study the quenched approximation, he introduced scalar-quark ghost fields which, upon functional integration, produce a determinant which cancels that from the integration over quarks. He showed how to calculate chiral logarithms in the combined strong coupling, large dimension limit, both for color gauge groups SU(2) and SU(3). In Ref. [10] I corrected errors in Morel's result, finding that, for three colors, there were *no* quenched chiral logarithms in f_{π} or m_{π} in the large *d*, strong-coupling limit.⁴ More importantly, I showed how the presence or absence of chiral logarithms could be understood by studying quark diagrams. It is the extension of this "diagrammatic method" that I use in this paper.

As will become clear in the following, this method is neither rigorous nor systematic. The lack of rigor is common to all calculations of quenched chiral logarithms. Such calculations use chiral perturbation theory, the validity of which is predicated on the existence of a welldefined underlying field theory. Quenched QCD is not a well-defined field theory. In particular, quenched amplitudes are neither unitary nor analytic. Nevertheless, as I discuss in Secs. II and III, it is reasonable to assume both that chiral logarithms are present and that they can be calculated using an adaptation of chiral perturbation theory.

The diagrammatic method is also not systematic. It amounts to a set of rules for discarding some of the chiral perturbation theory diagrams and for changing the masses of the pions that appear in some of the loops of the diagrams that are kept. These rules are developed in an *ad hoc* way, quantity by quantity. I do not know how to write them down in general. Bernard and Golterman [11] have recently outlined a more systematic approach in which chiral perturbation theory is extended to include states containing Morel's scalar ghosts. Their method appears to give the same answers for the quantities that I calculate here. I prefer to use the diagrammatic method, however, because it displays the underlying physics very clearly. For more complicated calculations Bernard and Golterman's method is likely to be superior.

A difficulty with calculations of quenched chiral logarithms stems from the fact that one must treat the flavor singlet η' as a pseudo Goldstone boson. This is in contrast with full QCD, where the η' is not a pseudo Goldstone boson because of the axial anomaly. It turns out that the contributions of η' loops do not satisfy the usual power counting rules of chiral perturbation theory, in which each new loop brings with it a power of m_{π}^2 . Thus corrections proportional to powers of $\ln(m_{\pi})$ are obtained from diagrams with any number of loops. This is potentially disastrous, for it might undermine the application of quenched chiral perturbation theory to all quantities. As a first attempt to understand the issue, I have summed the leading logarithms and find that they can be absorbed into a change of definition in the quark mass. The only observable effects are that m_{π} , $\langle \bar{\psi}\psi \rangle$, and related quantities should have a weak finite volume dependence. More study of this issue is required.

Fortunately, it turns out that there are a class of "safe" quantities which are insensitive, at leading order, to η' loops. Examples are the decay constants and kaon *B* parameters for degenerate quarks. I concentrate on these quantities for most of this paper.

Bernard and Golterman have suggested a more ambitious approach [11]. They point out that, although η' loops are not parametrically of higher order in chiral perturbation theory, they are nonleading in a combined chiral and large- N_c expansion. Furthermore, their contribution is numerically small. Thus, from a practical point of view, it may make sense to do a loop expansion, and they have begun such a program.

The work presented here was inspired by numerical results showing volume dependence in B_V and B_A [12]. It turns out that these results are described reasonably well by predictions given here. Present calculations of other quantities are not yet accurate enough to test the predictions.

³Strictly speaking this is only true if $m_u = m_d$, as discussed below.

⁴The reader may be confused by two points upon consulting Ref. [10]. (1) A class of diagrams involving η' loops, which are discussed at length in this article, were overlooked in Ref. [10]. This does not affect the main calculation of that reference, because these diagrams are not present in the large d, strong coupling limit [8]. It does, however, mean that the conclusion drawn in that paper, namely, that there are no leading-order chiral logarithms in quenched f_{π} and m_{π} in general, is false. This error is rectified here. (2) There are chiral logarithms in f_{π} and m_{π} for color SU(2), even in the large d, strong coupling limit. These may explain the numerical results of Ref. [9]. SU(2) is special, however, since there are Goldstone baryons as well as Goldstone pions. In this paper I consider only three or more colors.

The outline of this paper is as follows. In the next section, I recall some useful facts about the chiral Lagrangian, and discuss its extension to the quenched approximation. In Sec. III, I introduce the method of calculating chiral logarithms in the quenched approximation using the example of decay constants. Section IV extends this method to B_K and the related parameters B_V and B_A . In Sec. V, I discuss the contribution of η' loops, and in Sec. VI, I explain how to predict the leading finite volume dependence from the results for chiral logarithms. These predictions are confronted with numerical data in Sec. VII. Section VIII contains some conclusions.

Brief discussions of some of the results of this paper have been given previously in Refs. [13,14,12,15].

II. CHIRAL LAGRANGIAN

Chiral logarithms in full QCD are most simply calculated using the chiral Lagrangian. In this section, I first collect some standard results, and then explain my assumptions for how these are modified in the quenched approximation. I use the notation of Ref. [16] and work in Euclidean space. The Lagrangian is

$$\mathcal{L} = \frac{1}{4} f^2 \operatorname{Tr}(\partial_{\nu} \Sigma \partial_{\nu} \Sigma^{\dagger}) - \frac{1}{2} f^2 \mu \operatorname{Tr}[M(\Sigma + \Sigma^{\dagger})] + O(p^4, Mp^2, M^2) , \qquad (4)$$

where $\Sigma = \exp(2i\pi_a T_a/f)$, the π_a being the pseudo-Goldstone-boson fields. The group generators T_a are normalized such that $\operatorname{Tr}(T_a T_b) = \frac{1}{2} \delta_{ab}$. *M* is the quark mass matrix, in terms of which the pion mass matrix is $(m_{\pi}^2)_{ab} = 4\mu \operatorname{Tr}(MT_a T_b)$. The form of \mathcal{L} is identical for any number of flavors. Flavor dependence enters only through the number of pions $[N^2 - 1$ for flavor SU(*N*)], and through the constants *f* and μ .

In QCD the mass matrix is diag (m_u, m_d, m_s) . I will assume that $m_u = m_d$, which is a good approximation for the amplitudes I discuss, as the terms that I drop are suppressed by $(m_u - m_d)/m_s$ compared to those that I keep. In this isospin symmetric limit the flavor-off-diagonal eigenstates are the π^{\pm} , with $m_{\pi}^2 = 2\mu m_d$, and the kaons $K^{\pm}, K^0, \overline{K}^0$ with $m_K^2 = \mu(m_s + m_d)$, while the flavor-diagonal eigenstates are the π^0 $(T_a = T_3)$, which is degenerate with the π^{\pm} , and the η $(T_a = T_8)$, with $m_{\eta}^2 = \frac{1}{3}\mu(4m_s + 2m_d)$.

The left- and right-handed quark currents can be written in terms of Σ fields

$$L_{\mu a} = \overline{q} T_a \gamma_{\mu} \frac{1}{2} (1 + \gamma_5) q \longrightarrow \frac{1}{2} f^2 \operatorname{Tr}(T_a \Sigma \partial_{\mu} \Sigma^{\dagger}) , \qquad (5)$$

$$\boldsymbol{R}_{\mu a} = \overline{\boldsymbol{q}} T_a \boldsymbol{\gamma}_{\mu \frac{1}{2}} (1 - \boldsymbol{\gamma}_5) \boldsymbol{q} \rightarrow \frac{1}{2} f^2 \operatorname{Tr}(T_a \boldsymbol{\Sigma}^{\dagger} \partial_{\mu} \boldsymbol{\Sigma}) .$$
 (6)

The Euclidean space Dirac matrices are Hermitian and satisfy $\{\gamma_{\mu}, \gamma_{\nu}\} = \delta_{\mu\nu}$. The expansion of L_{μ} in terms of the pion field is

$$L_{\mu a} = -if \left[\operatorname{Tr}(T_a \partial_{\mu} \pi) - \frac{i}{f} \operatorname{Tr}(T_a [\partial_{\mu} \pi, \pi]) - \frac{2}{3f^2} \operatorname{Tr}(T_a [\pi[\pi, \partial_{\mu} \pi]]) + O(\pi^4) \right].$$
(7)

 R_{μ} has the same expansion except that $\pi \rightarrow -\pi$. The pion decay constant is defined by

$$0|\bar{q}T_{a}\gamma_{\mu}\gamma_{5}q|\pi_{b}\rangle = \langle 0|L_{\mu a} - R_{\mu a}|\pi_{b}\rangle = p_{\mu}f_{\pi}\delta_{ab} .$$
(8)

Comparing with Eq. (7) we see that $f_{\pi} = f$ at lowest order in chiral perturbation theory.

The chiral Lagrangian can be extended to include the flavor-singlet pseudoscalar meson [2]. Throughout this paper I refer to this meson as the η' , despite the fact that the physical η' is not exactly a flavor singlet. I use the label "pion" to refer to the nonsinglet pseudo Goldstone bosons, not including the η' . To include the η' in the chiral Lagrangian one begins by defining a U(N) matrix

$$U = \sum e^{i\phi_0}, \quad \phi_0 = 2\eta' / (\sqrt{2N}f) \quad , \tag{9}$$

in terms of which the extended Lagrangian can be written

$$\mathcal{L} = + \frac{1}{4} f^2 V_1(\eta') \operatorname{Tr}[\partial_{\nu} U \partial_{\nu} U^{\dagger}] - \frac{1}{2} f^2 V_2(\eta') \mu \operatorname{Tr}[MU] - \frac{1}{2} f^2 V_2(\eta')^* \mu \operatorname{Tr}[MU^{\dagger}] + V_0(\eta') + V_5(\eta') \partial_{\nu} \eta' \partial_{\nu} \eta' .$$
(10)

(The definition of the potentials differs slightly from those of Ref. [2].) The potentials are arbitrary aside from the following conditions: V_0 , V_1 , and V_5 are real and even functions of their arguments; V_2 satisfies $V_2(\eta')^* = V_2(-\eta')$; and $V_1(0) = V_2(0) = 1$. The arbitrariness implies that there are very few constraints on the couplings of the η' .

This Lagrangian simplifies when N_c , the number of colors, is taken to infinity. This limit is of interest here because it is similar to the quenched approximation. From the analysis of Ref. [2], it can be seen that, when $N_c \rightarrow \infty$, one can set $V_1 = V_2 = 1$, and $V_0 = V_5 = 0$. (V_0 does contribute a constant term proportional to N_c^2 , but this only affects the vacuum energy.) In other words, the Lagrangian has the same form as the original chiral Lagrangian, Eq. (4), except that Σ is replaced by U. Dependence on N_c enters only through the decay constant, which is proportional to $\sqrt{N_c}$. The symmetry group has thus enlarged to $\mathrm{U}(N)_L \times \mathrm{U}(N)_R$, with the η' being the additional pseudo Goldstone boson. The physical reason for this simplification is that the terms in Eq. (10) which differentiate the η' from the pions involve intermediate gluons, and are suppressed by powers of $1/N_c$.

Expanding the large N_c Lagrangian in terms of Σ and η' one finds

$$\mathcal{L}(N_c \to \infty) = \mathcal{L}_0 + \mathcal{L}_m , \qquad (11)$$

$$\mathcal{L}_{0} = \frac{1}{4} f^{2} \left[\operatorname{Tr}[\partial_{\mu} \Sigma \partial_{\mu} \Sigma^{\dagger}] + \frac{2}{f^{2}} \partial_{\mu} \eta' \partial_{\mu} \eta' \right], \qquad (12)$$

$$\mathcal{L}_{m} = -\frac{1}{2} f^{2} \mu \{ \cos(\phi_{0}) \operatorname{Tr}[\boldsymbol{M}(\boldsymbol{\Sigma} + \boldsymbol{\Sigma}^{\dagger})] + i \sin(\phi_{0}) \operatorname{Tr}[\boldsymbol{M}(\boldsymbol{\Sigma} - \boldsymbol{\Sigma}^{\dagger})] \} .$$
(13)

Three features of this result are important in the following.

(i) The η' couples to pions through the mass term \mathcal{L}_m ,

but does not couple through the derivative term \mathcal{L}_0 . This is not true at finite N_c because the potential V_1 will contain even powers of the η' field.

(ii) The masses of the pseudo Goldstone pions differ in general from those at $N_c = 3$. Diagonalizing the mass matrix, one finds that the η' , η , and π^0 mix. The eigenstates have flavor compositions \bar{ss} , \bar{dd} , and $\bar{u}u$, with masses $2\mu m_s$, $2\mu m_d$, and $2\mu m_u$, respectively. For $m_d = m_u$ one can use the usual π^0 and η linear combinations, since they are degenerate. Only if all three quarks are degenerate, however, is the flavor singlet η' an eigenstate.

(iii) The nonsinglet chiral currents that follow from the Lagrangian are the same as those of Eqs. (5) and (6), except that Σ is replaced by U. It is simple to see, however, that the η' drops out of the expressions, so that the currents are in fact unchanged. This can be seen explicitly in the expansion of $L_{\mu a}$, Eq. (7). The first term vanishes for the η' because of the flavor trace, and all subsequent terms involve commutators, which are zero for the η' .

The fundamental assumption of this paper is that the "nonhairpin" interactions involving pions and η 's are described by a Lagrangian which has, at leading order in the chiral expansion, the same form as the large N_c Lagrangian, Eq. (11). The constants f and μ will, however, differ from those in both full QCD and the $N_c \rightarrow \infty$ limit. There are additional "hairpin" interactions involving the η' , and possibly pions, but not involving pions alone. Hairpin interactions are those in which the quark and antiquark fields in at least one of the external η' mesons are contracted together. Examples appear in Figs. 3(d)-3(i). Before discussing these additional interactions, I explain the justification for the assumption just announced. This assumption has three parts.

(1) Interactions involving different numbers of nonsinglet pions, but no flavor singlet mesons, are related to one another in the same way as in full QCD. For example, the pion scattering lengths in the chiral limit can be expressed in terms of f_{π} according to the formulas of Weinberg [1]. In full QCD the form of the chiral Lagrangian follows from three inputs [17]: (a) Ward identities between correlation functions; (b) pion domination of low momentum amplitudes; and (c) analyticity and unitarity of the amplitudes. In the quenched approximation one has only the first two ingredients: the Ward identities in the quenched approximation are the same as in full theories, at least in lattice regularized theories [18,19], and pion dominance is a good approximation in practice. But the amplitudes are neither analytic nor unitary. In fact, it is only analyticity that is important for the $O(p^2)$ terms in the Lagrangian; the amplitudes from these terms are not themselves unitary. What is required is that multipion amplitudes have a polynomial dependence on the kinematic invariants. This is the real assumption being made concerning the quenched approximation. For staggered fermions, it is supported by numerical evidence for amplitudes involving up to four pions [20]. This is actually sufficient for the chiral logarithms discussed here, which require only the two- and four-point vertices from the $O(p^2)$ terms in the Lagrangian.

(2) Non-hairpin interactions of η' mesons with pions come entirely from \mathcal{L}_m , and are related to interactions of pions with each other by flavor factors. In the large- N_c limit this result follows because the quark diagrams are the same with or without η' mesons. This equivalence of diagrams is also true in the quenched approximation. In particular, there are no internal quark loops in either the large- N_c limit or the quenched approximation. Interactions involving intermediate gluons, which vanish when $N_c \rightarrow \infty$, involve hairpins and are discussed below. This distinction is useful because it turns out that, for some quantities, such hairpin interactions do not contribute at leading order. This will become clearer in the following sections. Thus I assume that the form of the effective Lagrangian describing the pion- η' interactions is the same in the quenched and $N_c \rightarrow \infty$ theories. The coefficients in the Lagrangians will differ, however, because the gluon interactions are different.

(3) The interaction of the η' with itself, through both the kinetic and mass terms, is related to the selfinteractions of pions by flavor factors. This applies only to nonhairpin interactions resulting from diagrams in which the η' does not annihilate into gluons. The justification for this assumption is as for the previous one: the result is true in the large- N_c limit, and the quenched diagrams involving η' mesons are related to those involving pions by the same flavor factors as in the large- N_c limit. In particular, the mass eigenstates are the same as in that limit.

An important consequence of the assumption that the quenched Lagrangian has the same form as that for $\mathcal{L}(N_c \to \infty)$ is that the nonsinglet chiral currents are also the same. As explained above, this means that the η' does not couple to these currents.

Although the form of the quenched Lagrangian is similar to that for full QCD, the calculation of loop corrections using chiral perturbation theory involves additional rules. According to these rules certain loop diagrams are kept, while others are discarded. The chiral perturbation theory diagrams kept are those corresponding to quark diagrams that do not contain internal quark loops. The method by which one determines the correspondence between pion and quark diagrams is best explained by example, and is the subject of the following section.

I now return to the additional hairspin interactions involving the η' . It is here that the quenched and large- N_c theories differ: such diagrams exist for the quenched approximation, but vanish in the $N_c \rightarrow \infty$ limit. For example, there is a "disconnected" contribution to the η' twopoint function in which the quarks annihilate into gluons. It is this diagram, and it iterates including any number of quark bubbles, that gives the η' its additional mass in full QCD. More generally, such diagrams give rise to the η' dependence in the potentials V_0 , V_1 , V_2 , and V_5 , and thus convert the Lagrangian $\mathcal{L}_0 + \mathcal{L}_m$ into one with the form of Eq. (10).

In quenched calculations one must include the contribution from η' loops since the η' is light. This means that the additional vertices coming from the V_i may be needed. Thus the coefficients of chiral logarithms due to η' loops are not, in general, predicted in terms of the leading-order coefficients.⁵ They depend on new parameters which even in full QCD are either poorly constrained by phenomenology, or not known at all. It turns out, however, that there is a class of quantities for which the η' does not enter at one-loop order, so that the additional η' interactions are not needed. This class includes f_{π} and the kaon *B* parameters, in the limit that all quarks are degenerate. These "safe" quantities are the main focus of this article.

For the remaining quantities one must include η' loops. As I explain in Sec. V, these loops do not obey usual power counting rules of chiral perturbation theory, according to which each loop brings an additional power of m_{π}^2 . Instead, the leading-order chiral logarithms are not suppressed by m_{π}^2 , and come from diagrams with any number of η' loops. These diagrams are of a simple form, however, and require only the additional η'^2 interactions.⁶ These can be parametrized as

$$\mathcal{L}_{\eta'} = \frac{A-1}{2} \partial_{\nu} \eta' \partial_{\nu} \eta' + \frac{m_0^2 A}{2} {\eta'}^2 , \qquad (14)$$

where the first term comes from V_5 , the second from V_0 . The most important feature of $\mathcal{L}_{\eta'}$ is that m_0 is a fixed scale, of $O(\Lambda_{\rm QCD})$, and does not vanish in the chiral limit. The parametrization of Eq. (14) is chosen so that, if this vertex could be iterated as in full QCD, the η' would get a mass

$$m_{\eta'}^2 = m_0^2 + \frac{2\mu m}{A} . \tag{15}$$

In the approximation that A = 1, the experimental η and η' masses imply that $m_0 \approx 0.9$ GeV. I use these values when making estimates of the quantitative importance of η' loops.

III. CHIRAL LOGARITHMS IN f_{π} AND f_{K}

I begin by summarizing the calculation of the chiral logarithmic corrections to the decay constant in a theory with N light flavors, with quark loops included. In chiral perturbation theory there are two one-loop diagrams, and they are shown, along with the leading-order diagram, in Fig. 1. The circle represents the four-pion vertex obtained by expanding Eq. (4). The square represents the axial-vector current $\frac{1}{2}(L_{\mu a} - R_{\mu a})$, the expansion of



FIG. 1. Diagrams contributing to f_{π} : (a) leading order; (b) wave-function renormalization; (c) vertex renormalization. The lines represent pseudo Goldstone bosons, the box is the axial-vector current, and the circle is the vertex coming from \mathcal{L} .

which is as in Eq. (7), except that only terms with an odd number of pions appear. The crucial point about these vertices is that they are completely determined in terms of the lowest-order parameters f and μ .

As explained in the preceding section, the leadingorder result is $f(\pi_d)=f$, for all flavors d. The corrections from Figs. 1(b) and 1(c) are multiplicative:

$$f(\pi_d) = f[1 + X(\text{Fig. 1(b)}) + X(\text{Fig. 1(c)}) + O(m_{\pi}^2)].$$
(16)

The nonlogarithmic corrections proportional to m_{π}^2 come from $O(p^4)$ terms in the chiral Lagrangian. They do not give rise to finite volume dependence (see Sec. VI), and I will drop them from subsequent equations.

Figure 1(b) gives rise to both mass and wave-function renormalization, but only the latter effects f_{π} . To contribute to wave-function renormalization, the vertex in Fig. 1(b) must contain derivatives, and thus must come from \mathcal{L}_0 . This is an important observation to which I return below. The contribution of Fig. 1(b) is

$$X(\text{Fig. 1(b)}) = \frac{1}{3f^2} \sum_{c} \text{Tr}([T_c, [T_c, T_d]]T_d) \int_{k} G(k, m_c) ,$$
(17)

where the measure is $\int_{k} = \int [d^{4}k/(2\pi)^{4}]$, and the integrand is the pion propagator $G(k,m) = 1/(k^{2}+m^{2})$. The sum runs over the $N^{2}-1$ pions, and is to be done in the mass eigenstate basis, in which the mass of the pion of flavor T_{c} is labeled m_{c} .

Figure 1(c) renormalizes the vertex; its contribution turns out to be proportional to the wave-function renormalization

$$X(Fig. 1(c)) = -4X(Fig. 1(b))$$
. (18)

Thus the fractional correction to the decay constant is

$$\frac{f(\pi_d) - f}{f} = -\sum_c \operatorname{Tr}([T_c, [T_c, T_d]] T_d) I_1(m_c) . \quad (19)$$

Here I have defined one of the standard integrals that appear at one loop:

$$I_1(m) = \frac{1}{f^2} \int_k G(k,m) .$$
 (20)

The integral must be regulated, and I use a sharp cutoff, $|k| < \Lambda$.⁷ The result is

⁵In full QCD this problem does not arise because the η' is heavy, and does not give rise to chiral logarithms. The uncertainties only enter through the $O(p^4)$ terms in the chiral Lagrangian.

⁶Strictly speaking there are also two terms involving V_2 which contribute. (1) A mass term proportional to $V''_2(0)\eta'^2 \text{Tr}(M)$. Using arguments similar to those of the following section, one finds that the Tr(M) corresponds to an additional quark loop. Since this loop is absent in the quenched approximation, this term should not be included. (2) A vertex proportional to $V'_2(0)\eta'\text{Tr}(M\eta'\pi^2)$. This vertex actually appears in Fig. 3(f), and does contribute to mass renormalization. By redefining the U field, however, one can absorb $V'_2(0)$ into the constant A [11]. Thus there is no loss of generality if one excludes this term.

⁷As is well known, such a momentum cutoff violates chiral symmetry. This can be corrected by adding a quartically divergent counterterm, but this does not contribute to any of the processes considered here [21].

$$I_{1}(m) = \frac{\Lambda^{2}}{(4\pi f)^{2}} + L(m) + O\left[\frac{m^{2}}{\Lambda^{2}}\right], \qquad (21)$$

where L(m) is the chiral logarithm defined in Eq. (3). The quadratic divergence can be absorbed by a redefinition of f.

For QCD, the general formula gives

$$f_{K} = f(1 - \frac{3}{8} [2I_{1}(m_{K}) + I_{1}(m_{\eta}) + I_{1}(m_{\pi})]), \qquad (22)$$

$$f_{\pi} = f(1 - \frac{1}{2} [I_1(m_K) + 2I_1(m_{\pi})]) , \qquad (23)$$

in agreement with the Refs. [4,2,22]. For N degenerate quarks the result is

$$f_{\pi} = f \left[1 - \frac{N}{2} I_1(m_{\pi}) \right],$$
 (24)

in agreement with Ref. [6].

I now turn to the quenched theory. The result of the following discussion is extremely simple: except for η' loops, which only contribute if the quarks are not degenerate, there are *no leading-order chiral logs* in decay constants. I explain this conclusion in gory detail, as the insight obtained can be applied directly to more complicated calculations.

The chiral logarithm comes from the infrared part of the loop integral, |p| < m. The coefficient of the logarithm is determined by the residue of the pole at $p^2 = -m^2$. In general this residue is itself determined by a physical, on-shell amplitude: for wave-function renormalization it is the part of the four-pion scattering amplitude proportional to the kinematic variable t, evaluated at threshold; for vertex renormalization it is part of the amplitude for the axial-vector current to convert one pion to two pions. In other words, to calculate the coefficient of the logarithm one can factorize the loop in Figs. 1(b) and 1(c) into two pieces: an on-shell pion propagator, and a physical scattering or conversion amplitude. Since these amplitudes are evaluated for small p, they are well described by lowest-order chiral perturbation theory, at least for small enough m. This is why the coefficient of the logarithm can be calculated in terms of only the coefficients of the lowest-order Lagrangian.

It is because of this factorization that one can hope to calculate the coefficient of the logarithm in the quenched approximation. As discussed in the preceding section, it is reasonable to assume that the leading-order expressions for both the scattering amplitudes and the chiral currents are the same in the quenched approximation as in the full theory, except for differences in the coefficients f and μ . Thus the coefficient of the chiral logarithms coming from each diagram is the same function of these coefficients. In other words, the calculation of individual chiral perturbation theory diagrams proceeds by the same method in both theories. The only differences in the quenched calculation are that the η' can propagate in the loops and that diagrams which involve internal quark loops should be discarded.

To determine what these differences mean in practice it is necessary to find the quark diagrams which correspond to the diagrams of chiral perturbation theory. Figure 2



FIG. 2. Quark diagram corresponding to Fig. 1(a). The lines represent quark propagators, as described in the text. The box is the axial current. The apex where the propagators meet is the operator used to create the pion.

shows the quark diagram corresponding to Fig. 1(a). The quark lines are to be interpreted as follows. If one were doing a lattice (or any other nonperturbative) calculation of f_{π} one would calculate the diagram of Fig. 2, with the lines corresponding to quark propagators in a background gauge field. The result would then be averaged over all the background fields, using either the full measure including quark loops, or the quenched measure. The resulting value of f_{π} would be correct to all orders in m_{π}^2 and, in particular, would include the chiral logarithms. In fact, the lines in Fig. 2 are almost the full propagators, except that the propagation is restricted so that there is no long-distance pion contribution. If the full measure is used, this restriction applies also to the quark loops contained in the background field. With this restriction, the diagram does not produce chiral logarithms. It does contribute terms of all orders in m_{π}^2 , but we are interested only in the lowest-order contribution.

This provides a definition of the meaning of the lines in quark diagrams that is both vague and difficult to implement. These shortcomings are not important, however, because the diagrams serve only as a method for tracing the flow of the flavors of the quark and antiquark. This aspect of the diagrams is completely well defined.

The quark diagrams contributing to wave-function renormalization [i.e., Fig. 1(b)] are shown in Fig. 3. These diagrams are obtained as follows. First, draw all possible quark diagrams for the pion scattering amplitude in which at least two of the pions have the flavor under consideration. One of these two is the external pion, the other couples to the axial-vector current. Next, join the remaining two pions from the scattering amplitude with a pion, or an η' , propagator. Figure 3 shows the topologies of the diagrams that result. They represent the different ways in which flavor can flow in the two step process of pion scattering followed by pion propagation. It is only because the coefficient of the chiral logarithm is deter-



FIG. 3. Quark diagrams corresponding to Fig. 1(b).

mined by this factorized limit that the flavor flow is well defined.

Having understood the meaning of the diagrams in Fig. 3, I now analyze them in turn. The results of this analysis are summarized in Table I. The conclusion is that, in the full theory, only Fig. 3(a) gives leading-order chiral logarithms in wave-function renormalization. This means that it alone must correspond to the chiral perturbation theory diagram, Fig. 1(b). In the quenched approximation, however, this diagram is absent. Furthermore, if the quarks are degenerate, none of the other diagrams contribute leading-order quenched chiral logarithms. Thus wave-function renormalization does not contain leading-order chiral logarithms in the quenched approximation.

I begin by analyzing the diagrams in full QCD. It turns out that Figs. 3(d)-3(i) do not contribute leadingorder chiral logarithms. This is because they all contain one or more mesons annihilating into gluons. This can only occur if the meson is the η' , because at leading order the gluonic amplitude is flavor symmetric. But in full QCD the η' is massive and its loops do not give chiral logarithms. The only possible contributions are thus from Figs. 3(a)-3(c). As I now explain, however, Figs. 3(b) and 3(c) give rise only to nonleading chiral logarithm proportional to $m_{\pi}^4 \ln(m_{\pi})$.

The extra power of m_{π}^2 comes from the four-pion vertex. The two types of diagram that contribute to pion scattering are shown in Fig. 4: diagram (a) is required for Fig. 3(a), while diagram (b) is required for Figs. 3(b) and 3(c). I claim that the amplitude from (a) is $O(p^2)$, while that from (b) is $O(p^4)$. (Here I am using p^2 loosely to refer to both p^2 and m_{π}^2 .) This is true both for the vertex coming from the derivative term in the chiral Lagrangian (which is that required for wave-function renormalization) and for the vertex from the mass term.

For a theory with four or more flavors this can be seen as follows. One can choose the flavors of the external pions such that only Fig. 4(b) contributes, e.g., $\bar{u}d$, $\bar{d}u$, $\bar{s}c$, and $\bar{c}s$. For these flavors the vertices from the leadingorder chiral Lagrangian give no contribution, because the flavor trace vanishes. Only terms with two flavor traces contribute, e.g., $Tr(\partial_{\mu}\Sigma\partial_{\mu}\Sigma^{\dagger})^2$, but these contributions are of $O(p^4)$. For general external flavors, there will be contributions to the scattering amplitude from both Figs. 4(a) and 4(b). Those from the disconnected diagram (b) will, however, be of $O(p^4)$, because the quark diagram is not affected if we simply change the flavor labels.⁸ Thus the only contributions of $O(p^2)$ can be from Fig. 4(a). That this diagram does contribute at $O(p^2)$ can be seen

TABLE I. Contributions of the diagrams of Fig. 3 to chiral logarithms in wave-function renormalization. "Leading" indicates a term proportional to $m_{\pi}^2 \ln(m_{\pi})$, "nonleading" a term proportional to $m_{\pi}^4 \ln(m_{\pi})$.

Diagram	Full theory $(N \ge 4)$	Quenched approximation	
(a)	Leading	Absent (no loop)	
(b)	Nonleading	Nonleading Absent (no loop)	
(c)	Nonleading		
(d),(e)	Absent (no η')	$\propto \delta m_a$	
(f)	Absent (no η')	Absent	
(g),(h),(i)	Absent (no η')	Absent (no loop)	

by choosing flavors such that it is the only contribution, e.g., $\overline{u}d$, $\overline{d}s$, $\overline{s}c$, and $\overline{c}u$. The $O(p^2)$ terms in the chiral Lagrangian do give a vertex with these flavors, because they can be connected together in a single flavor trace.

This argument requires only that the leading-order interactions are described by the chiral Lagrangian. Given the fundamental assumption of this paper, namely that the Lagrangian for quenched interactions has the same form as in the full theory, the argument of the preceding paragraph applies also to the quenched theory. In fact, it applies also if the external particles are flavor singlets. Since the background configurations in the quenched approximation are by construction independent of the number of flavors of the full theory one starts with, the argument also applies to quenched QCD.

An interesting question is whether Figs. 3(b) and 3(c)give rise to leading-order chiral logarithms in full QCD. Since QCD has only three light flavors, the argument given above does not work: none of the possible external flavors selects Fig. 4(b) alone. This means that the question cannot be asked in terms of correlation functions of operators, but only in terms of individual contractions. Thus one must have a nonperturbative regulator, such as the lattice. Although I do not know of a proof, I think it likely that the $N \ge 4$ result carries over to N = 3, for the following reasons. First, with staggered fermions, one can show, given certain assumptions, that Fig. 4(b) is $O(p^4)$ irrespective of the number of dynamical quarks [20]. Second, the contribution of wave-function renormalization in a theory with N degenerate quarks, Eq. (17), comes with an explicit factor of N for $N \ge 2$. This suggests that Fig. 3(a) alone is contributing, because there are N flavors of quark running in the loop. This is not conclusive, however, because there could be hidden Ndependence in f and μ . Finally, the claim that only Fig. 3(a) contributes is consistent with the $N_c \rightarrow \infty$ limit, in which both the vertex from the leading-order chiral Lagrangian and that from Fig. 3(a) are of $O(1/N_c)$, while



FIG. 4. Quark diagrams contributing to pion scattering.

⁸One might worry about the validity of this argument for nondegenerate quarks, for then changing the flavor labels also means changing the quark masses. This is not a problem, however, because there is always a theory in which, for external flavors chosen so that only the disconnected diagram contributes, the masses of the quarks are the same as those in the diagram under consideration. In all such theories the disconnected components are $O(p^4)$.

those from Figs. 3(b) and 3(c) are of $O(1/N_c^2)$.

I now turn to the analysis of the quark diagrams in the quenched approximation. Of the first three diagrams, only Fig. 3(b) is present in the quenched approximation, as the other two involve quark loops. As I have just argued, however, Fig. 3(b) gives nonleading chiral logarithms not only in the full theory (with $N \ge 4$), but also in the quenched approximation. Thus these three diagrams do not contribute leading-order chiral logarithms to quenched f_{π} .

The remaining diagrams involve η' propagation. Of these, Figs. 3(g)-3(i) require internal quark loops, and so are not present in the quenched approximation. Figures 3(d)-3(f), however, do give potential contributions of a type unrelated to those in full QCD. Fortunately, if the quarks are degenerate, this contribution vanishes at leading order. This is because, as explained in Sec. II, the η' does not appear in the four-pion vertex containing derivatives. (Recall that for wave-function renormalization the vertex must contain derivatives.) For degenerate quarks, one can use the usual basis of π_0 , η , and η' , and since the η' does not couple, Figs. 3(d)-3(f) simply vanish. There will, however, be contributions at higherorder chiral in perturbation theory.

In fact, Fig. 3(f) does not contribute leading-order chiral logarithms to f_{π} even for nondegenerate quarks. This is because only the V_2 term in the full chiral Lagrangian, Eq. (10), has the correct flavor structure to correspond to the vertex in Fig. 3(f). Since this term has no derivatives, however, it cannot contribute to wavefunction renormalization irrespective of the quark masses.

This is not true of Figs. 3(d) and 3(e). For nondegenerate quarks, the neutral eigenstates have flavors \overline{ss} , \overline{dd} , and $\overline{u}u$, and each of these individually does couple to the derivative term in the Lagrangian. It is simple to see that the contribution to wave-function renormalization vanishes for degenerate quarks because of a cancellation between the two diagrams. For nondegenerate quarks, this cancellation does not occur because the neutral eigenstates are different. Thus there are extra contributions proportional to the difference in mass of the two quarks making up the pion. For example, this extra contribution is proportional to $m_d - m_u$ for the physical f_{π} . These extra terms are discussed in Ref. [11], and I do not consider them further here.

For mass renormalization, Figs. 3(d)-3(f) all contribute even if the quarks are degenerate, as discussed further in Sec. V.

Finally, I turn to vertex renormalization, Fig. 1(c). The quark diagrams are in one-to-one correspondence with those in Fig. 3, and the analysis is almost identical. The result is that the summary of Table I applies both to vertex and wave-function renormalization. In Fig. 5, I show the diagrams analogous to Figs. 3(a)-3(f). I do not show those corresponding to Figs. 3(g)-3(i), since they contribute neither to full nor to quenched theories. Figure 5(a) is the only diagram contributing leading-order chiral logarithms in the full theory, but is absent in the quenched approximation. The three pion vertex in Figs. 5(b) and 5(c) is absent in leading order because it requires



FIG. 5. Quark diagrams corresponding to Fig. 1(c).

two flavor traces, whereas the axial current contains a single flavor trace [Eq. (7)]. Thus these diagrams contribute only nonleading chiral logarithms. This is true also in the quenched theory, for which only diagram (b) is present. The remaining three diagrams require an η' in the loop, and are thus absent in the full theory. They are present in general in the quenched theory, but vanish for degenerate quark masses because, as explained in Sec. II, the η' does not couple to the leading-order representation of the axial current.

In summary, the leading-order chiral logarithms in decay constants, which are present in full QCD, are absent in the quenched approximation. There are additional chiral logarithms from η' loops, which depend on parameters other than f and μ , but these are only present for nondegenerate quarks. These extra logarithms are discussed in Ref. [11].

IV. CHIRAL LOGARITHMS IN B PARAMETERS

A. B_K with dynamical quarks

To set the stage for the calculation of the chiral logarithms in B_K in the quenched approximation, I review the calculation in theories with dynamical quarks. The discussion will apply to theories with a d quark, an s quark, and any number of other light quarks.

The kaon B parameter is defined as a ratio of matrix elements:

$$\mathcal{M}_{K} = \langle \bar{K}^{0} | \mathcal{O}_{K} | K^{0} \rangle = B_{K} \mathcal{M}_{\text{vac}} , \qquad (25)$$

$$\mathcal{O}_{K} = \left[\overline{s}_{a} \gamma_{\mu} (1 + \gamma_{5}) d_{a}\right] \left[(\overline{s}_{b} \gamma_{\mu} (1 + \gamma_{5}) d_{b}) \right], \qquad (26)$$

$$\mathcal{M}_{\rm vac} = \frac{8}{3} \langle \overline{K}^0 | [\overline{s}_a \gamma_\mu (1 + \gamma_5) d_a] | 0 \rangle \\ \times \langle 0 | [(\overline{s}_b \gamma_\mu (1 + \gamma_5) d_b)] | K^0 \rangle , \qquad (27)$$

where a and b are color indices. The subscript "vac" indicates that this is the result for \mathcal{M}_K in vacuum saturation approximation. The definition can be written in the more familiar form

$$\mathcal{M}_{K} = \frac{16}{3} B_{K} f_{K}^{2} m_{K}^{2} .$$
(28)

The calculation of chiral logarithms is clearer using the definition in Eq. (25). What is required is the logarithms in both \mathcal{M}_K and \mathcal{M}_{vac} .

I begin with the latter. The matrix element \mathcal{M}_{vac} is the product of two matrix elements of the axial current, the corrections to which are discussed in the previous sec-

tion. The result can be read off from Eqs. (8) and (16):

$$\mathcal{M}_{\rm vac} = \frac{16}{3} f^2 p_{\overline{K}} \cdot p_K \{ 1 + 2X [\text{Fig. 1(b)}] + 2X [\text{Fig. 1(c)}] + O(m_K^2) \} .$$
(29)

The $O(m_K^2)$ term comes from nonleading operators, and does not contain chiral logarithms. As in the previous section, I will drop such terms henceforth. In the formula for X [Eq. (17)], one should use the flavor matrix appropriate to the kaon, $T_d = (1/\sqrt{2})(T_6 - iT_7)$. The momenta are both incoming, so that $p_{\overline{K}} \cdot p_K = -p_K^2 = m_K^2$, where m_K is the physical mass including one-loop corrections. The corrections in Eq. (29) are precisely those required to change f^2 to the one-loop corrected f_K^2 :

$$\mathcal{M}_{\rm vac} = \frac{16}{3} f_K^2 m_K^2 \ . \tag{30}$$

The chiral logarithmic corrections to \mathcal{M}_K in full QCD were first calculated in Ref. [5], and related results are given in Ref. [23]. I have checked these calculations, and in the following I present a slight generalization of the results. The $\Delta S = 2$ operator \mathcal{O}_K transforms as part of a (27,1) under SU(3)_L × SU(3)_R. In chiral perturbation theory it is represented by a sum of all operators having the same transformation property, with unknown coefficients. There are no such operators at O(1), and only one at $O(p^2, m_a)$ [5]:

$$\mathcal{O}_K \to \mathcal{O}_K^{\gamma} = \frac{4}{3} B f^4 (\Sigma \partial_{\mu} \Sigma^{\dagger})_{ds} (\Sigma \partial_{\mu} \Sigma^{\dagger})_{ds} , \qquad (31)$$

where the subscripts indicate the *ds*th element of the matrix. I have chosen the coefficients so that, at tree level, $B_K = B$, as can be seen by combining the tree-level result

$$\mathcal{M}_K = \frac{16}{3} B f^2 m_K^2 \tag{32}$$

with that for \mathcal{M}_{vac} from Eq. (29). Chiral perturbation theory does not determine the value of B, which is another parameter like f and μ .

Chiral logarithms in \mathcal{M}_K come from the five types of diagrams shown in Fig. 6. I have represented the operator by two squares, each corresponding to a left-handed current, because it can be written as

$$\mathcal{O}_{K}^{\chi} = \frac{16}{3} B \left(L_{\mu 6} - i L_{\mu 7} \right) \left(L_{\mu 6} - i L_{\mu 7} \right) \,. \tag{33}$$

This factorization, which plays an important role in the calculation, is only true at leading order in the chiral expansion. The result from these diagrams is a multiplicative correction

$$\mathcal{M}_{K} = \frac{16}{3} B f^{2} m_{K}^{2} \{ 1 + X [Fig. 6(a)] + X [Fig. 6(b)] + X [Figs. 6(c) - 6(c)] \} .$$
(34)



FIG. 6. Chiral perturbation theory diagrams for $\mathcal{M}_{\mathcal{K}}$. The pair of squares represents the operator $\mathcal{O}_{\mathcal{K}}^{\chi}$, while the circle is the vertex from the chiral Lagrangian.

Here m_K is the one-loop-corrected kaon mass. Strictly speaking, some of the results from Figs. 6(c)-6(e) are proportional to the tree-level value of m_K^2 (or other tree-level meson masses squared). But since these are correction terms, one is justified in replacing m_K (tree) with m_K (1 loop).

The calculation of Figs. 6(a) and 6(b) is identical to that of Figs. 1(b) and 1(c) respectively, except for an overall factor of 2 since there are two left-handed currents in \mathcal{O}_{k}^{χ} :

$$X[Fig. 6(a)] = 2X[Fig. 1(b)],$$

X[Fig. 6(b)] = 2X[Fig. 1(c)]. (35)

This part of the correction to \mathcal{M}_K is the same as the complete correction to \mathcal{M}_{vac} , Eq. (29), and so cancels in the ratio which defines B_K . In other words, the corrections from Figs. 6(a) and 6(b) are precisely those needed to convert the factor of f^2 in Eq. (34) to the one-loop value f_K^2 , and thus do not affect B_K . This is true for any number of flavors in addition to the s and d quarks.

The only new work is that required to calculate Figs. 6(c)-6(e). The result can be written in a form which is valid for any number of flavors:

$$X[\text{Fig. } 6(c)-6(e)] = I_2(m_K) - \sum_a \left[1 + \frac{m_a^2}{m_K^2} \right] I_1(m_a) [(T_a)_{dd} - (T_a)_{ss}]^2 ,$$
(36)

where I_2 is the standard integral

$$I_{2}(m) = \frac{m^{2}}{f^{2}} \int_{k} G(k,m)^{2}$$
(37)

$$= -L(m) - \left[\frac{m}{4\pi f}\right]^{2} [1 + O(m^{2}/\Lambda^{2})]. \quad (38)$$

This integral results from Figs. 6(c) and 6(d), both of which have a K^0 in the loop. In Fig. 6(e) there are only flavor-diagonal pions in the loop. This diagram gives rise to the second term in X, in which the sum is over flavor-diagonal pions in the mass eigenstate basis.⁹

An important consistency check on this result is that divergent parts can be absorbed into *B*. This requires that there be no quartic divergences, for these would imply that $\mathcal{M}_K \propto \Lambda^4$, while chiral symmetry dictates that the leading term is proportional to m_K^2 . All three diagrams do contain quartic divergences, but they cancel in the sum. There are quadratic divergences in Eq. (36) because $I_1 \propto \Lambda^2$ at leading order. These can be absorbed into *B* as long as they do not depend on the quark masses. The factor of m_a^2/m_K^2 multiplying $I_1(m_a)$ suggests that this may not be true, but in fact the mass dependence cancels because of the result

⁹In fact, the sum can be extended to all pions, because the last factor is nonzero only for flavor-diagonal pions. The same is true in Eq. (40) below.

$$\sum_{a} m_a^2 [(T_a)_{dd} - (T_a)_{ss}]^2 = m_K^2 .$$
(39)

Here again the sum is in the mass eigenstate basis, and only flavor-diagonal states contribute. This result holds for any number of flavors.

Combining Eqs. (30), (34), (35) and (36), the corrected B_K is

$$B_{K} = B \left[1 + I_{2}(m_{K}) - \sum_{a} \left[1 + \frac{m_{a}^{2}}{m_{K}^{2}} \right] I_{1}(m_{a}) [(T_{a})_{dd} - (T_{a})_{ss}]^{2} \right].$$
(40)

In QCD (with $m_u = m_d$) this gives

ſ

$$B_{K} = B \left[1 + I_{2}(m_{K}) - \frac{3}{4} \left[1 + \frac{m_{\eta}^{2}}{m_{K}^{2}} \right] I_{1}(m_{\eta}) - \frac{1}{4} \left[1 + \frac{m_{\pi}^{2}}{m_{K}^{2}} \right] I_{1}(m_{\pi}) \right].$$
(41)

In the limit $m_u = m_d \rightarrow 0$ the chiral logarithmic part is

QUENCHED CHIRAL LOGARITHMS

where the quadratic divergences have been absorbed into B. It is also useful to expand about $m_s = m_d$, in terms of $\delta = (m_s - m_d)^2 / (m_s + m_d)^2$:

$$B_{K} = B[1 - L(m_{K})(3 + \frac{1}{3}\delta) + O(\delta^{2})], \qquad (43)$$

a result previously quoted in Ref. [15]. Finally, for N degenerate flavors one finds

$$B_{K} = B[1 + I_{2}(m) - 2I_{1}(m)]$$

= B[1 - 3L(m) + · · ·]. (44)

Unlike the correction to f_{π} , this result has no explicit dependence on N.

B. B_K in the quenched approximation

I now turn to the calculation of the chiral logarithms in quenched B_K . The conclusion of the following analysis is simple: the result for B_K is unchanged from the full theory except that (1) there are contributions from η' loops when $m_s \neq m_d$, and (2) in the formula Eq. (40) the mass eigenstates are \bar{ss} and \bar{dd} (with masses $m_{ss} = 2\mu m_s$ and $m_{dd} = 2\mu m_d$, respectively). Thus the quenched result is

$$B_{K}(\text{quenched}) = B\left[1 + I_{2}(m_{K}) - \frac{3m_{d} + m_{s}}{2m_{d} + 2m_{s}}I_{1}(m_{dd}) - \frac{m_{d} + 3m_{s}}{2m_{d} + 2m_{s}}I_{1}(m_{ss}) + (m_{s} - m_{d})X(\eta')\right], \qquad (45)$$

where $X(\eta')$ comes from η' loops. In the limit $m_d = m_s$ the result is identical to that in the full theory, Eq. (44).

To begin the quenched calculation, the operator \mathcal{O}_K must be written in terms of pion fields. As before, I assume that the form of the operator is the same as for the full theory, Eq. (31), except that the constant *B* can differ. In other words, I assume that the amplitudes for the operator to create two and four pions are related in the same way as in the full theory. In the continuum, these relations are derived from the chiral Ward identities. With staggered fermions, analogous Ward identities hold in both full and quenched theories [19,15]. With Wilson fermions, the Ward identities are regained in the continuum limit of both full and quenched theories [18]. Dynamical fermions play no role in the derivations.

Given this assumption, the calculation proceeds as in the continuum, except for the usual two complications. First, one must keep only those diagrams in Fig. 6 which correspond to quark diagrams without internal loops. Second, one must determine whether there are extra contributions from η' mesons propagating in the loops.

For Figs. 6(a) and 6(b) the analysis from the preceding section can be used, because the operator \mathcal{O}_{K}^{χ} is a product of two factors of L_{μ} . The calculation of Figs. 6(a) and 6(b) is identical to that of Figs. 1(b) and 1(c). This means that, just as in the full theory, any chiral logarithms from Figs. 6(a) and 6(b) cancel in a ratio that defines B_{K} . In

fact, as shown in the preceding section, the chiral logarithms in quenched f_K differ from those in the full theory. In particular there are no logarithms in f_K for $m_s = m_d$, and there are extra logarithms from η' loops if $m_s \neq m_d$. But these differences are irrelevant for B_K .

Just as in the full theory, then, the logarithms for B_K come entirely from Figs. 6(c)-6(e). The quark diagrams corresponding to these are shown in Fig. 7. The operator \mathcal{O}_K is shown as two squares, each representing a left-handed current. Each diagram actually represents two contractions, one in which the color indices are contracted within each current, the other having the indices contracted between currents. The pairings of quark and antiquark propagators into bilinears can be interchanged by a Fierz transformation. I have chosen the pairing so that each bilinear couples only to one of the mesons in the loop. This makes it easier to see the correspondences with pion diagrams.

None of the diagrams in Fig. 7 contain internal quark loops. Thus all are present in both quenched and full theories. The correspondences between quark and pion diagrams is as follows. Figures 7(a) and 7(b) correspond to Fig. 6(c), since both have a four pion scattering amplitude. Figure 7(c) corresponds to Fig. 6(d), since the particle in the loop is a kaon. And, finally, Figs. 7(d) and 7(e) correspond to Fig. 6(e), in which the mesons in the loop have flavors \bar{ss} and \bar{dd} .

(42)



FIG. 7. Quark diagrams corresponding to Figs. 6(c)-6(e).

Since Figs. 7(a)-7(c) are the same in both full and quenched theories, and since the particles in the loops are flavor off diagonal, the result from the chiral perturbation theory diagrams Figs. 6(c) and 6(d) is identical in full and quenched theories. This is the first example in which there are chiral logarithms in a quenched quantity coming from the same diagrams as those in the full theory. In fact, Fig. 7(b) does not contribute at leading order because the four-pion vertex is $O(p^4)$.

The guenched approximation can, however, affect the evaluation of Fig. 6(e), because the loop contains flavordiagonal mass eigenstates, which, in general, are different in quenched and full QCD. There are also potential η' loop contributions in the quenched theory not present in the full theory, for example that of Fig. 7(e). The effect of these differences is that, in the expression for X [Figs. 6(c)-6(e)], Eq. (36), the sum in the quenched theory runs over both the pions and the η' , while in full QCD the η' is excluded. The difference between these two sums vanishes, however, for degenerate quarks. This is for two reasons. First, the nonsinglet pion mass eigenstates are the same in quenched and full theories, so that the contribution to the sums from these pions agree. Second, the η' is a mass eigenstate, and its extra contribution to quenched sum vanishes, because $T_{ss} - T_{dd} = 0$ for a flavor singlet. Another way of seeing this is to recall that the η' does not couple to L_{μ} , so the extra diagram 7(e) vanishes.

It follows from this discussion that Eq. (45) describes the leading-order chiral logarithms in quenched B_K .

C. B_V and B_A

The matrix element $\mathcal{M}_{\mathcal{K}}$ picks out the positive-parity part of $\mathcal{O}_{\mathcal{K}}$, which consists of a sum of eight different terms: four "vector-vector" $(\gamma_{\mu} \cdot \gamma_{\mu})$ and four "axialvector-axial-vector" $(\gamma_{\mu} \gamma_5 \cdot \gamma_{\mu} \gamma_5)$. There are two different color contractions for each of these eight terms, with one or two traces over color indices. When doing a numerical calculation one calculates all sixteen terms separately, so it is interesting to predict their chiral behavior and finite volume dependence. In this subsection I calculate the leading-order chiral logarithms for the four Lorentz-invariant combinations in which the index μ is summed.

The precise definition of the breakup of \mathcal{M}_K into sixteen pieces is the following.

(1) Write down all contractions (i.e., products of quark

propagators).

(2) Fierz transform the operator such that both external kaons are connected to one or other bilinear. All terms will thus have two traces over spinor indices, and either one or two traces over color indices.

(3) Break up each of the resulting terms into the eight vector-vector and axial-vector-axial-vector pieces. Clearly, this definition applies at the level of contractions. In order to use chiral perturbation theory, however, one must have a definition in terms of matrix elements of operators. For three or less flavors it is not possible to write down operators whose matrix element give each of the sixteen terms separately. As I show shortly, to do this requires at least four valence flavors. In quenched QCD, this presents no problem, for one has already decoupled the number of valence and dynamical flavors. It does mean that the results derived below are not directly applicable to full QCD. They should, however, be as valid for full QCD as they are for quenched QCD.

Consider, then, a theory with $N \ge 4$ light flavors: an s and an s', both of mass m_s ; a d and a d', both of mass m_d ; and N - 4 other quarks with arbitrary (though small) masses. Let the K be the $d\overline{s}$ pion, and the K' be the corresponding state made out of primed quarks. Define \mathcal{M}_K and B_K in this theory exactly as in QCD:

$$M_{K} = \langle \overline{K} | [\overline{s}_{a} \gamma_{\mu} (1 + \gamma_{5}) d_{a}] [(\overline{s}_{b} \gamma_{\mu} (1 + \gamma_{5}) d_{b})] | K^{0} \rangle \qquad (46)$$

$$= \frac{16}{2} B_K f_K^2 m_K^2 \ . \tag{47}$$

Then, by comparing contractions, we can write \mathcal{M}_K as

$$\mathcal{M}_{K} = 2(\mathcal{M}_{V} + \mathcal{M}_{A}) , \qquad (48)$$

$$\mathcal{M}_V = \mathcal{M}_{V1} + \mathcal{M}_{V2} , \qquad (49)$$

$$\mathcal{M}_A = \mathcal{M}_{A1} + \mathcal{M}_{A2} , \qquad (50)$$

$$\mathcal{M}_{V1} = \langle \,\overline{K}' | [\,\overline{s}'_a \gamma_\mu d'_b\,] [\,\overline{s}_b \gamma_\mu d_a\,] | K \,\rangle \,\,, \tag{51}$$

$$\mathcal{M}_{V2} = \langle \,\overline{K}' | [\overline{s}'_a \gamma_\mu d'_a] [\overline{s}_b \gamma_\mu d_b] | K \rangle \,, \tag{52}$$

$$\mathcal{M}_{A1} = \langle \,\overline{K}' | [\overline{s}'_{a} \gamma_{\mu} \gamma_{5} d'_{b}] [\overline{s}_{b} \gamma_{\mu} \gamma_{5} d_{a}] | K \,\rangle \,, \tag{53}$$

$$\mathcal{M}_{A2} = \langle \, \overline{K} \, ' | [\overline{s} \, '_a \gamma_\mu \gamma_5 d'_a \,] [\overline{s}_b \gamma_\mu \gamma_5 d_b \,] | K \, \rangle \ . \tag{54}$$

The point of the extra flavors is to restrict contractions; each of the matrix elements in the last four lines has only one contraction. The factor of 2 on the first line appears because the two bilinears in \mathcal{M}_K are identical, and so there are twice as many contractions as in \mathcal{M}_V and \mathcal{M}_A . The subscripts 1 and 2 indicate the number of traces over color indices.

The corresponding B parameters are defined by

$$\mathcal{M}_{V1} = \frac{8}{3} f_K^2 m_K^2 B_{V1}, \quad \mathcal{M}_{V2} = \frac{8}{3} f_K^2 m_K^2 B_{V2} , \quad (55)$$

and identical equations with $V \rightarrow A$. These definitions are such that

$$B_{K} = B_{V} + B_{A} = B_{V1} + B_{V2} + B_{A1} + B_{A2} .$$
 (56)

It is straightforward to show that in vacuum saturation approximation $B_{A1} = 0.25$, $B_{A2} = 0.75$, and thus $B_A = 1$, while $B_V = B_{V1} = B_{V2} = 0$. The numerical results, shown

in Sec. VII, do not, in fact, agree with these predictions.

I calculate the chiral logarithms in these B parameters in the full theory, and then argue that the results apply with only minor changes in the quenched approximation. The first step is to transcribe the problem into the language of chiral perturbation theory. As explained in Sec. II, the chiral Lagrangian itself takes the same form for any number of flavors. To transcribe the operators, however, is more complicated. It turns out that, at the order I am working, one needs the set

$$\mathcal{O}_1 = f^4 (\Sigma \partial_\mu \Sigma^\dagger)_{d's'} (\Sigma \partial_\mu \Sigma^\dagger)_{ds} , \qquad (57)$$

$$\mathcal{O}_2 = f^4 (\Sigma \partial_\mu \Sigma^\dagger)_{d's} (\Sigma \partial_\mu \Sigma^\dagger)_{ds'} , \qquad (58)$$

$$\mathcal{O}_3 = -2\mu^2 f^4 \Sigma_{d's} \Sigma_{ds'}^{\dagger} , \qquad (59)$$

$$\mathcal{O}_4 = f^4(\partial_\mu \Sigma_{d's})(\partial_\mu \Sigma_{ds'}^\dagger) , \qquad (60)$$

$$\mathcal{O}_{5} = f^{4} (\Sigma \partial_{\mu} \Sigma^{\dagger})_{d's'} (\Sigma^{\dagger} \partial_{\mu} \Sigma)_{ds} .$$
(61)

The overall factors are chosen for later convenience.

To transcribe the quark operators, they must be written in terms of representations of the $SU(N)_L \times SU(N)_R$ chiral group. The first step is to rewrite the operators in terms of left- and right-handed bilinears

$$[V'V]_{i} = \frac{1}{2}([L'L]_{i} + [L'R]_{i}) + \mathcal{O}^{-},$$

$$[A'A]_{i} = \frac{1}{2}([L'L]_{i} - [L'R]_{i}) + \mathcal{O}'^{-},$$
(62)

where i = 1, 2 labels the number of color traces. The negative-parity operators \mathcal{O}^- and \mathcal{O}'^- do not contribute to the matrix elements under study and can be ignored. I am using the compact notation

$$[V'V]_1 = [\overline{s}'_a \gamma_\mu d'_b] [\overline{s}_b \gamma_\mu d_a] , \qquad (63)$$

$$[A'A]_1 = [\overline{s}'_a \gamma_\mu \gamma_5 d'_b] [\overline{s}_b \gamma_\mu \gamma_5 d_a], \qquad (64)$$

$$[L'L]_{1} = [\bar{s}'_{a} \gamma_{\mu} (1 + \gamma_{5}) d'_{b}] [\bar{s}_{b} \gamma_{\mu} (1 + \gamma_{5}) d_{a}], \qquad (65)$$

$$[L'R]_1 = [\overline{s}'_a \gamma_\mu (1+\gamma_5) d'_b] [\overline{s}_b \gamma_\mu (1-\gamma_5) d_a], \qquad (66)$$

with similar definitions for operators with two color traces.

The left-handed bilinears L and L' transform as $[(N^2-1), 1]$, i.e., as adjoints under $SU(N)_L$ and singlets under $SU(N)_R$. Under $SU(N)_L, [L'L]$ lies in the symmetric product of two adjoints,

$$[(\mathbf{N}^{2}-1)\otimes(\mathbf{N}^{2}-1)]_{\text{symm}} = 1 + (\mathbf{N}^{2}-1) + \mathcal{S} + \mathcal{A} .$$
 (67)

The representations \mathscr{S} and \mathscr{A} are both traceless, and are respectively symmetric and antisymmetric under interchange of quark indices. They have dimensions $(N-1)N^2(N+3)/4$ and $(N-3)N^2(N+1)/4$ respectively. Since L' and L have no flavors in common, their product must be a combination of \mathscr{S} and \mathscr{A} . The combinations belonging to one or other representation are

$$[L'L]_{2} \pm [L'L]_{1} = [\overline{s}'_{a}\gamma_{\mu}(1+\gamma_{5})d'_{a}][\overline{s}_{b}\gamma_{\mu}(1+\gamma_{5})d_{b}] \pm [\overline{s}'_{a}\gamma_{\mu}(1+\gamma_{5})d_{a}][\overline{s}_{b}\gamma_{\mu}(1+\gamma_{5})d'_{b}] \in \begin{cases} \mathscr{S} \\ \mathscr{A} \end{cases}$$
(68)

In chiral perturbation theory the lowest dimension operators transforming in these representations are

$$(S,1): O_S = O_1 + O_2,$$

$$(\mathcal{A},\mathbf{1}): \mathcal{O}_{\mathcal{A}} = \mathcal{O}_1 - \mathcal{O}_2$$

Thus the transcription of the continuum operators is

$$[L'L]_{2} + [L'L]_{1} = \alpha_{S} \mathcal{O}_{S} + O(p^{4}) ,$$

$$[L'L]_{2} - [L'L]_{1} = \alpha_{A} \mathcal{O}_{A} + O(p^{4}) ,$$
(69)
(70)

where α_S and α_A are new, unknown coefficients. One can estimate their values by taking matrix elements between a K and a \overline{K}' , using the vacuum saturation approximation for the left-hand side (LHS), and tree-level chiral perturbation theory for the RHS. The resulting estimates are $\alpha_S = \frac{4}{3}$ and $\alpha_A = \frac{2}{3}$. The same results are obtained using external $\overline{d}'s$ and $\overline{s}'d$ kaons.

The [L'R] operators transform as $[(N^2-1), (N^2-1)]$. They can thus be written as linear combinations of the three operators \mathcal{O}_{3-5} , which also transform in this way

$$[L'R]_i = \beta_i \mathcal{O}_3 + \gamma_i \mathcal{O}_4 + \delta_i \mathcal{O}_5 + O(p^4) .$$
⁽⁷¹⁾

There are six unknown coefficients, three for each color contraction. To estimate their size I first take the matrix element between a K and a \overline{K}' . At tree level, only \mathcal{O}_5 has a nonvanishing matrix element. Using vacuum saturation for the matrix element of $[L'R]_i$ yields the estimates $\delta_1 = \frac{1}{3}$, $\delta_2 = 1$.

To estimate the β_i , I take the matrix element between \overline{d} 's and \overline{s} 'd kaons. In chiral perturbation theory, one finds, at tree level,

$$\langle K(\overline{d}'s)|[L'R]_i|K(\overline{s}'d)\rangle = -4\beta_i\mu^2 f^2 + O(\gamma_i m_K^2).$$
⁽⁷²⁾

To use the vacuum saturation approximation I first Fierz transform the operators, e.g.,

$$[L'R]_{1} = -2[\bar{s}'_{a}(1-\gamma_{5})d_{a}][\bar{s}_{b}(1+\gamma_{5})d'_{b}], \qquad (73)$$

and then use

$$\langle 0|\overline{s}_{b}\gamma_{5}d_{b}'|K(\overline{s}'d)\rangle = i\mu\sqrt{2}f_{K} , \qquad (74)$$

to give

$$\langle K(\overline{d}'s)|[L'R]_1|K(\overline{s}'d)\rangle_{\rm vac} = -4\mu^2 f^2 , \qquad (75)$$

$$\langle K(\vec{d}'s)|[L'R]_2|K(\vec{s}'d)\rangle_{\rm vac} = -\frac{4}{3}\mu^2 f^2 .$$
⁽⁷⁶⁾

The resulting estimates are $\beta_1 = 1$ and $\beta_2 = \frac{1}{3}$. This method gives no estimate for the γ_i .

In summary, each quark operator can be written as $\sum_j c_j O_j$ with corrections of $O(p^4)$. The coefficients are collected in Table II, where I use the definitions

$$\alpha_{\pm} = \frac{1}{2} (\alpha_S \pm \alpha_A) . \tag{77}$$

In vacuum saturation approximation these are $\alpha_{+} = 1$ and $\alpha_{-} = \frac{1}{3}$.

All four operators in Eqs. (51)-(54) contain "left-left" and "left-right" parts. Equation (72) shows that the matrix elements of "left-right" operators are generically of order 1, unlike those of "left-left" operators which are of $O(p^2)$. Nevertheless, at tree level, and for the external flavors K and K', all four operators have matrix elements of $O(p^2)$. The choice of external flavors has removed the order-1 contributions. As we will see, however, these contributions do have an effect when loops are included.

The calculation of the chiral logarithms parallels that for B_K . The same diagrams contribute and the general result is

$$\langle \overline{K'^{0}} | c_{j} \mathcal{O}_{j} | K^{0} \rangle = (c_{1} - c_{5}) 2m_{K}^{2} f_{K}^{2} + c_{2} f^{2} [2m_{K}^{2} I_{2}(m_{K}) - (m_{K}^{2} + m_{ss}^{2}) I_{1}(m_{ss}) - (m_{K}^{2} + m_{dd}^{2}) I_{1}(m_{dd})] + c_{3} 2\mu^{2} f^{2} [I_{1}(m_{ss}) + I_{1}(m_{dd}) - 2I_{1}(m_{K}) - 2I_{2}(m_{K})] + c_{4} f^{2} [4m_{K}^{2} I_{1}(m_{K}) - 2m_{K}^{2} I_{2}(m_{K}) - (m_{K}^{2} + m_{ss}^{2}) I_{1}(m_{ss}) - (m_{K}^{2} + m_{dd}^{2}) I_{1}(m_{dd})] .$$
(78)

I have simplified the result using $m_{s'}=m_s$ and $m_{d'}=m_d$. Only the operators \mathcal{O}_1 and \mathcal{O}_5 contribute to Figs. 6(a) and 6(b), and the corrections simply convert f^2 into the 1-loop result f_K^2 . The general expression for f_K at 1-loop is given in Sec. III. These are the only contributions of \mathcal{O}_1 and \mathcal{O}_5 . The remaining three operators contribute only to Figs. 6(c)-6(e). The pions in the loops in these diagrams always consist of a primed quark and an unprimed antiquark (or vice versa). These states are automatically mass eigenstates, for any number of flavors, so the result is valid for all $N \ge 4$.

There are no quartic divergences in Eq. (78). The only quadratic divergences at this order are proportional to $c_2 m_K^2 \Lambda^2$. These can be absorbed into the coefficient c_1 , and correspond to mixing between \mathcal{O}_1 and \mathcal{O}_2 . The quadratic divergences proportional to c_3 cancel; this is essential, because they are not proportional to m_K^2 , and could not be absorbed by a redefinition of the coefficients. The quadratic divergences proportional to c_4 cancel because $2m_K^2 = m_{ss}^2 + m_{dd}^2$. This is not required by chiral symmetry, however, as a divergence could have been absorbed into c_5 . Finally, there are also quadratic divergences hidden in f_K^2 , which can be absorbed into f, as discussed previously.

I now collect the results for the *B* parameters, using the general formula Eq. (78), the decompositions given in Table II, and the definitions Eqs. (48)–(55). I present the results for $m_s = m_d$, since it turns out that these also apply in the quenched approximation. The generalizations to nondegenerate quarks are easy to obtain from the above. For the vector B parameters I find

$$B_{V1} = \frac{3}{8}(\alpha_{-} - \delta_{1}) + \frac{3}{8}\alpha_{+}(I_{2} - 2I_{1}) - \frac{3}{4}\beta_{1}\frac{\mu^{2}}{m_{K}^{2}}I_{2} - \frac{3}{8}\gamma_{1}I_{2} ,$$
(79)

$$B_{V2} = \frac{3}{8}(\alpha_{+} - \delta_{2}) + \frac{3}{8}\alpha_{-}(I_{2} - 2I_{1}) - \frac{3}{4}\beta_{2}\frac{\mu^{2}}{m_{K}^{2}}I_{2} - \frac{3}{8}\gamma_{2}I_{2} .$$
(80)

Since $m_K = m_{ss} = m_{dd}$, the arguments of the I_i are all the same, and I have dropped them. The results for B_A are the same except that terms proportional to β_i , γ_i , and δ_i have the opposite sign. A check on these results is that they combine to give the same result for $B_K = B_{V1} + B_{V2} + B_{A1} + B_{A2}$ as found previously [Eq. (44)],

TABLE II. Decomposition of quark operators into chiral operators $\sum_j c_j \mathcal{O}_j$.

Operator	<i>c</i> ₁	<i>c</i> ₂	<i>c</i> ₃	C4	c 5
$2[V'V]_1$	lpha	α_+	$oldsymbol{eta}_1$	γ ₁	δ_1
$2[V'V]_2$	α_+	α_{-}	β_2	γ_2	δ_2
$2[A'A]_1$	α	$lpha_+$	$-oldsymbol{eta}_1$	$-\gamma_1$	$-\delta_1$
$\frac{2[A'A]_2}{2}$	α_+	α	$-\beta_2$	$-\gamma_2$	$-\delta_2$

$$B_K = \frac{3}{4} \alpha_S (1 + I_2 - 2I_1) , \qquad (81)$$

providing one identifies $B = \frac{3}{4}\alpha_S$.

The corrections to B_V and B_A include a term of the form I_2/m_K^2 which is proportional to $\ln(m_K)$. This is enhanced by a factor of μ^2/m_K^2 over the usual chiral logaarithms. The general possibility of enhanced chiral logarithms was noticed long ago by Pagels and Langacker [4]. The enhanced logarithm here is the remnant of the fact that \mathcal{O}_3 is of order 1 in the chiral expansion, rather than of $O(p^2)$. For small enough m_K the enhanced logarithm dominates over the "leading-order" terms, and it diverges in the chiral limit. The vector-vector and axialvector-axial-vector matrix elements do not themselves diverge, but contain a chiral logarithm proportional to $m_K^2 \ln(m_K)$, rather than the usual $m_K^4 \ln(m_K)$. The unknown scale in the logarithm can be absorbed into δ_i , and corresponds to mixing between \mathcal{O}_3 and \mathcal{O}_5 .

These results can be adapted to the quenched approximation with the by now standard procedure. The reasoning is identical to that for B_K , because the quark diagrams are the same. The only diagrams containing internal quark loops are those that change f^2 to f_K^2 , and these cancel in B parameters. The other diagrams, those of Fig. 7, are included in the guenched approximation. There are only two possible differences between quenched and full theories. The first come from the fact that the mass eigenstates are not the same for flavor diagonal states. This is not relevant here, however, since all pions in the loops are flavor off diagonal. The second comes from η' loops, which do not contribute to the full theory, but might be present in the quenched approximation. The quark diagram is that shown in Fig. 7(e). As for B_K , however, this diagram does not contribute if $m_s = m_d$. For \mathcal{O}_1 , \mathcal{O}_2 , and \mathcal{O}_5 , this is because the η' does not couple to these operators. For \mathcal{O}_3 and \mathcal{O}_4 , it is because the η' , although it couples to the operators, cannot appear in a loop such as in Fig. 7(a) because the flavor indices are wrong. The η' does appear in two loop diagrams. The conclusion is that, for $m_s = m_d$, the results of Eqs. (78)-(80) apply without change to the quenched approximation. For $m_s \neq m_d$, however, there will be additional contributions from Fig. 7(e).

There is one complication in the numerical calculation of the B parameters which arises if one uses staggered fermions. I describe this briefly here; for more details see, for example, Refs. [19,15]. Each lattice staggered fermion represents four continuum flavors, so the quark fields have an additional flavor index. The bilinears that are actually used have these extra indices contracted as in

$$\overline{s}\gamma_{\mu}d \rightarrow \overline{s}_{1}\gamma_{\mu}d_{1} + \overline{s}_{2}\gamma_{\mu}d_{2} - \overline{s}_{3}\gamma_{\mu}d_{3} - \overline{s}_{4}\gamma_{\mu}d_{4} , \qquad (82)$$

where the color indices are implicit. This particular flavor is that of the lattice Goldstone pion, and in numerical calculations the external kaons have this flavor. The analysis given above is valid for any number of flavors, and thus applies directly to these generalized operators. One simply has to break each matrix element up into parts, each of which is like those in Eqs. (51)-(54). Keeping track of flavor factors, one finds that the form of the final result is unchanged. The only difference is that most of the pions in the loops have flavors such as \overline{s}'_1d_2 , and thus are not Goldstone pions. In fact, it is straightforward to show that only one out of sixteen pions is a Goldstone pion. This is important at finite lattice spacing, because non-Goldstone pions are slightly more massive than the Goldstone pion, and one must use the appropriate mass in the loop integrals. In addition, the vertices of the non-Goldstone pions will differ from those of the Goldstone pions by terms of O(a). The effect of this is to make the replacement

$$I_2(m_K)/m_K^2 \rightarrow \frac{1}{16}I_2(m_G)/m_G^2 + \frac{15}{16}R_{\rm NG}I_2(m_{\rm NG})/m_{\rm NG}^2$$
,
(83)

where $m_{\rm G}$ and $m_{\rm NG}$ are, respectively, the mass of the Goldstone and non-Goldstone pions, and $R_{\rm NG}$ is the ratio of non-Goldstone to Goldstone coefficients. (I am assuming here that all the non-Goldstone pions are degenerate, which, numerically, is a good approximation.) Similar replacements should be made in all the loop integrals in Eqs. (79) and (80).

V. η' LOOPS

The major new feature of quenched chiral perturbation theory is the appearance of the light η' in loops. Thus far I have considered only quantities which do not have η' contributions at one-loop order. Clearly it would be preferable to understand such contributions, since this would increase the number of quantities whose finite volume dependence could be predicted. In addition, it is potentially dangerous to leave an unsettled problem in one part of the theory, as it may feed back at higher loop order into the quantities I have studied.

I consider the effects of η' loops in theory with N degenerate valence quarks. Using degenerate quarks simplifies the discussion in two ways. First, the η' can be chosen as one of the mass eigenstates, which allows the effects of η' loops to be clearly distinguished. Second, as discussed above, the η' only couples to the pions through the mass term in the chiral Lagrangian, \mathcal{L}_m . This means, in particular, that η' loops only affect mass renormalization, and I focus on this in the following discussion.

In the quenched approximation, the only diagrams that renormalize the pion mass at leading order are those of Figs. 3(d)-3(f). Note that only the η' appears in these loops, since nonsinglet states cannot annihilate into gluons. As discussed at the end of Sec. II, the contribution from Fig. 3(f), which arises from the vertex $V'_2(0)\eta' \text{Tr}(M\eta'\pi^2)$, can be rotated away by a field redefinition. Its effect appears in a change in the value of the constant A in $\mathcal{L}_{\eta'}$ [Eq. (14)]. Thus I need only consider Figs. 3(d) and 3(e). These are represented in chiral perturbation theory by Fig. 8(a), where the cross is the gluon annihilation vertex, $\mathcal{L}_{\eta'}$. The renormalized mass at one loop from Fig. 8(a) is

$$m_{\pi}^2 = 2\mu m \{1 + X[\text{Fig. 8(a)}]\}$$
 (84)



FIG. 8. Diagrams contributing to mass renormalization in the quenched approximation. The circle represents the vertex from the mass term in the chiral Lagrangian, the cross an insertion of the gluonic amplitude $\mathcal{L}_{\eta'}$. Lines with a cross are η' propagators; the others can be either pion or η' propagators.

$$X[\text{Fig. 8(a)}] = \frac{1}{Nf^2} \int_k G(k, m_{\eta'})^2 [Am_0^2 + (A-1)k^2]$$

$$= \frac{1}{N} \left[A \frac{m_0^2}{m_{\eta'}^2} - (A-1) \right] I_2(m_{\eta'})$$

$$+ \frac{1}{N} (A-1) I_1(m_{\eta'}) . \qquad (86)$$

The mass in the loop is m'_{η} , which equals m_{π} at tree level. This result agrees with Ref. [11]. The k^2 term in the integrand leads to a correction of the usual type: it is proportional to $m^2_{\eta'}\ln(m_{\eta'})$, and vanishes in the chiral limit. The m^2_0 term, by contrast, leads to a correction proportional to $m^2_0\ln(m_{\eta'})$ which diverges in the chiral limit. This enhanced chiral logarithm is similar to those appearing in B_V and B_A . The additional dimensionful parameter m_0 violates the standard power counting rules of chiral perturbation theory, according to which each loop comes with a factor of m^2_{π} , possibly multiplied by logarithms. For small enough m_{π} , the one-loop correction proportional to $m^2_0 \ln(m_{\eta'})$ is not small, and one must sum all such terms.

In practice, as stressed in Ref. [11], this summation may not be numerically important. This is because in present lattice calculations the pions are not very light, $m_{\pi} \ge 400$ MeV. To estimate the size of the correction, I extract the leading logarithm

$$X[\text{Fig. } 8(a)] = -\delta \ln \left[\frac{m_{\eta'}^2}{\Lambda^2} \right].$$
(87)

Here I have introduced the parameter

$$\delta = \frac{Am_0^2}{N(4\pi f)^2} \approx 0.2 , \qquad (88)$$

where the estimate uses N=3, A=1, and $m_0=900$ MeV. Taking $\Lambda \approx m_p$, the magnitude of the logarithm is roughly unity for $m_{\pi}=400$ MeV. Thus the correction increases m_{π}^2 by only 20%, and resummation is not required. Since the logarithm is not large, however, it does not dominate over constants and power corrections, and the full expression of Eq. (86) should be used.

Bernard and Golterman have begun a program of calculations, in which they include η' loops only at one-loop order, for the general case of nondegenerate quarks [11]. Their predictions can be compared to numerical results for various quantities. My interest in this section is not phenomenological, but rather to understand the apparent divergence of the correction in the chiral limit. To address this question one must resum the leading logarithms of the form $m_0^{2n} \ln(m_{\eta'})^n / f^{2n}$. This is relatively straightforward, because the leading logarithms come only from diagrams in which $\mathcal{L}_{\eta'}$ appears on a propagator both ends of which come from the same vertex. An example is shown in Fig. 8(b). Diagrams involving η' propagators connecting two or more vertices have fewer logarithms. For example, the diagram of Fig. 8(c) gives a correction proportional to $m_0^4 m_{\eta'}^2 / f^6$, with no logarithms at all.

The bubbles exemplified by Figs. 8(a) and 8(b) sum up to an exponential

$$m_{\pi}^2 = 2\mu m \exp\{X[\text{Fig. 8(a)}]\},$$
 (89)

the leading-logarithmic part of which is

$$m_{\pi}^{2} = 2\mu m \left[\frac{\Lambda^{2}}{m_{\eta'}^{2}} \right]^{\delta} .$$
⁽⁹⁰⁾

It is straightforward to show that the same exponential factor corrects any vertex coming from \mathcal{L}_m , so that the entire mass term is replaced by

$$\mathcal{L}'_{m} = \left[\frac{\Lambda^{2}}{m_{\eta'}^{2}}\right]^{\delta} \mathcal{L}_{m} , \qquad (91)$$

with the proviso that no η' fields coming from this vertex be joined together by an insertion of $\mathcal{L}_{\eta'}$. One corollary of this result is that the η' mass itself is corrected in the same way as m_{π} , so that they remain equal. This means that Eq. (90) should be interpreted as a self-consistent equation for the common pion and η' mass, with $m_{\eta'}$ on the right-hand side replaced by m_{π} . Solving this equation yields the final result

$$m_{\pi}^{2} = (2\mu m)^{1/(1+\delta)} (\Lambda^{2})^{\delta/(1+\delta)} .$$
(92)

Solving Eq. (90) self-consistently corresponds to summing all the diagrams of the form Fig. 9(a), in which each internal η' propagator is decorated with any number of bubble sums. In terms of quarks, these "cactus" diagrams look like Fig. 9(b).

The result Eq. (92) seems very odd at first sight. For one thing, m_{π}^2 has a nonanalytic dependence on the quark mass. For another, it appears to contradict the strong numerical evidence that $m_{\pi}^2 \propto m$ in quenched as well as full theories [24]. To investigate further, I substitute Eq. (92) back into Eq. (91) to give [using also Eq. (13)]



FIG. 9. An example of a diagram corresponding to selfconsistent leading-logarithm evaluation of m_{π} in the quenched approximation: (a) in chiral perturbation theory; (b) in terms of quarks. Notation as in Fig. 8.

$$\mathcal{L}'_{m} = -\frac{f^{2}}{4} (2\mu m)^{1/(1+\delta)} (\Lambda^{2})^{\delta/(1+\delta)} \{\cos(\phi_{0}) \operatorname{Tr}[\Sigma + \Sigma^{\dagger}] + i \sin(\phi_{0}) \operatorname{Tr}[\Sigma - \Sigma^{\dagger}] \} .$$
⁽⁹³⁾

This says that, at leading-logarithmic order, the effect of η' loops amounts to a change in the coefficient multiplying the mass term. The quenched chiral Lagrangian looks the same as that in the full theory¹⁰ as long as one uses a new mass m':

$$m' = m \left[\frac{\Lambda^2}{2\mu m} \right]^{\delta(1+\delta)}.$$
 (94)

Looked at this way, the result is less strange. All the relationships between amplitudes involving different numbers of pions remain valid, e.g., the Weinberg scattering length formulas. These relations involve only the physical parameters f_{π} and m_{π} , and not the quark masses. In particular, there is no contradiction with exact lattice Ward identities that can be derived with staggered fermions [19,20].

Associated with the enhanced chiral logarithms is an "enhanced" finite volume dependence. The precise form of this dependence is discussed in Sec. VII. The important point for present purposes is that it too is as a function of physical quantities alone, and not of m (or m'). Thus one cannot use the finite volume behavior to observe the non-analytic dependence on m.

Is there any way of observing this dependence? The remaining hope rests with the numerical lattice results for the quark mass dependence of m_{π}^2 and the condensate $\langle \bar{\psi}\psi \rangle$. The predicted dependence of m_{π}^2 on *m* is given in Eq. (92). That of the condensate may be deduced from Eq. (93) (or from the formula of Gell-Mann, Oakes, and Renner):

$$\langle \bar{\psi}\psi \rangle \propto m^{-\delta/(1+\delta)}$$
 (95)

These predictions are in apparent conflict with numerical simulations which find that $m_{\pi}^2 \propto m_q$ and $\langle \bar{\psi}\psi \rangle = \langle \bar{\psi}\psi \rangle_0 + O(m_q)$, where m_q is the bare lattice quark mass [24]. The results are probably good enough to rule out $\delta \approx 0.2$, assuming that we identify m_q with m.

In full QCD this identification is needed for the following reason. The bare quark mass m_q is part of a scalar source s. Functional derivatives with respect to s yield correlators involving scalar bilinears in the presence of the source. Setting the source to zero we obtain the correlators in the massless theory, the condensate being the simplest example. We expect these correlators to exist in a regulated theory. For this to be so, the chiral Lagrangian must contain only positive integral powers of s. This means that, up to an overall constant, and up to corrections of $O(m^2)$, we should identify m_q with m.

In quenched QCD, by contrast, one does not have a Lagrangian framework, and I see no reason why the con-

nection between m and m_q should be analytic. I propose that m_q should be identified with m' (up to an overall factor), rather than with m. This would imply the relationship

$$m_a \propto m^{1+\delta} . \tag{96}$$

As far as I can tell, once this identification is made there are no observable effects of the nonanalytic dependence on m.

VI. FINITE VOLUME DEPENDENCE

The discussion thus far concerns quantities calculated in infinite volume. Numerical simulations, by contrast, are done in finite boxes. These have finite extent in both space and in Euclidean "time." I denote the length of the lattice in the spatial directions by L_i , i=1-3, and the length in the time direction by L_4 . The latter is related to the temperature by $T=1/L_4$. In the following discussion, I assume that $L_4 \gg L_i$, so that the dominant finitesize effect is due to the spatial volume.

Gasser and Leutwyler have shown that the same loop graphs which give rise to the leading chiral logarithms also allow one to predict the leading finite volume dependence [6]. They have presented results for f_{π} and m_{π} in a theory with N degenerate flavors. It is simple to extend their method to any quantity for which the chiral logarithms are known. The only subtlety concerns the applicability of their method to the quenched approximation.

Their method depends upon the result that the volume dependence in chiral perturbation theory enters only through the pion propagators. The chiral Lagrangian itself does not depend on the volume, and so the vertices in chiral perturbation theory are unchanged. This is also true of external sources, such as that representing \mathcal{O}_K . Thus the constants f, μ , and B are volume independent. The volume dependence of the pion propagator enters through the boundary conditions, which are periodic in all four Euclidean directions. This means that the integral over momenta is replaced by a sum, in which p_{μ} is an integer multiple of $2\pi/L_{\mu}$. In position space, the propagator is a sum over periodic images

$$G_L(x,m) = G(x,m) + \sum_{n_\mu \neq 0} G(x + n_\mu L_\mu, m) , \qquad (97)$$

where $n_{\mu} = (1,0,0,0)$, (-1,0,0,0), (0,1,0,0), etc. Volume dependence comes from replacing G with G_L in loop integrals such as I_1 and I_2 .

The volume independence of the chiral Lagrangian is clearly crucial to the method. The arguments for this are presented in detail in Ref. [25]. The ingredients are¹¹ (a)

¹⁰Of course, in the application of this Lagrangian, one must follow the rules discussed in Secs. III and IV for discarding diagrams containing internal quark loops.

¹¹In addition, one must work at volumes large enough that chiral symmetry is not restored.

that chiral Ward identities remain valid in finite volume, (b) causality, (c) that the pion propagator at fixed distance should approach the infinite volume form as the volume increases to infinity, and (d) the permutation symmetry between the four directions $\mu = 1-4$. The first three ingredients establish the existence of a finite volume chiral Lagrangian, and that the coefficients of its leading terms are the same as at infinite volume. Volume dependence suppressed by powers of $1/L^2$ is allowed, so there can be terms such as $\sum_i \text{Tr}(\partial_i \Sigma \partial_i \Sigma^{\dagger})/L_i^2$. Such terms are ruled out by the fourth ingredient, permutation symmetry. This symmetry implies that, if there is any dependence on the spatial lengths L_i , there must also be a dependence on L_4 , and thus on T. But this contradicts a standard field theoretic result, namely that the temperature enters the functional integral representation of the partition function only through the boundary conditions. Thus permutation symmetry rules out L_i dependence in the vertices at any order in the chiral expansion. For the symmetry to be present, the boundary conditions must be the same in all four directions. In particular, since the finite temperature functional integral has antiperiodic boundary conditions on fermions in the time direction, the boundary conditions must also be antiperiodic in the spatial directions.

In the quenched approximation only the first and third ingredients are present. Concerning the former, one can derive the Ward identities on the lattice independent of the guenched approximation [18,19]. As for the latter, it is reasonable to assume that the boundaries become unimportant as they recede to infinity. But the lack of dynamical guarks means that one cannot even formulate the second ingredient, causality. Causality requires that the anticommutator of fermion fields vanish outside the forward light cone. In the quenched approximation, one cannot construct the transfer matrix and obtain an operator formulation in which quarks are represented by anticommuting operators. A less formal way of seeing this is to note that, without quark loops, there is no Pauli exclusion principle, so that, in some sense, fermion fields no longer anticommute. The lack of quark loops also means that the fourth ingredient, permutation symmetry, is not relevant. The requirement of antiperiodic boundary conditions applies to internal quark loops, for it is these loops which build up the contribution of states containing quarks to the partition function. Since these loops are absent, one is free to choose the boundary conditions on the valence quarks as one pleases.

The absence of two of the four ingredients means that one cannot rigorously argue for the volume independence of the chiral Lagrangian in the quenched approximation. In fact, as already discussed, even the existence of such a Lagrangian is an assumption. I will simply extend this assumption by taking the quenched chiral Lagrangian to have volume-independent coefficients. I think this is reasonable, and in the following I give a qualitative argument in its support. Ultimately, however, it must be tested by numerical simulations.

The physical origin of the finite size effects from chiral loops is that particles are surrounded by a cloud of virtual pions. The pions in this cloud can propagate "around the world" due to the periodicity, the propagator falling roughly like $\exp(-m_{\pi}L_i)$. If there were no interactions, then there would be no pion cloud, and the only effect of finite volume would be to restrict the momenta to discrete values. In particular, masses would remain unchanged. Lüscher has given a general analysis of finite volume effects arising from polarization clouds on the masses of stable particles [7]. For the pion, his analysis implies that the leading-order effect comes from a "tadpole" diagram, similar to that of Fig. 1(b), except that the vertex and propagator are fully dressed. The crucial result is that volume dependence enters only through the pion propagator, which goes "around the world" in each of the six spatial directions. Lüscher's general formula simplifies when applied to pions, because one can do a chiral expansion of the propagator and vertices.¹² The result coincides with that obtained from chiral loops, up to exponentially small corrections which are beyond the accuracy of Lüscher's formula [6]. This agreement would not hold if there were any volume dependence in the coefficients of the chiral Lagrangian, such as terms of $O(1/L^2)$. This provides independent confirmation of the arguments of Ref. [25].

Lüscher's result can be paraphrased as follows. The pion only "knows" about the finite size of the box through loop diagrams, and in particular through the pion propagating in the loop. The vertex that emits and absorbs the pion does not itself "know" about the finite volume. Such knowledge would require another pion loop, which would bring an additional exponential suppression. This physical picture applies also to properties of the pion other than its mass, for example f_{π} and B_{K} . It also seems reasonable to apply it to the quenched approximation.

The previous discussion has ignored the possibility of finite volume effects due to particles being "squeezed" by the box. To avoid such squeezing the box must be larger than diameter of the particle. It is not very clear what box size this requires. Numerical calculations of various definitions of "wave functions" show that they fall approximately exponentially in the range 0.5-1.5 fm, with a scale of roughly $m_{\rm wf} = 750$ MeV [26,27]. This may imply a correction factor from squeezing that falls as $\exp(-m_{\rm wf}L)$. As long as $m_{\pi} < m_{\rm wf}$, the contribution from squeezing should be suppressed. At large enough distances the wave function should fall faster than an exponential, so the effect of squeezing should be further suppressed.

Another source of finite volume dependence is loops of heavier mesons, such as the ρ . Such contributions would fall roughly as $\exp(-m_{\rho}L)$. As long as $m_{\pi} \ll m_{\rho}$, they

¹²The reader consulting Ref. [7] might be misled by the fact that the leading term in Lüscher's formula falls like $exp(-\sqrt{3/4}mL)$, i.e., more rapidly than exp(-mL). This term, however, is absent for pions because it is proportional to the three pion coupling, which is zero. The dominant term falls as exp(-mL), and the leading corrections are proportional to $exp(-\sqrt{2}mL)$.

should be strongly suppressed relative to the contributions from pion loops.

I now turn to the calculation of finite volume dependence. The loop integrals that appear are I_1 and I_2 , proportional respectively to $\int_k G$ and $\int_k G^2$. They are related by

$$I_2(m) = -m^2 \frac{d}{dm^2} I_1(m) , \qquad (98)$$

so it is sufficient to consider I_1 . It is simplest to work in position space

$$I_1(m) = \frac{1}{f^2} \int_{|k| < \Lambda} G(k,m) = \frac{1}{f^2} G_{\Lambda}(x=0,m) .$$
 (99)

The subscript Λ indicates that the position space propagator is smeared over a range $1/\Lambda$, due to the cutoff on the momentum integral. At finite volume we simply substitute G_L [Eq. (97)] for G. We are interested in the difference

$$I_1(m,L) = I_1(m,L) - I_1(m) , \qquad (100)$$

which is given by

$$f^{2}\overline{I}_{1}(m,L) = \sum_{n_{\mu}\neq 0} G_{\Lambda}(n_{\mu}L_{\mu},m)$$
(101)

$$\approx \sum_{n_{\mu} \neq 0} G(n_{\mu}L_{\mu}, m) . \qquad (102)$$

The smearing of the propagator has been dropped in the second line; this is a good approximation if $L >> 1/\Lambda$, as is true in practice. The precise form of the cutoff is irrelevant. Using Eq. (98), the corresponding result for I_2 is

$$f^{2}\overline{I}_{2}(m,L) = \sum_{n_{\mu}\neq 0} -m^{2} \frac{d}{dm^{2}} G(n_{\mu}L_{\mu},m) .$$
(103)

In the notation of Gasser and Leutwyler these results are [6]

$$f^{2}\overline{I}_{1}(m,L) = g_{1}(m^{2},L_{\mu}) ,$$

$$f^{2}\overline{I}_{2}(m,L) = m^{2}g_{0}(m^{2},L_{\mu}) .$$
(104)

The finite volume dependence of f_{π} , f_K , B_K , B_V , and B_A can now be predicted by substituting the results for \overline{I}_1 and \overline{I}_2 into the formulas presented in previous sections.

To actually calculate the \overline{I}_i , it is convenient to use the heat-kernel representation

$$G(x,m) = \frac{1}{(4\pi)^2} \int_0^\infty d\alpha \, \alpha^{-2} e^{-x^2/4\alpha} e^{-\alpha m^2} \,.$$
(105)

Using this, the functions g_0 and g_1 turn out to be particular cases of the general function

$$g_{r}(m,L_{\mu}) = \frac{1}{(4\pi)^{2}} \int_{0}^{\infty} d\alpha \, \alpha^{r-3} \sum_{n_{\mu} \neq 0} e^{-(n_{\mu}L_{\mu})^{2}/4\alpha} e^{-\alpha m^{2}} ,$$
(106)

the properties of which have been discussed extensively in Ref. [28]. The crucial point for present purposes is that the sum over n_{μ} converges rapidly, as long as $mL_{\mu} \gg 1$. This is because each term falls like $\exp(-mn_{\mu}L_{\mu})$, as can be seen by evaluating the α integrals in the saddle-point approximation

$$G(x,m) = \frac{m^2}{(4\pi)^2} \left[\frac{8\pi}{(mx)^3} \right]^{1/2} \times e^{-mx} \left[1 + \frac{3}{8mx} + O\left[\frac{1}{(mx)^2} \right] \right], \quad (107)$$

$$-m^{2} \frac{d}{dm^{2}} G(x,m) = \frac{m^{2}}{(4\pi)^{2}} \left[\frac{2\pi}{mx} \right] \times e^{-mx} \left[1 - \frac{1}{8mx} + O\left[\frac{1}{(mx)^{2}} \right] \right].$$
(108)

To illustrate the properties of the integrals, I set $L_1 = L_2 = L_3 = L$ and assume $L_4 >> L$, which is the geometry used in simulations. This means that in the sum over n_{μ} only the n_i are nonzero. It is useful to normalize the results by dividing by the saddle-point approximations to the terms involving propagation only to the six adjacent periodic images, i.e., $n_i = (1,0,0), (-1,0,0), (0,1,0)$, etc. These are

$$S_{1}(m,L) = 6 \frac{m^{2}}{(4\pi f)^{2}} \left(\frac{8\pi}{(mL)^{3}} \right)^{1/2} e^{-mL}, \qquad (109)$$

$$S_{2}(m,L) = 6 \frac{m^{2}}{(4\pi f)^{2}} \left[\frac{2\pi}{mL} \right]^{1/2} e^{-mL} .$$
 (110)

Figure 10 shows the ratios \overline{I}/S , computed numerically, for different numbers of terms in the sum over n_i . The notation is that $\overline{I}_i(N)$ includes the first N distances, where "distance" means $d = \sqrt{n_i n_i}$. Thus for N = 1 only the six d = 1 terms are kept, while for N = 4, the twelve $d = \sqrt{2}$, eight $d = \sqrt{3}$, and the six d = 2 terms are also included. Various points are noteworthy.

(1) It is straightforward to show that the ratios depend only on the product mL.

(2) The results for N = 1 show that the saddle-point approximation is accurate to better than 10% for mL > 3. If one includes the 1/mL corrections from Eqs. (107) and (108) then the agreement is better than 1%.

(3) The full result $(N = \infty)$ differs substantially from the leading-order term (N = 1) at small *mL*. Typical simulations work down to $m_{\pi}L = 3-4$, and for these it is clear that the leading term is not sufficient. This was not appreciated in Refs. [13,12,15].

(4) The bulk of the difference is taken up by the N=2 terms, with N=4 being a good approximation to the full result.

(5) The ratios for I_1 and I_2 at $N = \infty$ are very similar.

The need to include not only the d = 1 term in the sum over n_i raises the question of whether there are other corrections of comparable size. Higher loop diagrams give rise to terms which are also proportional to $\exp(-m_{\pi}L)$, but are suppressed by an additional power of $m_{\pi}^2/16\pi^2 f^2$. These may be substantial for present lat-



FIG. 10. The ratios of the finite volume part of the integrals \overline{I}_i to the saddle point approximants S_i for (a) I_1 and (b) I_2 .

tice pion masses. There are also corrections from two loop diagrams suppressed both by a power of m_{π}^2 , and by an additional power of $\exp(-m_{\pi}L)$. These are presumably small.

I close this section with a comment on the effect of the nonzero lattice spacing a. This cuts off the momenta at $p \approx 2\pi/a$, which should be much larger than the chiral cutoff Λ . This condition is satisfied in present simulations, for which $2\pi/a \ge 12$ GeV. Such a large cutoff should have little effect on the propagator over distances of L or greater. To check this, I replace the continuum propagator by that for a lattice scalar field for which the second derivative is discretized using the standard nearest-neighbor form. This is for illustrative purposes only: it is not clear how the pion field should be discretized, or indeed whether it should be discretized at all. My choice is really the worst case one can imagine, since discretizations involving more neighbors will give a better approximation to the continuum derivative. With the nearest-neighbor form, the lattice propagator is

$$G(x,y,z,t) = \int_0^\infty d\alpha \, I_x(2\alpha) I_y(2\alpha) I_z(2\alpha)$$
$$\times I_z(2\alpha) e^{-8\alpha} e^{-\alpha m^2}. \tag{111}$$

Here x, y, z, and t are integers giving the distances in lattice units, and I_x is the xth modified Bessel function (and should not be confused with the integrals I_1 and I_2 under study). The ratio of the integrals using this lattice propagator to those using the continuum propagator are shown in Fig. 11. The results are for N=1, but there is no



FIG. 11. Ratios of the $I_i(N=1)$ evaluated using the lattice propagator to those evaluated using the continuum propagator.

significant dependence on N. The ratio is close to unity, the difference not exceeding 10% for the values of m_{π} and L used in the simulations discussed in the following section. The effect is probably smaller than that of higher-order chiral corrections.

VII. NUMERICAL PREDICTIONS AND RESULTS

In this section, I compare the predictions for chiral logarithms and the associated volume dependence with numerical results. For the most part the predicted volume dependence is smaller than the statistical errors of present simulations. My aim in such cases is to show how close we are to testing the predictions. Only for B_V and B_A is there clear numerical evidence of finite volume dependence.

I use the numerical results from quenched simulations at $\beta=6$ presented in Refs. [13] and [24]. These calculations use staggered fermions, and work on $16^3 \times 40$ and $24^3 \times 40$ lattices. Finite size effects can be studied by comparing results from the two lattices. Using the lattice spacing obtained from f_{π} , 1/a=1.7 GeV, the smaller lattice has L=1.9 fm, the larger L=2.8 fm. Both lattices are probably large enough so that the squeezing of the pion wave function is a small effect. The lattice pions have masses ranging from $m_{\pi}a=0.24$ to 0.41, or 0.41-0.70 GeV in physical units. The convergence of chiral perturbation theory is controlled by the ratio $m_{\pi}^2/16\pi^2 f_{\pi}^2$, and this varies from 0.12 to 0.36. Thus higher-order corrections are likely to be sizable for the heavier pions.

The pions in these simulations are composed of quarks of masses $m_q a = 0.01$, 0.02, and 0.03. These are combined in all possible ways, so there are results for pions with both degenerate and nondegenerate quarks. The predictions for quenched chiral logarithms given above are only for degenerate quarks. It turns out, however, that the results for m_{π}^2 , f_{π} , B_K , B_V and other quantities depend mainly on the average quark mass. When I present results I will simply treat all pions as though they contained two degenerate quarks each with the average mass.

Typical chiral logarithmic corrections are numbers of O(1) multiplied by the integrals I_1 and I_2 . The finite

volume-dependent parts of these integrals are shown in Fig. 12. These are exact results, and are thus a sum of exponentials, multiplied by powers of m_{π} . Nevertheless, as the figure shows, the results are well approximated by single exponentials for the range of quark masses considered. The magnitude of the \bar{I}_i should be compared to the typical statistical errors in present numerical simulations, which are rarely as small as 1%. For the larger lattice, and for $m_{\pi}a = 0.24 - 0.41$, \bar{I}_1 and \bar{I}_2 are always smaller than 0.4%. Thus, for practical purposes, L = 24 corresponds to infinite volume. For the smaller lattice, however, \bar{I}_2 reaches 3% at $m_{\pi}a = 0.24$, which is a level that might be measurable.

To give an idea of the typical size of finite volume corrections I first consider f_{π} in a theory with four degenerate dynamical quarks. This is the theory represented by a single species of staggered fermions in the continuum limit. From Eq. (24) we see that

$$f_{\pi}(L) \approx f_{\pi}(L = \infty) [1 - 2\bar{I}_{1}(m_{\pi}, L)]$$
 (112)

It is clear from Fig. 12 that to observe this effect will require calculations with a precision of less than 1%. In the quenched approximation, the prediction is that there is no finite size dependence in f_{π} at leading order. Such dependence enters only due to the diagrams of Figs. 3(b) and 3(c), and is suppressed by $m_{\pi}^2/16\pi^2 f_{\pi}^2$. In fact, the numerical results of Ref. [24] show no evidence for finite size dependence. The errors are large, 5-10%, so that this does not constitute a test of the predicted difference between full and quenched chiral logarithms.

I note in passing that it will be even more difficult to see the volume dependence in m_{π} calculated with dynamical quarks. The result for N degenerate flavors is [6]

$$m_{\pi}^{2}(L) \approx m_{\pi}^{2}(L = \infty) \left[1 + \frac{1}{N} \overline{I}_{1}(m_{\pi}, L) \right].$$
 (113)

For the condensate, on the other hand, the situation is slightly better than for f_{π} [6]

$$\langle \bar{\psi}\psi \rangle(L) \approx \langle \bar{\psi}\psi \rangle(L = \infty) \left[1 - \frac{N^2 - 1}{N} \bar{I}_1(m_{\pi}, L) \right].$$
(114)



FIG. 12. \overline{I}_1 and \overline{I}_2 , for $N = \infty$, using 1/a = 1.7 GeV.

Since the condensate can be calculated more accurately than f_{π} and m_{π} , this equation may offer the best chance of testing the predictions of chiral loops in theories with dynamical quarks. The only problem is that there is a quadratically divergent perturbative contribution to the condensate, to which the prediction does not apply, and which must be subtracted.

As discussed in Sec. V, quenched chiral logarithms in m_{π}^2 come only from η' loops. Using Eqs. (89) and (86), the volume dependence is predicted to be

$$m_{\pi}^{2}(L) = m_{\pi}^{2}(L = \infty) \exp\left[\frac{A}{N} \frac{m_{0}^{2}}{m_{\pi}^{2}} \overline{I}_{2}(m_{\pi}, L)\right],$$
 (115)

where I have dropped terms suppressed by powers of m_{π}^2 . For N = 3, the phenomenological estimate from the physical η' mass gives $Am_0^2/3 \approx (0.5 \text{ GeV})^2$. Thus the factor multiplying I_2 ranges from 1.6 to 0.55 for $m_{\pi} = 0.41 - 0.70$ GeV. This means that, although the finite volume correction is parametrically enhanced by $1/m_{\pi}^2$, and is also increased by the exponentiation, the actual numerical enhancement is small. The volume dependence is somewhat larger than that for the full theory, because it is the integral I_2 which appears, rather than I_1 . The result is that, at $m_{\pi}a = 0.24$, the correction increases m_{π} (not its square) by 2.4% for L = 16, and by 0.2% for L = 24. The correction decreases rapidly with increasing pion mass. The numerical results for m_{π} show no significant finite size dependence, but the errors are about 2%, i.e., at same level of the expected effect [24]. Future quenched calculations may be able to test this prediction.

The form of the volume dependence of the condensate in the quenched approximation is the same as that for m_{π}^2 , because $m_{\pi}^2 f_{\pi}^2 \propto m \langle \bar{\psi}\psi \rangle$. Again the numerical results show no significant finite volume dependence, but the errors are too large (5-10% for the value extrapolated to $m_q=0$) for this to represent a disagreement with the prediction.

I now turn to the kaon B parameters. For degenerate quarks, the predicted volume dependence has the same functional form in full and quenched theories. From Eq. (45) I find

$$\delta B_K(L) = B_K(L) - B_K(L = \infty)$$

$$\approx B_K(L = \infty) [\overline{I}_2(m_K, L) - 2\overline{I}_1(m_K, L)] . \quad (116)$$

I use m_K rather than m_{π} to be consistent with the name of the *B* parameter; this does not imply that the quarks in the "kaon" have different masses. In Fig. 13 I show the fractional size of this correction. There is a large cancellation between \overline{I}_2 and \overline{I}_1 in the mass range of interest, and the correction does not exceed 0.2%. This is unfortunate as B_K can be calculated with errors as small as 1%. The prediction is thus that there be no finite volume dependence, and the results of Ref. [13] are consistent with this.

As discussed in Sec. IV, the chiral logarithms in B_V and B_A are enhanced by $1/m_K^2$ over those in B_K , and one might hope that the associated finite size effects can be seen. From Eqs. (79) and (80), I find

46



FIG. 13. Predictions for the fractional correction $\delta B_K(L)/B_K(L = \infty)$.

$$B_{V} = \frac{1}{2} B_{K} - \frac{3}{4} (\beta_{1} + \beta_{2}) \frac{\mu^{2}}{m_{K}^{2}} I_{2}(m_{K}) - \frac{3}{8} (\delta_{1} + \delta_{2}) - \frac{3}{8} (\gamma_{1} + \gamma_{2}) I_{2}(m_{K}) + O(m_{K}^{2}) , \qquad (117)$$

where B_K is one-loop corrected result, Eq. (45). The enhanced chiral logarithm is that proportional to I_2/m_K^2 . The important features of this prediction are the following.

(a) Since $I_2/m_K^2 \propto \ln(\Lambda/m_K)$, B_V should diverge logarithmically in the chiral limit. This is unlike B_K , for which the chiral logarithms, being proportional to $m_K^2 \ln(m_K)$, vanish in the chiral limit.

(b) The coefficient of this divergence is not related to B_K ; it is a new constant proportional to $\beta_1 + \beta_2$. In vacuum saturation approximation $\beta_1 + \beta_2 = \frac{4}{3}$. Typically vacuum saturation gives the correct sign and order of magnitude; if so, B_V is predicted to diverge negatively.

(c) The extent of the enhancement is much greater than that in the correction to the quenched m_{π}^2 , Eq. (115). The constant $\mu = m_K^2 / (m_s + m_d)$ is about 5 GeV on the lattice, roughly ten times the corresponding quantity $(Am_0^2/N)^{1/2}$.

(d) Since I_2 increases with L, B_V is predicted to become more negative as L decreases.

Since $B_A = B_K - B_V$, the enhanced logarithm in B_A differs only in sign from that in B_V . Thus B_A is predicted to have a positive logarithmic divergence, and to increase as *L* decreases. Since the numerical results for B_K behave as expected, i.e., they have a smooth chiral limit and show no volume dependence, and since the errors in B_K are much smaller than those in B_V and B_A , there is no additional information to be gained from studying B_A . Thus I concentrate on B_V in the following.

The numerical results of Ref. [13] for B_V are shown in Fig. 14. There are two striking features: the results appear to be diverging as the chiral limit is approached, with the sign expected from the chiral logarithm; there is a significant finite volume dependence, which also has the expected sign. It was, in fact, these features that sparked the present study.

To provide a quantitative test of Eq. (117), I fit the data using



FIG. 14. Results for B_V vs m_{π}^2 in lattice units. The fits are described in the text.

$$B_{V}(L) = -\frac{\mu^{2}}{16\pi^{2}f^{2}} \left[c_{1} \ln \left[1 + \frac{1}{x} \right] + c_{2} \frac{1}{1+x} + c_{3} + c_{4} \frac{x^{2}}{1+x^{2}} \right] - c_{1} \frac{\mu^{2} \overline{I}_{2}(m_{K}, L)}{m_{K}^{2}} , \qquad (118)$$

where $x = m_K^2 / \Lambda^2$, Λ being the cutoff on chiral loop integrals. This function contains the chiral logarithm and its associated finite volume dependence, together with a power series in m_K^2 . The power series is supposed to represent all terms in Eq. (117) aside from the enhanced chiral logarithm. This includes the normal chiral logarithms proportional to $m_K^2 \ln(m_K)$, which can be well approximated by a polynomial for the range of masses under consideration. I choose the particular form of the first two terms in Eq. (118) because the loop integral is

$$16\pi^2 \int_{|k| < \Lambda} G(k,m)^2 = \ln(1+1/x) - 1/(1+x) .$$
 (119)

The denominator of the c_4 term is chosen so that the function is smooth for $x \sim 1$. Finally, the flavor symmetry breaking for staggered fermions is accounted for by applying the prescription (83), with $R_{\rm NG} = 1$, to all terms in Eq. (118).

Figure 14 shows the best fit. I have taken Λ to be 0.8 GeV; the precise value is unimportant since, to good accuracy, changes can be absorbed into the other constants. These constants are

$$c_1 = 0.29, c_2 = -1.25,$$

 $c_3 = 0.63, c_4 = -0.37.$
(120)

The fit is reasonable, with a χ^2 of 1.4 per degree of freedom. I therefore conclude that the numerical results are consistent with the expected chiral logarithm. The coefficient c_1 is smaller than the result expected from vacuum saturation approximation, $c_1=1$, but I do not consider this to be a serious problem for the following reasons.

(i) Vacuum saturation can be a poor approximation.



FIG. 15. Finite size dependence of B_V . The line shows the result of the same fit as in Fig. 14.

For example, the matrix \mathcal{M}_K may be as much as 50% smaller than the prediction of vacuum saturation [29]. What matters here, however, are the matrix elements of a pseudoscalar-pseudoscalar operator [cf. Eq. (71)], and these are likely to be better approximated by vacuum saturation [30].

(ii) I have assumed (by setting $R_{NG}=1$) that the vertices of the non-Goldstone and Goldstone pions are the same, whereas in fact the former could be smaller. The matrix elements of some non-Goldstone bilinears are smaller by as much as 25% [20]. Since the loop diagrams involve the matrix element of operators consisting of two bilinears, and also a four pion scattering amplitude, a more substantial reduction is possible for c_1 .

(iii) There could be substantial higher-order chiral perturbation theory corrections to the coefficient of the enhanced logarithm. That higher-order terms can be substantial is illustrated by the fact that $f_K^2/f^2=1.6$ for the heaviest lattice kaon.

The fit does not, however, demonstrate the existence of the chiral logarithm. In particular, the coefficient c_1 is determined not by the nature of the divergence, but by the finite size dependence. If one allows the coefficient of the logarithm to differ from that of the \overline{I}_2 term, the fit favors much smaller values for the former. In fact, a fit as good as that displayed above is obtained with no logarithmic term at all. This flexibility is possible because x is not small, and varies over a small range, 0.5-0.8, so the nonlogarithmic terms in the fit function can mock up a logarithm.

It is thus the finite volume dependence which is the strongest evidence for quenched chiral logarithms. The functional dependence of $B_V(16) - B_V(24)$ on m_K is predicted, leaving only the overall factor to be determined by the fit. As shown in Fig. 15, the fit does a reasonable job of representing the dependence on m_K . The shape can be altered by higher-order chiral terms, which are proportional to m_K^2 , so one would not expect a perfect match. To provide a more rigorous test of the prediction results from smaller kaon masses are needed.

VIII. CONCLUSIONS

Chiral logarithms can provide powerful consistency checks on numerical simulations. The theoretical calculations are well founded if dynamical quarks are included. I have argued that the methods of chiral perturbation theory, suitably modified, can also be applied in the quenched approximation. One can imagine a program in which one first tests the results of quenched chiral perturbation theory using simple quantities such as f_{π} , m_{π} , and $\langle \bar{\psi}\psi \rangle$, and then applies the predictions to other quantities in order to test that the physics of pion loops is being properly included. Furthermore, as I argued in the Introduction, if the calculations are reliable one can use them to estimate the size of the errors introduced by the quenched approximation.

Such a program requires numerical results with errors of 1% or less, calculated on a number of lattice sizes at a variety of pion masses. This should be possible in the fairly near future, as computer power approaches the TeraFlop milestone. The numerical results presented here are encouraging, but not conclusive. What is particularly needed is for the calculations to be pushed to smaller pion masses. On a given lattice there is a minimal mass, $m_{\pi}(\min)$, below which chiral symmetry is restored [6]. This minimal mass scales with size as $m_{\pi}(\min)^2 \propto L^{-3}$. This means that $m_{\pi}(\min)L \propto 1/\sqrt{L}$, so that, by working on larger lattices, one can push to smaller $m_{\pi}L$ and thereby increase the finite size effects.

A program of testing finite volume dependence also requires that quenched chiral logarithms be calculated for other quantities. Examples for which there is hope of good numerical results in the next few years include the matrix elements of the $\Delta S = 1$ operators responsible for kaon decays, both CP conserving and violating, and the properties of heavy-light systems, such as B and D mesons. Calculations of the finite size effects in f_B and f_D in full QCD have already been done [31]. For some quantities the quenched calculations appear to be a straightforward extension of those presented here. For others, one must face the problem of how best to include η' loops. It may be that the perturbative approach of Bernard and Golterman [11] will be adequate. Or it may be that a resummation, such as that presented here for the pion mass and the condensate, is required.

It does not appear to be simple to apply the methods presented here to all quantities. In particular, in the calculation of chiral corrections to baryon properties, it may not be possible to uniquely assign quark diagrams to chiral perturbation theory diagrams. This problem is under study.¹³

ACKNOWLEDGMENTS

I am very grateful to Claude Bernard, Heiri Leutwyler, and Apoorva Patel for many discussions and helpful comments, Claude Bernard and Maarten Golterman for comments on the manuscript, and to Rajan Gupta, Greg Kilcup, and Apoorva Patel for their collaboration in obtaining the numerical results. I also thank Norman

¹³Cohen and Leinweber [32] have recently begun the study of quenched chiral logarithms in baryon properties.

Christ and Paul Mackenzie for helpful comments. Parts of this work were done at the Santa Fe Workshop on QCD, at the Aspen Center for Physics, and at the Institute for Theoretical Physics, Santa Barbara, all of which I thank for their excellent hospitality. Finally I owe partic-

- [1] S. Weinberg, Phys. Rev. Lett. 17, 616 (1966).
- [2] J. Gasser and H. Leutwyler, Nucl. Phys. B250, 465 (1985).
- [3] L.-F. Li and H. Pagels, Phys. Rev. Lett. 26, 1204 (1971).
- [4] P. Langacker and H. Pagels, Phys. Rev. D 8, 4595 (1973).
- [5] J. Bijnens, H. Sonoda, and M. Wise, Phys. Rev. Lett. 53, 2367 (1985).
- [6] J. Gasser and H. Leutwyler, Phys. Lett. B 184, 83 (1987); 188, 477 (1987).
- [7] M. Lüscher, Commun. Math. Phys. 104, 177 (1986).
- [8] A. Morel, J. Phys. (Paris) 48, 1111 (1987).
- [9] A. Billoire, R. Lacaze, E. Marinari, and A. Morel, Nucl. Phys. B271, 461 (1986).
- [10] S. Sharpe, Phys. Rev. D 41, 3233 (1990).
- [11] C. Bernard and M. Golterman, in Lattice '91, Proceedings of the International Symposium, Tsukuba, Japan, 1991, edited by M. Fukugita, Y. Iwasaki, M. Okawa, and A. Ukawa [Nucl. Phys. B (Proc. Suppl.) 26, 360 (1992)]; Phys. Rev. D 46, 853 (1992).
- [12] S. Sharpe, in *Lattice '89*, Proceedings of the International Symposium, Capri, Italy, 1989, edited by R. Petronzio et al. [Nucl. Phys. B (Proc. Suppl.) 17, 146 (1990)].
- [13] G. Kilcup, S. Sharpe, R. Gupta, and A. Patel, Phys. Rev. Lett. 64, 25 (1990).
- [14] S. Sharpe, in *Lattice '90*, Proceedings of the International Symposium, Tallahassee, Florida, 1990, edited by U. M. Heller, A. D. Kennedy, and S. Sanielevici [Nucl. Phys. B (Proc. Suppl.) 20, 429 (1991)].
- [15] S. Sharpe, in Standard Model, Hadron Phenomenology and Weak Decays on the Lattice, edited by G. Martinelli (World Scientific, Singapore, in press).
- [16] H. Georgi, Weak Interactions and Modern Particle Theory, (Benjamin/Cummings, Menlo Park, CA, 1984).
- [17] S. Weinberg, Physica A96, 327 (1975).

ular thanks to Nathan Isgur, for giving me the time and facilities necessary to finish this project. This work was supported in part by the DOE under Contract DE-AC05-84ER40150 and Grant DE-FG09-91ER40614, and by the Alfred P. Sloan Foundation.

- [18] M. Bochiccio, L. Maiani, G. Martinelli, G. Rossi, and M. Testa, Nucl. Phys. **B262**, 331 (1985); L. Maiani, G. Martinelli, G. Rossi, and M. Testa, *ibid*. **B289**, 505 (1987).
- [19] G. Kilcup and S. Sharpe, Nucl. Phys. B283, 493 (1987).
- [20] S. Sharpe, R. Gupta, and G. Kilcup, University of Washington Report No. DOE/ER/40423-23, 1992 (unpublished).
- [21] A. Redfield, Phys. Lett. 109B, 311 (1982); J. Honerkamp and K. Meetz, Phys. Rev. D 3, 1996 (1971); J. M. Charap, *ibid.* 3, 1998 (1971).
- [22] W. Bardeen, A. Buras, and J.-M. Gérard, Phys. Lett. B 192, 138 (1987).
- [23] W. Bardeen, A. Buras, and J.-M. Gérard, Phys. Lett. B 221, 343 (1988).
- [24] See, for example, R. Gupta, G. Guralnik, G. Kilcup, and S. Sharpe, Phys. Rev. D 43, 2003 (1991).
- [25] J. Gasser and H. Leutwyler, Nucl. Phys. B307, 763 (1988).
- [26] S. Sharpe, in *Particles, Strings and Cosmology*, Proceedings of the First International Symposium, Boston, Massachusetts, 1990, edited by P. Nath and S. Reucroft (World Scientific, Singapore, 1991), pp. 270-291.
- [27] M.-C. Chu, M. Lissia, and J. Negele, Nucl. Phys. B360, 31 (1991).
- [28] P. Hasenfratz and H. Leutwyler, Nucl. Phys. **B343**, 241 (1990).
- [29] S. Sharpe, in [11], p. 197.
- [30] C. Bernard and A. Soni, in *Lattice '88*, Proceedings of the International Symposium, Batavia, Illinois, 1988, edited by A. S. Kronfeld and P. B. Mackenzie [Nucl. Phys. B (Proc. Suppl.) 9, 155 (1989)].
- [31] J. Goity, Phys. Lett. B 249, 495 (1990).
- [32] T. D. Cohen and D. B. Leinweber, University of Maryland Report No. 92-189, 1992 (unpublished).