

Studies with Improved Renormalization Group Techniques

by

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Dedication

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Contents

Chapter

1	Introduction	1
1.1	Motivation	1
1.2	Organization	3
1.3	Notation and Conventions	5
2	Strongly Coupled Physics Beyond the Standard Model	7
2.1	The Standard Model	7
2.2	Higgs Mechanism	10
2.3	Beyond Standard Model Physics	14
2.4	Composite Higgs	17
2.4.1	Technicolor	17
2.4.2	Extended Technicolor	20
2.4.3	Conformal Window	21
2.4.4	Lattice Studies	24
3	Lattice Field Theory	29
3.1	The Lattice	29
3.2	Gauge Fields	29
3.2.1	Pure Gauge	29
3.2.2	Negative Adjoint Term	32

3.3	Fermionic Fields	32
3.3.1	Naive Fermions	32
3.3.2	Wilson Fermions	33
3.3.3	Staggered Fermions	34
3.3.4	Improved Fermions	35
4	Monte Carlo Renormalization Group	38
4.1	Introduction	38
4.2	Method	39
4.2.1	Blocking	40
4.2.2	Two-lattice matching procedures and the need for optimization	43
4.2.3	Chirally Broken Theories	46
4.2.4	Conformal Theories	46
4.3	8 and 12 Flavor Results	47
4.3.1	8 Flavors	49
4.3.2	12 Flavors	49
4.3.3	The Case of the Wandering Fixed Point	50
5	Wilson Flow	57
5.1	Introduction	57
5.2	Improving the Gradient Flow Step Scaling	59
5.3	4 Flavor Test	61
5.4	12 Flavor Results	65
5.5	Summary	67
6	Wilson Flow MCRG	71
6.1	Introduction	71
6.2	Wilson Flow Optimized MCRG	72

6.3	12 Flavor Results	73
7	Conclusion	79
	Bibliography	82
	Appendix	

Tables

Table

2.1	The standard model fermions are spin 1/2 fundamental particles in the color singlet. The masses for the electron, muon, and tau are known to the precision given. The neutrinos are known to have masses but their masses are not precisely measured. Left handed leptons have weak hypercharge of -1 while right handed leptons have weak hypercharge of -2. Right handed neutrinos have not been observed.	11
2.2	The standard model quarks are spin 1/2 fundamental particles that are charged under $SU(3)_C$. Left handed quarks have 1/3 hypercharge. Right handed quarks with 2/3 electric charge have weak hypercharge of 4/3 while the right handed quarks with -1/3 electric charge have weak hypercharge of -2/3.	11
2.3	The standard model bosons. All particles are spin 1 except for the Higgs particle which is spin 0.	11
2.4	This table summarizes some commonly used groups, their dimension, first Casimir invariant, and second Casimir invariant. F, S_2, A_2 and G are the is the fundamental, 2-index symmetric, 2-index antisymmetric, and adjoint representations respectively. .	22

Figures

Figure

2.1 These are the Feynman Diagrams of the allowed QCD verticies. All interactions in the theory are built by combining these verticies. Quarks are represented by the solid lines and gluons are represented by curly lines. The three and four gluon vertecies are unique to non-Abelian gauge theories. 8

2.2 There is only one allowed vertex in QED which couples a fermion charged under U(1) with the photon. 10

2.3 The Higgs potential with $\mu^2 > 0$ has a spontaneously broken global symmetry. We can think of this classically. The origin is a local maximum of potential energy and therefore an unstable equilibrium. If we place a particle at the origin any perturbation will push the particle ‘down hill’ to a minimum of the potential. Since there are a number of degenerate minima, choosing one spontaneously breaks the symmetry. 13

2.4 In a confining theory, such as QCD, the running of the coupling starts out negative. As g^2 grows the running becomes increasingly negative. Chiral symmetry is broken and the theory is described by confined bound states. 27

2.5 In a conformal field theory the β function starts out negative but turns around and crosses the $\beta = 0$ axis. At this crossing the theory has an infrared fixed point. The renormalized flow for the theory will run to the IR fixed point and the theory is conformal. 27

- 2.6 If the fermionic degrees of freedom overwhelm those of the bosons it the beta function will start positive. This is the loss of asymptotic freedom. The beta function for such a theory resembles that of QED. 28
- 2.7 In a walking theory the beta function looks similar to the conformal picture of figure 2.5. The crucial difference is that before the β function develops a zero, shown by the dashed line, chiral symmetry is spontaneously broken driving the beta function back down. Near the would be IRFP where the dashed line has a zero, the coupling runs very slowly and is said to be walking. 28
- 3.1 The phase diagram of SU(3) gauge theory in the fundamental-adjoint coupling plane. The solid lines are first order phase transitions. The phase transition ends in a second order critical point. Although the critical point lies above the β_F axis, the presence of the critical point influences the RG flow of actions where $\beta_A = 0$. Therefore we choose an action with $\beta_A < 0$ to avoid effects from the critical point. The phase diagram for confining SU(3) gauge theories with fermions is similar, only the location of the transitions changes. 37
- 4.1 The original 6x6 lattice on the left possesses a discrete scaling symmetry of $s = 2$ and $s = 3$. The shaded orange square is a $s = 2$ block variable. The resulting orange 3×3 blocked lattice in the bottom right formed by replacing each block variable with a single site in the upper left corner of the block. The cyan shaded region shows a $b = 3$ block variable. Performing a block transformation that replaces each block with a point in the upper left of the block produces the 2×2 blocked lattice shown in the upper right. 51

- 4.2 This figure shows how links are blocked on the lattice. Two adjacent links in the same direction are block transformed to form one link of twice the lattice spacing. We perform all possible block transformation shown on the left as the red, blue, green, and orange block tilings of the unblocked lattice. We then store the links of the blocked lattice as shown on the right hand side of the figure. 51
- 4.3 Here I show coupling space for a system with one relevant direction K_0 . All irrelevant directions are collected in K' . We simulate at a point P in parameter. As we block the system it the effective couplings will change. In the diagram here the couplings reach the renormalized trajectory after 3 blocking steps. Further blocking steps move the couplnigs along the renormalized trajectory. 52
- 4.4 For matching we pick two points in coupling space P_1 and P_2 . After 3 blocking steps, shown as circles the effective action of P_2 has reached the renormalized trajectory. P_1 requires 4 blocking steps shown as stars but reahes the same point on the renormalized trajectory. Because P_2 took one less blockign step its correlation length ξ is a factor of s smaller than the correlation length of the first ensemble that started at P_1 . By choosing pairs of points in coupling space that block to the same point on the renormalized trajectory in coupling space, we can construct a discrete step scaling function. 52
- 4.5 The RG flow of a confining theory on the $m = 0$ critical surface. β is the relevant gauge coupling, β' are irrelevant couplings. 53
- 4.6 The RG flow of a conformal theory on the $m = 0$ critical surface. β is the relevant gauge coupling, β' are irrelevant couplings. 53
- 4.7 Optimization of the HYP-smearing parameter α in the RG blocking transformation, for $\beta_F = 5.0$. The uncertainties on the data points are dominated by averaging over the different observables as described in the text. 54

- 4.8 Results for the bare step-scaling function s_b from traditional MCRG two-lattice matching with $24^3 \times 48$, $12^3 \times 24$ and $6^3 \times 12$ lattice volumes for $N_f = 8$. The blue dashed lines are perturbative predictions for asymptotically weak coupling. 54
- 4.9 Results for the bare step-scaling function s_b from traditional MCRG two-lattice matching with $24^3 \times 48$, $12^3 \times 24$ and $6^3 \times 12$ lattice volumes for $N_f = 12$. The blue dashed lines are perturbative predictions for asymptotically weak coupling. 55
- 4.10 An illustration of how optimizing the block transformation can result in difficulties locating an IRFP, β is the relevant gauge coupling and β' are irrelevant couplings. The upper figure shows the renormalized trajectory in red, green, and blue found by optimizing the RG transformation at β_1, β_2 , and β_3 respectively. The location of the IRFP changes in each renormalized trajectory, in this picture the IRFP is moved to the coupling we perform MCRG at. The resulting s_b is consistent with zero across a wide range of couplings. 56
- 5.1 Continuum extrapolations of the discrete β_{lat} function of the $N_f = 4$ system at $\tilde{g}_c^2(L) = 2.2$ with several different values of the t -shift coefficient τ_0 . The dotted lines are independent linear fits at each τ_0 , which predict a consistent continuum value. . 62
- 5.2 Continuum extrapolations of the discrete β_{lat} function of the $N_f = 4$ system for several different $\tilde{g}_c^2(L)$ values. For $\tilde{g}_c^2(L) = 1.8, 2.2$ and 2.6 $\tau_0 = -0.02$ is near-optimal, while the larger couplings $\tilde{g}_c^2(L) = 3.0$ and 3.4 require $\tau_0 = -0.01$ to remove most $\mathcal{O}(a^2)$ effects. The colored points at $(a/L)^2 = 0$ are the continuum extrapolated results, while the black crosses at $(a/L)^2 = 0$ show the corresponding two-loop perturbative predictions. 63
- 5.3 The $N_f = 12$ running coupling $g_c^2(L)$ versus the bare coupling β_F on several volumes, for $c = 0.2$. Crossings between results from different volumes predict the finite volume IRFP coupling $g_\star^2(L)$ in this scheme. 66

- 5.4 Continuum extrapolations of the 12-flavor finite volume IRFP $g_\star^2(L)$, with several different t -shift coefficients τ_0 for fixed scale change $s = 2$. The dotted lines are a joint linear fit constrained to have the same $(a/L)^2 = 0$ intercept, which gives $g_\star^2 = 6.21(25)$ 68
- 5.5 Continuum extrapolations of the 12-flavor finite volume IRFP $g_\star^2(L)$, with several different scale changes for the near-optimal improvement coefficient $\tau_{\text{opt}} \approx 0.04$. The $s = 4/3$ and $3/2$ data points are horizontally displaced for greater clarity. The dashed lines are a joint linear fit constrained to have the same $(a/L)^2 = 0$ intercept, which gives $g_\star^2 = 6.18(20)$ 69
- 6.1 In Wilson Flow MCRG we use the Wilson Flow (blue) to approach the renormalized trajectory. An optimization step similar to that used in MCRG allows us to locate the flow time that gets us closest to the renormalized trajectory (orange). We then block our lattice using a fixed block transformation (green). 76
- 6.2 Optimization of the Wilson flow time t_f with fixed $\alpha = 0.5$, for $\beta_F = 4.5$. The uncertainties on the data points are dominated by averaging over the different observables. 76
- 6.3 The bare step-scaling function s_b predicted by three-lattice matching with 6^4 , 12^4 and 24^4 lattices blocked down to 3^4 , comparing three different renormalization schemes. The error bars come from the standard deviation of predictions using the different observables. 77
- 6.4 As in Fig. 6.3, the bare step-scaling function s_b for three different renormalization schemes from three-lattice matching, now using 8^4 , 16^4 and 32^4 lattices blocked down to 4^4 77
- 6.5 The bare step-scaling function s_b for scheme 1, comparing three-lattice matching using different volumes: 6^4 , 12^4 and 24^4 lattices blocked down to 3^4 (black \times s) as well as 8^4 , 16^4 and 32^4 lattices blocked down to 4^4 (blue bursts) and 2^4 (red crosses). 78

Chapter 1

Introduction

1.1 Motivation

The standard model (SM) of particle physics is among the most rigorously tested theories describing the physical world we live in. It describes three of the four fundamental forces of nature in terms of quantum interactions of fundamental particles. Recently, on July 4, 2012, the discovery of the Higgs particle was announced. This is an exciting discovery for many reasons. First of all, the Higgs is predicted by the standard model as are many of its properties. In this sense its discovery is a good test of the standard model. Additionally, the Higgs was the last ‘missing’ piece of the standard model, now we have a complete theory. Finally, the discovery of the Higgs puts very hard constraints on theories of Beyond Standard Model (BSM) physics.

The Higgs particle in the standard model is an excitation of the Higgs field. The Higgs field is responsible for electroweak symmetry breaking (EWSB). EWSB explains how the W^\pm and Z bosons acquire a mass that would otherwise be forbidden. Explicitly putting a mass term in the Electroweak theory violates local gauge invariance, a critical component of the theory. By allowing the W^\pm and Z bosons to couple to the Higgs field, they are allowed to become massive without violating local gauge invariance. Additionally, in the EWSB process, the Higgs develops a vacuum expectation value (VEV). By coupling to the Higgs VEV, the standard model fermions acquire mass as well. In this way the Higgs provides an explanation of how all the fundamental particles gain their mass.

The Higgs field in the standard model is a scalar field. This leads to the hierarchy problem

(also equivalent to the naturalness or fine tuning problem) as the Higgs mass will be dependant on the UV cutoff of the standard model. If the standard model is the complete story and there is no new physics just above the electroweak scale then the SM cutoff will be at a very high scale. Navelly, we expect the Higgs mass to be at the same order as the cutoff. The fact the Higgs mass is 126 GeV requires fine tuning in the electroweak sector to cancel the UV divergences.

There have been many solutions proposed as an alternative to the Standard Model that don't require fine tuning. Supersymmetry (SUSY) is one such proposal that achieves cancellations by introducing a new symmetry between bosonic and fermionic fields. Since the bosons and fermions enter loop calculations with opposite signs there is no need for fine tuning. Most models of SUSY predict a spectrum of super partner masses near the Higgs mass. The fact that we have not yet observed any particles besides the Higgs puts constraints on SUSY models. Additionally most versions of SUSY that have not been ruled out replace the standard model hierarchy problem with a new 'little' hierarchy problem where fine tuning is still required but it is not as severe as it is in the standard model.

Another natural possibility is that EWSB is the result of new strong dynamics at the TeV scale. Under this prescription, EWSB happens much like spontaneous chiral symmetry breaking in quantum chromodynamics (QCD). QCD is the only strongly coupled sector of the standard model and is responsible for the strong nuclear force. Theories of EWSB that rely on strong dynamics are generally called Technicolor models. Technicolor is also highly constrained by the Higgs mass, the top quark mass, and precision electroweak measurements that show flavor changing neutral currents are highly suppressed.

It is natural to draw insights from the parallels of Technicolor and QCD or directly from perturbation theory, however such insights do not paint an accurate picture. The strongly coupled nature of these theories demands that they are studied through fully nonperturbative means. Therefore, lattice gauge theory is the only controlled way to study these theories. In lattice gauge theory we discretized euclidean space time into a regular grid of sites connected by links. The Lagrangian describing the theory is also discretized in such a way that in the continuum limit the

original theory is recovered. The continuum limit is the limit that spacing between lattice sites is infinitesimally small. Lattice QCD, the study of QCD on the lattice, has been extremely successful in understanding QCD. Many nonperturbative effects such as confinement, hadronization, chiral symmetry breaking, and instantons have been studied extensively. Many of these properties are completely inaccessible through analytic techniques while others could not be fully understood.

The ultimate goal of efforts in this field is to find a theory can explain electroweak symmetry breaking, as it is observed in nature, through new strong dynamics. Finding such a theory is highly nontrivial and a great deal of theoretical and computational effort has gone into looking for such a theory. The lattice community has set out to do this in a controlled and systematic way: map out the space of all possible theories and see what we find. To date a lot of the focus of the lattice community has been on locating the conformal window and developing techniques to distinguish conformal from chirally broken theories. The conformal window describes a class of theories that develop an attractive infrared fixed point. These theories are scale invariant and are not candidates for Technicolor.

This thesis has two primary objectives. The first goal is a physics goal: map out the behavior of $SU(3)$ gauge theories with N_f fermions in the fundamental representation. For technical reasons we choose to study theories with 4, 8, 12, and 16 flavors. 4 and 16 flavor theories are already well understood so the focus of our work is on 8 and 12 flavor systems.

My second objective is to develop tools that make the first goal possible. Towards this end I introduce improvements to lattice step scaling calculation techniques. I explore how these techniques are different and how to implement them in a calculation. Finally these techniques are applicable to any lattice calculation and therefore will be of use to any lattice practitioner.

1.2 Organization

This thesis is organized into five chapters not including the introduction and conclusion. Briefly, I use chapter 2 to set the stage and chapter 3 to give a flavor for lattice calculations. Chapters 4-6 discuss my contributions, each chapter is its own calculation and reflects approximately

a years worth of work.

I begin chapter 2 with a brief overview of the standard model. The Higgs mechanism is given its own section due to its relevance for the remainder of the thesis. I then summarized flaws in our current understanding of the standard model with emphasis given to the hierarchy problem. Finally I discuss how a composite Higgs can resolve the hierarchy problem in a natural way. Because a composite Higgs arises from a strongly interacting quantum field theory we cannot understand such a theory with the standard analytical tools of perturbation theory. Fortunately there is a well developed nonperturbative approach, lattice gauge theory. I conclude the chapter with an overview of current lattice studies.

Chapter 3 provides the reader with a brief overview of lattice gauge theory. Like any mature discipline, this topic is a subject of many books. My goal is not to cover this subject comprehensively; rather I want to provide the reader with a flavor of how a lattice calculation is performed and highlight areas that are useful in later sections. To that end I develop the pure gauge action and discuss the negative adjoint term that our group uses in our simulations. I also discuss the problems associated with the discretization of fermions on the lattice. I end the discussion of lattice fermions with a description of staggered fermions and the nHYP smearing we use in our calculations.

Chapter 4 - 6 document the body of work that is this thesis. Chapter 4 concerns my early work using Monte Carlo Renormalization Group (MCRG) on $SU(3)$ $N_f = 8$ and $SU(3)$ $N_f = 12$ theories. In chapter 5, I introduce the gradient flow and how it can be used to calculate the renormalized step scaling function. I also use this chapter to discuss an optimization improvement that our group developed. Results for $SU(3)$ $N_f = 4$ and $SU(3)$ $N_f = 12$ are given. Finally chapter 6 shows how the Wilson Flow can be used as an optimization step for MCRG. Results from $SU(3)$ $N_f = 12$ are given.

In my concluding chapter I make some final remarks about interpreting our results. I will also summarize results from other methods that our group used to study this system. All of the methods our group has used to study there systems have converged to the same answer making a strong case that $SU(3)$ gauge theories with 12 flavors of fundamental fermions has an infrared fixed

point.

1.3 Notation and Conventions

Unless otherwise noted these are the standard conventions used in this thesis. I use standard particle theory units of $\hbar = c = 1$. The four vector is given by Greek indices

$$x^\mu = \begin{pmatrix} x_0 \\ x_1 \\ x_2 \\ x_3 \end{pmatrix}, \quad (1.1)$$

where x_0 is the temporal component t and the remaining three components are the spatial components of \vec{x} . In the continuum, I use the west coast convention for the metric

$$\eta_{\mu\nu} = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & -1 & 0 & 0 \\ 0 & 0 & -1 & 0 \\ 0 & 0 & 0 & -1 \end{pmatrix}. \quad (1.2)$$

On the lattice we use the Euclidian metric

$$\eta_{\mu\nu} = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{pmatrix}. \quad (1.3)$$

Standard Einstein summation notation applies to contract indices.

I work in the chiral representation for the Dirac spin matrices γ^μ

$$\gamma^0 = \begin{pmatrix} 0 & I \\ I & 0 \end{pmatrix}, \quad \gamma^i = \begin{pmatrix} 0 & \sigma^i \\ -\sigma^i & 0 \end{pmatrix}, \quad (1.4)$$

where I is the 2×2 identity matrix and σ^i are the Pauli matrices:

$$I = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}, \quad \sigma^1 = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \quad \sigma^2 = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}, \quad \sigma^3 = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}. \quad (1.5)$$

I use the standard ‘slash’ notation to contract four vectors with γ^μ :

$$\not{p} \equiv p_\mu \gamma^\mu. \quad (1.6)$$

The momentum operator I use the notation

$$p^\mu = i\partial^\mu, \quad (1.7)$$

where

$$\partial^\mu = \left(\frac{\partial}{\partial x^0}, -\nabla \right) = \left(\frac{\partial}{\partial x^0}, \frac{\partial}{\partial x^1}, \frac{\partial}{\partial x^2}, \frac{\partial}{\partial x^3} \right). \quad (1.8)$$

The Gell-Mann matrices λ_a are:

$$\begin{aligned} \lambda_1 &= \begin{pmatrix} 0 & 1 & 0 \\ 1 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix} & \lambda_2 &= \begin{pmatrix} 0 & -i & 0 \\ i & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix} \\ \lambda_3 &= \begin{pmatrix} 1 & 0 & 0 \\ 0 & -1 & 0 \\ 0 & 0 & 0 \end{pmatrix} & \lambda_4 &= \begin{pmatrix} 0 & 0 & 1 \\ 0 & 0 & 0 \\ 1 & 0 & 0 \end{pmatrix} \\ \lambda_5 &= \begin{pmatrix} 0 & 0 & -i \\ 0 & 0 & 0 \\ i & 0 & 0 \end{pmatrix} & \lambda_6 &= \begin{pmatrix} 0 & 0 & 0 \\ 0 & 0 & 1 \\ 0 & 1 & 0 \end{pmatrix} \\ \lambda_7 &= \begin{pmatrix} 0 & 0 & 0 \\ 0 & 0 & -i \\ 0 & i & 0 \end{pmatrix} & \lambda_8 &= \frac{1}{\sqrt{3}} \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & -2 \end{pmatrix} \end{aligned} \quad (1.9)$$

Chapter 2

Strongly Coupled Physics Beyond the Standard Model

2.1 The Standard Model

In our current understanding of the universe, there are four fundamental forces in nature: gravity, electromagnetism, the strong nuclear force, and the weak nuclear force. The latter three forces are understood in terms of quantum field theories which form the Standard Model (SM). The Standard Model is a $SU(3)_C \times SU(2)_W \times U(1)_Y$ gauge theory. Y is electroweak hypercharge and is related to electric charge and the weak neutral current.

The $SU(3)_C$ component of the theory explains the strong nuclear force, also called quantum chromodynamics (QCD). QCD is a $SU(3)$ Yang Mills theory with fermions in the fundamental representation. This force explains how spin 1/2 quarks possessing ‘color’ charge interact by exchanging gluons. It is important to note that this color charge has nothing to do with visible colors which are actually a narrow band of electromagnetic radiation that we can see. Rather the color charge is analogous to electric charge. Unlike electric charge which can be expressed by only one number, color charge is expressed by three numbers commonly labeled R, G, B, for red, green, and blue respectively. The Lagrangian for QCD is

$$\mathcal{L} = \sum_j \bar{\psi}_j (i \not{D}) \psi_j - \frac{1}{4} F_{\mu\nu}^a F_a^{\mu\nu}, \quad (2.1)$$

where

$$F_{\mu\nu}^a = \partial_\mu A_\nu^a - \partial_\nu A_\mu^a + g f^{abc} A_\mu^b A_\nu^c, \quad (2.2)$$

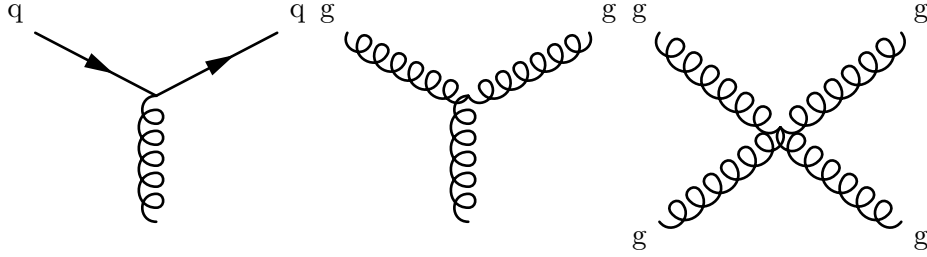


Figure 2.1: These are the Feynman Diagrams of the allowed QCD vertices. All interactions in the theory are built by combining these vertices. Quarks are represented by the solid lines and gluons are represented by curly lines. The three and four gluon vertices are unique to non-Abelian gauge theories.

and

$$\not{D} = D_\mu \gamma^\mu = (\partial_\mu - ig A_\mu^a \lambda^a) \gamma^\mu. \quad (2.3)$$

A^μ are the gluon fields, ψ_j is the j th quark flavor. Greek μ and ν are the usual spacetime indices while a , b , and c are color indices. The structure constant f^{abc} is given by the commutation relation between the generators of the group $[\lambda^a, \lambda^b] = \frac{i}{2} f^{abc} \lambda^c$ where λ^a are the Gell-Mann matrices.

The Lagrangian allows the three vertices shown in figure 2.1. The first vertex couples a fermion to a gluon and is analogous to diagrams in QED that couple an electron to the photon. The other two vertices are the three and four gluon vertices and have no analogue in QED. The gluon vertices are a result of the non-abelian nature of the theory.

QCD exhibits two interesting properties: confinement and asymptotic freedom. Confinement means that the force between two quarks does not diminish as they are pulled apart. In fact if you try to pull two quarks apart eventually there will be enough energy in the field to produce a new quark anti quark pair. As a result we never observe free quarks in nature, they only exist in colorless bound states called hadrons. Hadrons come in two varieties $q\bar{q}$ pairs called mesons and qqq triplets called baryons. Despite the fact that confinement is easy to demonstrate on the lattice [120], there is an outstanding Millennium Prize for an analytic proof[1]. Asymptotic freedom reflects that at very large momentum transfers the quarks and gluons interact weakly. The 2004 Nobel prize in physics was awarded for the discovery of asymptotic freedom by Frank Wilczek, David Gross, and David Politzer in 1973 [60, 101].

The strong nuclear force is responsible for the proton, composed of two up and one down

quark, and neutron, composed of two down and one up quark. Additionally, the strong nuclear force binds these particles together to form atomic nuclei. Most of the understood mass in the universe is a result of the strong nuclear force. In QCD the up and down quarks have masses on the order of an MeV while the proton has a mass of 938 GeV. Almost all of the proton mass results from its own binding energy.

The remaining $SU(2)_W \times U(1)_Y$ part of the standard model is the unified electroweak force [105, 116, 61, 73, 74, 44, 58]. This force accounts for the quantum theory of electrodynamics and the weak interactions. The electroweak force has three bosonic force carriers: the massless photon and the massive W^\pm and Z . As a result of being carried by the massless photon, electrodynamics is an infinite range force. However, the weak force has an interaction range of $\mathcal{O} \approx 10^{-16}\text{m}$ due to the heaviness of the W^\pm and Z . In addition to affecting the quark sector, the electroweak force also interacts with leptons. Leptons are spin 1/2 fermions that are not charged under the $SU(3)_C$ group. The electron, muon, tau, and their corresponding neutrinos are all leptons.

An essential feature of the electroweak theory is that $SU(2)_W \times U(1)_Y$ spontaneously breaks down to $U(1)_{EM}$, see section 2.2 for further details. This breaking gives the W^\pm and Z bosons their mass and also accounts for the standard model fermion masses. $U(1)_{EM}$ is the theory of quantum electrodynamics in which spin 1/2 fermionic matter charged with electric charge interact by exchanging photons. Unlike in QCD, photons are not charged under $U(1)$ and therefore don't interact at tree level. This yields a much simpler theory than QCD which can be understood perturbatively.

The Lagrangian for QED is similar to 2.1 however the field strength tensor $F_{\mu\nu}^a$ is replaced by $F_{\mu\nu}$, dropping the color indices. The simpler Abelian $U(1)$ symmetry also means that the field strength tensor is defined as

$$F_{\mu\nu} = \partial_\mu A_\nu - \partial_\nu A_\mu \quad (2.4)$$

Where A^μ is the four potential defined as $A^\mu \equiv (\phi, \vec{A})$. Only one vertex is allowed that couples a fermion to a photon.

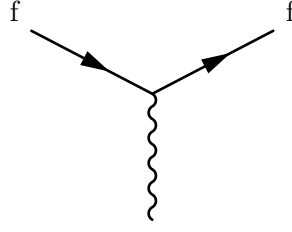


Figure 2.2: There is only one allowed vertex in QED which couples a fermion charged under $U(1)$ with the photon.

In the standard model, matter is divided into three generations. The first generation of matter is the most familiar. It consists of the up quark down quark, electron, and electron neutrino. All of the matter that we experience in our daily lives is made of this generation of matter. The other two generations of matter are essentially heavier replicas of the first generation. These heavier generations rapidly decay via the weak nuclear force to the first generation of matter and therefore are only detected in high energy events such as cosmic rays or particle physics experiments. The second generation consists of the strange quark, charm quark, muon, and muon neutrino. Finally the third consists of the top quark, bottom quark, tau, and tau neutrino. The properties of all the standard model particles is summarized in tables 2.1 and 2.1.

2.2 Higgs Mechanism

In the standard model the W^\pm and Z bosons are massive but the photon is massless. Since gauge invariance dictates that a mass term is forbidden in the Lagrangian, this seems to pose a problem. Disaster is averted because $SU(2)_W \times U(1)_Y$ is spontaneously broken to $U(1)_{EM}$. In the standard model this is accomplished by adding a complex scalar doublet field to the theory. The field is named the Higgs field after one of its discoverers. The following discussion shows how electroweak symmetry breaking is facilitated by the Higgs [102, 103, 98].

Lets introduce a complex elementary scalar doublet $\Phi = \begin{pmatrix} \phi^+ \\ \phi_0 \end{pmatrix}$ that transforms in the $(1, 2, \frac{1}{2})$ representation of $SU(3)_C \times SU(2)_L \times U(1)_Y$. We allow all terms in the Lagrangian that have mass

Particle Name	Symbol	Mass (MeV)	Generation	charge $(U(1)_{EM})$
electron	e	0.510999	1	-1
electron neutrino	ν_e	$< 2.2 \times 10^{-6}$	1	0
muon	μ	105.659	2	-1
muon neutrino	ν_μ	< 0.170	2	0
tau	τ	1.776×10^3	3	-1
tau neutrino	ν_τ	< 15.5	3	0

Table 2.1: The standard model fermions are spin 1/2 fundamental particles in the color singlet. The masses for the electron, muon, and tau are known to the precision given. The neutrinos are known to have masses but their masses are not precisely measured. Left handed leptons have weak hypercharge of -1 while right handed leptons have weak hypercharge of -2. Right handed neutrinos have not been observed.

Particle Name	Symbol	Mass (MeV/ c^2)	Generation	charge $U(1)_{EM}$
up	u	$2.3^{+0.7}_{-0.5}$	1	$\frac{2}{3}$
down	d	$4.8^{0.5}_{-0.3}$	1	$-\frac{1}{3}$
strange	s	95 ± 5	2	$-\frac{1}{3}$
charm	c	$(1.275 \pm 0.025) \times 10^3$	2	$\frac{2}{3}$
top (truth)	t	$(173.07 \pm 1) \times 10^3$	3	$\frac{2}{3}$
bottom (beauty)	b	$(4.180 \pm 0.03) \times 10^3$	3	$-\frac{1}{3}$

Table 2.2: The standard model quarks are spin 1/2 fundamental particles that are charged under $SU(3)_C$. Left handed quarks have 1/3 hypercharge. Right handed quarks with 2/3 electric charge have weak hypercharge of 4/3 while the right handed quarks with -1/3 electric charge have weak hypercharge of -2/3.

Particle Name	Symbol	mass (MeV)	charge $U(1)_{EM}$
gluon	g	0	0
photon	γ	0	0
W^+	W^+	$(80.385 \pm 0.015) \times 10^3$	1
W^-	W^-	$(80.385 \pm 0.015) \times 10^3$	-1
Z	Z	$(91.1876 \pm 0.0021) \times 10^3$	0
Higgs	H	$(125.9 \pm 0.4) \times 10^3$	0

Table 2.3: The standard model bosons. All particles are spin 1 except for the Higgs particle which is spin 0.

dimension ≤ 4 , thus ignoring irrelevant operators,

$$\mathcal{L}_H = (D_\mu \Phi)^\dagger (D^\mu \Phi) + \mu^2 \Phi^\dagger \Phi - |\lambda| (\Phi^\dagger \Phi)^2, \quad (2.5)$$

with covariant derivative

$$D_\mu \Phi = (\partial_\mu - ig A_\mu^a \tau^a - ig' Y B_\mu) \Phi. \quad (2.6)$$

A_μ^a and B_μ are the $SU(2)$ and $U(1)$ gauge bosons respectively. The coupling constant g belongs to $SU(2)$ and the coupling constant g' belongs to $U(1)$. Finally Y is the generator of hypercharge, τ^a are the generators of $SU(2)$ and are related to the Pauli matrices

$$\tau^a = \frac{1}{2} \sigma^a. \quad (2.7)$$

We can now identify the potential $V(\Phi)$ as

$$V(\Phi) = -\mu^2 \Phi^\dagger \Phi + |\lambda| (\Phi^\dagger \Phi)^2, \quad (2.8)$$

this potential is shown in figure 2.3. As long as μ^2 is positive the potential has a spontaneously broken symmetry. Gauge invariance allows us to choose the vacuum state to correspond to the vacuum expectation value (VEV)

$$\langle \Phi \rangle_0 = \frac{\nu}{\sqrt{2}} \begin{pmatrix} 0 \\ \nu \end{pmatrix}, \quad (2.9)$$

where $\nu = \sqrt{\frac{\mu^2}{|\lambda|}}$. Moreover, ν is related to the Fermi constant, G_F , by

$$\nu = \frac{1}{2^{1/4} \sqrt{G_F}} \approx 246 \text{ GeV}. \quad (2.10)$$

If we assign our theory to have hypercharge $Y = +\frac{1}{2}$, a complete gauge transformation in this theory is

$$\Phi \rightarrow e^{i\alpha^a(x)\tau^a} e^{i\beta(x)/2} \Phi. \quad (2.11)$$

Making the particular choice of $\alpha^1 = \alpha^2 = 0$ and $\alpha^3 = \beta$ we see that $\langle \Phi \rangle$ is invariant. Therefore the theory still contains an unbroken $U(1)$ symmetry which we identify with electromagnetism. This unbroken $U(1)$ symmetry contains one massless gauge boson which is the photon. The other three

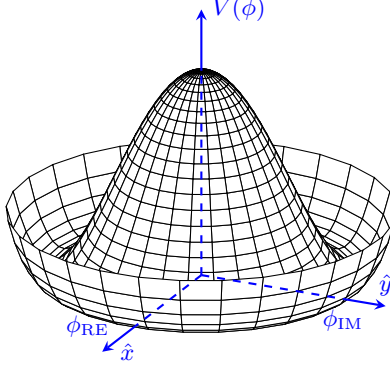


Figure 2.3: The Higgs potential with $\mu^2 > 0$ has a spontaneously broken global symmetry. We can think of this classically. The origin is a local maximum of potential energy and therefore an unstable equilibrium. If we place a particle at the origin any perturbation will push the particle ‘down hill’ to a minimum of the potential. Since there are a number of degenerate minima, choosing one spontaneously breaks the symmetry.

gauge bosons corresponding to the broken generators of the symmetry group become massive. The massive gauge bosons get their mass from the square of the kinetic term evaluated at the VEV

$$(D_\mu \Phi)^\dagger (D^\mu \Phi) = \frac{1}{2} \begin{pmatrix} 0 & \nu \end{pmatrix} \left(g A_\mu^a \tau^a + \frac{1}{2} g' B_\mu \right) \left(g A^{b\mu} \tau^b + \frac{1}{2} g' B^\mu \right) \begin{pmatrix} 0 \\ \nu \end{pmatrix}. \quad (2.12)$$

Substituting τ^a and taking the product we get

$$(D_\mu \Phi)^\dagger (D^\mu \Phi) = \frac{\nu^2}{8} \left[g^2 (A_\mu^1)^2 + g^2 (A_\mu^2)^2 + (-g A_\mu^3 + g' B_\mu)^2 \right]. \quad (2.13)$$

We can now perform a field redefinition and recover

$$W_\mu^\pm = \frac{1}{\sqrt{2}} (A_\mu^1 \mp A_\mu^2) \quad \text{with mass } m_W = g \frac{\nu}{2} \quad (2.14)$$

$$Z_\mu = \frac{1}{\sqrt{g^2 + g'^2}} (g A_\mu^3 - g' B_\mu) \quad \text{with mass } m_Z = \sqrt{g^2 + g'^2} \frac{\nu}{2} \quad (2.15)$$

$$A_\mu = \frac{1}{\sqrt{g^2 + g'^2}} (g' A_\mu^3 + g B_\mu) \quad \text{with mass } m_A = 0, \quad (2.16)$$

which are the standard model W^\pm, Z bosons and photon.

The change in basis from the original massless bosons to massive bosons and photon is

characterized by the weak mixing angle θ_w :

$$\begin{pmatrix} Z_\mu \\ A_\mu \end{pmatrix} = \begin{pmatrix} \cos \theta_w & -\sin \theta_w \\ \sin \theta_w & \cos \theta_w \end{pmatrix} \begin{pmatrix} A_\mu^3 \\ B_\mu \end{pmatrix}, \quad (2.17)$$

with

$$\cos \theta_w = \frac{g}{\sqrt{g^2 + g'^2}}, \quad \sin \theta_w = \frac{g'}{\sqrt{g^2 + g'^2}}. \quad (2.18)$$

When the W^\pm and Z bosons become massive they each gain an additional longitudinal degree of freedom. These new degrees of freedom aren't free; they come from the three components of the Higgs doublet Φ associated with the original broken symmetry. The Goldstone mechanism facilitates this transfer of degrees of freedom. The vernacular is that the massless Goldstone bosons of the original theory are “eaten” by the W^\pm and Z .

Additionally, the Higgs mechanism provides a mass for the standard model fermions. Fermion mass terms are forbidden in the standard model because they would violate local gauge invariance since left and right fermions transform differently. Now that we have a scalar field Φ we can couple it to our fermion with a Yukawa coupling, $-c\bar{f}_L\Phi f_R$. In such a term f_L and f_R are the right and left components of the fermion, c is an arbitrary coupling characterizing the strength of the interaction with the Higgs VEV, and Φ is the Higgs VEV. Heavy particles like the top have a large value of c while light particles like the up quark have a small value of c .

Finally, the Higgs Boson, a particle excitation of the Higgs field, was recently discovered at the Large Hadron Collider (LHC). The discovery was announced to the public July 4, 2012. Current experimental results pin the Higgs mass at 126 GeV. Peter Higgs and Francois Englert shared the Nobel prize in 2013 for their work in developing the theory of EWSB through the Higgs mechanism.

2.3 Beyond Standard Model Physics

While the Standard Model is the pinnacle of our current understanding of the universe it is not the last word on the subject. There are many phenomena that the standard model does not

explain. Furthermore there are theoretically troubling aspects to the standard model that leaves more to be desired.

Perhaps the biggest omission in the standard model is a quantum field theory of Gravity. Our current understanding of gravity is Einstein's theory of General Relativity (GR). GR is a thoroughly verified theory and famously explained the perihelion of mercury and the deflection of light rays massive objects. Unfortunately we don't have a testable theory that combines general relativity with quantum mechanics, and to date a force carrying boson, the graviton, has eluded detection. A related problem is that gravity is so much weaker than the other three fundamental forces.

Another unexplained phenomenon is neutrino oscillations. In nature we have observed the flavor of neutrinos to change. These oscillations prove that the neutrino's are not massless particles as the standard model requires.

Finally we currently understand from astrophysical observations that the standard model only accounts for 4% of the energy budget of the universe. Dark matter, which has eluded detection, composes of another 21% of the energy budget. Dark Energy, which is responsible for the acceleration of the expansion of the universe, composes 75% of our energy budget. Taking this view the standard model actually explains a relatively small part of the universe.

In addition to not explaining the entire range of physical phenomena that we observe (or fail to observe) there are theoretical shortcomings that indicate a lack of complete understanding. The first indication that something is awry is the number of parameters. The standard model in its most general form has nineteen free parameters. All of these parameters are fixed by experiment and cover a large range of scales. We expect that a more nuanced understanding of the universe will require less parameters.

Another theoretical quirk of the standard model is the strong CP problem [42]. Unlike the electroweak interactions, QCD does not violate CP symmetry. CP symmetry is the symmetry of charge conjugation (C) and parity (P). The QCD Lagrangian allows terms that violate CP symmetry but these terms appear to be zero which is a type of fine tuning.

Finally there is the hierarchy problem. In the standard model the coupling constants of the theory change with energy scale. In principle, the standard model is valid to the Plank scale. However, if we take the cutoff of the theory to be the Plank scale, it is removed from several orders of magnitude from the weak scale. This is a problem because the Higgs is an elementary scalar field and therefore has a relevant quartic self coupling. The renormalized Higgs mass is

$$m_H^2 = m_0^2 + \frac{3}{4\pi}\lambda\Lambda^2, \quad (2.19)$$

where m_0 is the bare Higgs mass, λ is the quartic coupling, and Λ is the cutoff scale. The renormalized mass has a quadratically divergent additive renormalization to its mass that is proportional to the cutoff. To keep the renormalized Higgs mass at its physical value an unnatural degree of fine tuning to its bare mass, m_0 , and quartic term, λ , is required. A more natural solution to this problem is for Λ to be at a scale similar to the Higgs mass.

These issues have pushed researchers to look for extensions or alternatives to the standard model that answer one or more of these unresolved questions. The state of affairs for almost 40 years was that we knew there had to be a Higgs mechanism to complete the standard model. We also had a very good guess where to look for the Higgs boson because of the mass of the W^\pm and the Z .

Most methods of solving the hierarchy problem introduce new physics at the electroweak scale. This moves the SM cutoff much closer to the electroweak scale, removing the need for fine tuning. Because the Tevatron and more recently the LHC could reach energies high enough to probe the electroweak scale it was possible to build and test falsifiable theories.

Supersymmetry is one such approach. It postulates a symmetry between fermions and bosons. This symmetry results in a cancellation of the quadratic divergence that the scalar Higgs theory has. Another approach which I will describe further in the next section is for the Higgs particle to be a composite particle. A consequence of all of these theories is new particles. Now we know the value of the Higgs mass. Additionally, no other particles have yet been discovered by the LHC. This has proven to be a deadly combination, ruling out the simplest incarnations of these models.

2.4 Composite Higgs

One solution to the hierarchy problem is for the Higgs to be a composite composed of particles from a new strongly interacting sector [81, 55, 117, 112, 118, 43, 41]. This new sector is responsible for electroweak symmetry breaking. There are many types of theories that use such a modus operandi. While these theories have the benefits of solving the hierarchy and fine tuning problems with the standard model Higgs they typically favor a heavier Higgs mass than what has been observed and pose other theoretical challenges that I will elaborate on below. In the next three subsections I will briefly discuss technicolor, extended technicolor, introduce the conformal window, and discuss the state of Lattice endeavors in this area.

2.4.1 Technicolor

Technicolor seeks to replace the Higgs sector of the electroweak theory with a new strongly interacting sector. In this section I will show how this is accomplished. Some good reviews on the subject can be found in references [75, 85, 108, 90, 106] For simplicity lets consider a $SU(3)_{TC}$ gauge group that couples to a pair of massless fermions $\Psi = (\psi_1, \psi_2)$ in the fundamental representation. It is worth noting that so far this theory is identical to QCD with massless up and down quarks. The Lagrangian for our theory is

$$\mathcal{L}_{TC} = -\frac{1}{4}F_{\mu\nu}^a F^{\mu\nu,a} + \bar{\Psi}_L(i\not{D})\Psi_L + \bar{\Psi}_R(i\not{D})\Psi_R. \quad (2.20)$$

Recall that fermions have a right handed and left handed component $\psi = (\psi_L, \psi_R)$. We know that such a theory possesses a global $SU(2)_L \times SU(2)_R$ symmetry and that this symmetry is spontaneously broken to the vector subgroup $SU(2)_V$. This symmetry manifests itself in the nonzero VEV of the chiral condensate

$$\langle \bar{\Psi}\Psi \rangle = \langle \bar{\Psi}_L\Psi_R + \bar{\Psi}_R\Psi_L \rangle \neq 0. \quad (2.21)$$

To see how spontaneous chiral symmetry breaking translates to spontaneous electroweak symmetry breaking we have to use chiral perturbation theory [54, 59]. Chiral perturbation the-

ory is an effective field theory whose fundamental degrees of freedom are the Goldstone bosons associated with the symmetry breaking. In our case the broken $SU(2)$ symmetry has 3 degrees of freedom and therefore our effective theory will have three massless Goldstone bosons. Continuing our analogy to QCD, these bosons are the pions.

The chiral Lagrangian is nonlinear and possesses an infinite number of terms. The standard approach is to perform an expansion about momentum, p , that are small with respect to the cutoff Λ . We are then free to choose what order in $\frac{p}{\Lambda}$ we work with. The lowest order in the expansion is

$$\mathcal{L}_\chi = \frac{F^2}{4} \text{Tr} \left[(D^\mu U)^\dagger (D_\mu U) \right]. \quad (2.22)$$

U is a non-linear function of the Goldstone fields ϕ_a

$$U \equiv e^{\sigma^a \phi_a \frac{2i}{f}}. \quad (2.23)$$

The derivative D_μ is coupled to the electroweak and gauge field because our technifermions are charged under $SU(2) \times U(1)$

$$D_\mu = \partial_\mu - ig \frac{\sigma^a}{2} A_\mu^a + ig' \frac{\sigma^3}{2} B_\mu. \quad (2.24)$$

We can substitute our derivative into our lowest order effective Lagrangian and find that

$$\mathcal{L}_\chi = \frac{F^2}{4} \text{Tr} \left[\frac{g^2}{4} \left[\left(A_\mu^1 - \frac{4}{Fg} \partial_\mu \phi_1 \right)^2 + \left(A_\mu^2 - \frac{4}{fg} \partial_\mu \phi_2 \right)^2 + \left(A_\mu^3 - \frac{g'}{g} B_\mu - \frac{4}{fg} \partial_\mu \phi_3 \right)^2 \right] \right]. \quad (2.25)$$

We can now make the field redefinitions

$$\begin{aligned} W_\mu^{1,2} &\equiv A_\mu^{1,2} - \frac{4}{fg} \partial_\mu \phi_{1,2} \\ Z_\mu &\equiv \frac{g}{\sqrt{g^2 + g'^2}} \left(A_\mu^3 - \frac{g'}{g} B_\mu - \frac{4}{fg} \partial_\mu \phi_3 \right) \\ A_\mu &\equiv \frac{g}{\sqrt{g^2 + g'^2}} \left(\frac{g'}{g} A_\mu^3 + B_\mu \right). \end{aligned} \quad (2.26)$$

The field redefinition completely removes ϕ_a from our theory to first order. The field A_μ is the massless photon field. Our Lagrangian now reads

$$\mathcal{L}_\chi = \mathcal{L}_{\text{kinetic}} + \frac{1}{2} m_W^2 [(W_\mu^1)^2 + (W_\mu^2)^2] + \frac{1}{2} m_Z^2 Z_\mu^2, \quad (2.27)$$

where the masses for the heavy bosons are

$$\begin{aligned} m_W &= \frac{1}{2} F g \\ m_Z &= \frac{1}{2} F \sqrt{g^2 + g'^2}. \end{aligned} \tag{2.28}$$

With the substitution $F = \nu$ we reproduce m_W and m_Z that we found in section 2.2 exactly! This is quite a shock, as I mentioned earlier so far everything we are doing is simply QCD with 2 flavors of massless quarks, why did we introduce the Higgs field in the first place? Simply put the effect in QCD is much too small. In QCD, $F = F_\pi \approx 93$ MeV, this is more than a factor of a thousand too small to account for the observed W^\pm and Z masses.

Despite the fact that spontaneous chiral symmetry breaking in QCD is too small to account for the observed electroweak sector, there is no reason that we can't introduce a new strongly interacting theory where the equivalent $F = \nu \approx 246$ GeV. Furthermore we can choose a different gauge group, number of fermions, and representation if we wish. The only constraint is that chiral symmetry is broken and the results are phenomenologically consistent with experimental observation. It is worth noting that more flavors of fermions will produce additional pions that will acquire masses. In general $SU(N_f)_L \times SU(N_f)_R \rightarrow SU(N_f)_V$ will produce $N_f^2 - 1$ Goldstone bosons, three can be eaten to form the W^\pm and Z and the remaining $N_f^2 - 4$ become massive pseudo-Goldstone bosons. The masses of the pseudo-Goldstone bosons in a viable theory need to be large enough to not have been discovered yet.

In summary we have shown how strong dynamics can explain electroweak symmetry breaking. Because of asymptotic freedom, such a theory does not have UV divergences, solving the fine tuning problem. One failure of our model is that we have lost an explanation for the fermion masses. It is possible to construct an effective theory of fermion masses using a four fermion operator and adjusting the couplings for each quark and lepton flavor by hand. Unfortunately a four-Fermi interaction is non-renormalizable, for a UV complete technicolor that gives fermion masses we need extended Technicolor.

2.4.2 Extended Technicolor

In order for a theory of electroweak symmetry breaking to replace the Higgs mechanism it must also generate masses for the quarks and leptons of the standard model. This is accomplished through a framework called extended technicolor (ETC)¹. Under this framework an extended technicolor group is introduced $SU(3 + N_{TC})_{ETC}$. This extended group is also asymptotically free and therefore UV complete. Both the usual standard model content as well as $SU(N_{TC})$ are charged under ETC. At some scale $\Lambda_{ETC} \gg \nu$ the ETC group spontaneously breaks to $SU(N)_{TC} \times SU(3)_c \times SU(2)_L \times U(1)_Y$. The gauge bosons of ETC theories have masses $M_{ETC} \approx g_{ETC}\Lambda_{ETC}$. Because of the wide range of masses for standard model fermions, most models ETC models spontaneously break at multiple scales. Most models have three breaking scales, corresponding to the three generations of quarks and leptons. ETC is an ambitious theory, a successful model of ETC will solve almost all the outstanding theoretical issues with the SM. Both the flavor hierarchy problem and the number of free parameters in the standard model would be solved dynamically. It is not surprising then that no ‘reasonable’ ETC model has been produced yet.

Extended technicolor faces many practical problems. One major hurdle is flavor changing neutral currents (FCNC). Flavor changing neutral currents such as $\mu \rightarrow e\bar{e}e$ and mixing between neutral mesons are highly constrained in the standard model. Most ETC theories typically introduce FCNC that are proportional to M_{ETC}^2 .

Another issue with ETC theories is size of the top quark mass. The top quark is so massive, $m_t = 172$ GeV, that its associated ETC scale is $M_{ETC} \approx 3$ TeV. This is problematic because it is so close to the electroweak scale. Additionally incorporating both the top and the bottom with the same ETC breaking would require an accompanying large isospin breaking. These issues have forced many models to give the top special treatment in so called top assisted ETC.

Skirting these and other issues without producing particles light enough to have already been discovered is a daunting task. There are many models that have been able to trade one

¹ The first rule of ETC is no scalar fields, the second rule of ETC is no scalar fields.

defect for another, but to date no complete model exists. There is a general consensus that any viable ETC model probably will have a slowly running (walking) coupling. In a walking theory $\gamma(\alpha(\mu))$ is large between Λ_{TC} and M_{ETC} . It has been shown that walking behavior addresses all of the problems I have discussed to a point. However any model that breaks at multiple scales and has the peculiarities of walking between those scales is likely to be very baroque. I discuss how a walking theory can be generated in more detail below.

2.4.3 Conformal Window

A general class of strongly interacting theories that are of interest for Technicolor and extended Technicolor dynamics are Yang-Mills gauge theories [32, 34]. Pure gauge Yang-Mills theories have $SU(N_C)$ interactions while more general theories include N_f flavors of fermions in some representation. At low energies QCD is effectively a $SU(3)$ gauge theory with $N_f=2$. This description includes only the up and down quarks which are nearly massless and respect an approximate $SU(2)$ isospin symmetry. The most general Yang-Mills Lagrangian with a $SU(N_c)$ local gauge symmetry and N_f flavors of massless fermions in a representation R is:

$$\mathcal{L}_{YM} = -\frac{1}{4g^2} \sum_{a=1}^{N_c} F_{\mu\nu}^a F^{\mu\nu,a} + \sum_{i=1}^{N_f} \bar{\Psi}_i (i\not{D}) \Psi_i. \quad (2.29)$$

$F_{\mu\nu}^a$ is the field strength tensor, shown here for an arbitrary group

$$F_{\mu\nu}^a = \partial_\mu A_\nu^a - \partial_\nu A_\mu^a + gf^{ijk} A_\mu^i A_\nu^j, \quad (2.30)$$

where A_μ^a is the field, a is a group index and the structure coefficient f^{ijk} is determined by the commutation relationship

$$[T^i, T^j] = if^{ijk} T^k, \quad (2.31)$$

and T^a are the generators of the group. The Lagrangians may look similar however their behavior in both UV and IR can be dramatically different from theory to theory. While the Lagrangian is classically scale invariant; the quantum theory is not. This is understood through the beta function,

$$\beta(g^2) = -\mu \frac{\partial g^2}{\partial \mu}. \quad (2.32)$$

Representation	$\dim(R)$	$T(R)$	$C_2(R)$
F	N	$\frac{1}{2}$	$\frac{N^2-1}{2N}$
S_2	$\frac{N(N+1)}{2}$	$\frac{N+2}{2}$	$\frac{(N+2)(N-1)}{N}$
A_2	$\frac{N(N-1)}{2}$	$\frac{N-2}{2}$	$\frac{(N-2)(N+1)}{N}$
G	$N^2 - 1$	N	N

Table 2.4: This table summarizes some commonly used groups, their dimension, first Casimir invariant, and second Casimir invariant. F, S_2, A_2 and G are the fundamental, 2-index symmetric, 2-index antisymmetric, and adjoint representations respectively.

The beta function describes how the gauge coupling evolves as the renormalization scale μ is changed. This function can be expanded in perturbation theory and is universal to two loops:

$$\beta(g^2) = -\frac{b_1}{16\pi^2}g^4 - \frac{b_2}{(16\pi^2)^2}g^6 \quad (2.33)$$

Any terms beyond two loops are renormalization scheme dependant and are not relevant for the following discussion. The values for b_1 and b_2 are:

$$\begin{aligned} b_1 &= \frac{11}{3}N_c - \frac{4}{3}N_f T(R) \\ b_2 &= \frac{34}{3}N_c^2 - \frac{4}{3}T(R)N_f \left[5N_c + 3C_2(R) \right]. \end{aligned} \quad (2.34)$$

$T(R)$ and $C_2(R)$ are the first and second Casimir invariants and depend on the representation R of the group. Table 2.4.3 give these invariants for a few common representations.

Clearly, tuning N_c , N_f and R offers a great deal of freedom in specifying the gauge theory. We can see from equation 2.34 that if the coefficients b_1 and b_2 are both positive than the beta function is negative. In this scenario the theory asymptotically free, confining, and spontaneously breaks chiral symmetry. The dynamics in the IR will be strongly coupled and non-perturbative. Figure 2.4 shows what this scenario looks like. This occurs in QCD ($SU(3)$ $N_f = 2$ or 3 depending on your treatment of the light quarks), which is an example of a theory where the beta function is negative.

If we keep b_1 positive and allow b_2 to become negative we can force the two terms to compete. In perturbation theory this occurs at a critical value of fermions $N_f^{crit} < \frac{17N_c^2}{2T(R)[5N_c+C(R)]}$. Such a beta function would start out negative, then at some coupling it would pass through a local minimum after which it would grow. Eventually, at some g_*^2 , the beta function would have a zero where $\beta(g_*^2) = 0$. Perturbatively this zero occurs at $g_*^2 = -b_1/b_2$ and is called a Banks-Zaks infrared fixed point. This is illustrated in figure 2.5, such a theory is governed by the conformal dynamics at the infrared fixed point and is scale invariant. If the fixed point is at very weak coupling it is possible to gain insights from perturbation theory, however many theories are known to have strongly coupled IRFP's where insights from perturbation theory are unreliable. Conformal theories, like the scenario described above, do not support bound states of particles. As such they are often referred to as ‘unparticle theories’.

Allowing both b_1 and b_2 to be negative results in a trivial theory. Such a theory, shown in figure 2.6 is not asymptotically free. The beta function is very similar to that of QED. Perturbatively this occurs when $N_f < \frac{11N_c}{4T(R)}$.

The region in theory space between where the IRFP appears and where asymptotic freedom is lost is referred to as the conformal window. Limiting our consideration to theories using the fundamental representation fermions and $N_c = 3$ we see that perturbation theory predicts the conformal window is between $N_f \approx 9.4$ and $N_f = 16.5$. It is important to note that all of the discussion in this section are perturbative results. On the upper end of the conformal window before asymptotic freedom is lost, the IRFP of many theories is weak enough that perturbation theory is reasonable. However, as we pass through the conformal window and consider theories with less fermionic degrees of freedom, the IRFP becomes strongly coupled. Here perturbation theory is not to be trusted. The location of the lower bound of the conformal window is an area of active research.

The conformal window itself is not a very interesting system to study but understanding the lower bound of the conformal window is of great interest for potential Technicolor theories. Right below the conformal window it is believed that a walking theory can exist. In such a theory,

shown in figure 2.7 the beta function would start out like the conformal scenario. However, as the beta function approaches the IRFP, chiral symmetry is spontaneously broken and the beta function would turn around away from the ‘would be’ IRFP. If this occurs, the coupling g will run very slowly (walk) over a wide range of scales where the β function approaches zero. Walking is necessary to achieve a wide separation of scales in Technicolor theories without generating flavor changing neutral currents.

2.4.4 Lattice Studies

Today there is a rich ecosystem of lattice studies of BSM studies, a survey of the current state of the field can be found in [?]. It is extremely important that different ways of probing for IR conformality on different lattice discretization schemes converge to common conclusions. In general this is accomplished but there are still systems that cause debate because different groups draw different conclusions from their findings. Below I summarize the state of the field both in terms of calculation types and methods of theories studied.

Many studies that seek to locate the conformal window have proceeded by calculating the RG flow of the theory under consideration. The first example of this type of study is [10]. If an IRFP is found then the theory is conformal while a lack of an IRFP at strong coupling indicates that chiral symmetry is broken and the theory confines. One benefit of running coupling calculations is that they are not as computationally intensive as many other alternatives. There are multiple types of calculations that fall under this umbrella. Schrödinger functional, MCRG (section ??), and now the Wilson flow step scaling (section 5) are examples.

Another common method uses lattice spectroscopy calculations. In these methods the masses of light bound states are measured at a variety of input quark masses. The fit of the bound state as a function of the bare quark mass will determine if the theory is best described by a chirally broken theory or conformal one. References [78, 79, 45] are examples of studying the conformal window using spectroscopy. In practice these fits are extremely difficult, in particular fits for volume squeezed conformal theories has proved troublesome.

Several groups have studied the finite temperature phase diagram [77, 107]. These studies identify an IRFP with a deconfining bulk transition. Additionally techniques using the Dirac eigenvalue spectrum [27] has proven useful in calculating the mass anomalous dimension over a wide range of scales.

The majority of calculations have been for theories with $N_c = 3$ with fermions in the fundamental representation. This is in part because the wide availability of optimized code bases for three colors. Within these studies staggered fermions, a formulation of lattice fermions that has four doublers and preserves a remnant of chiral symmetry, are commonly employed. As a consequence theories with $N_f = 4, 8, 12, 16$ are most widely studied although others have been studied as well. The consensus is that $N_f \leq 8$ are confining, $N_f = 12$ is probably conformal, and $N_f = 16$ is conformal.

There is also a growing number of calculations with $SU(2)$. The two flavor model in the adjoint representation is widely studied and is clearly in the conformal window [23, 37, 36, 35]. Studies with fundamental fermions exist for $N_f = 2, 4, 6, 8, 10$ with the conformal window being identified as existing between $4 < N_f < 8$. Studies of $SU(3)$ and $SU(4)$ in the 2-index symmetric representation also exist.

Clearly, beyond standard model physics is inherently a difficult subject to study on the lattice. Unlike studies of QCD we don't know the answer before hand. Taking a holistic approach and finding consensus is vital to our understanding. Additionally, as more of these theories have been studied, vital improvements have been made. This is true of the lattice code base, which prior to BSM studies was highly optimized and sometimes only available for QCD simulations. This is also true for improvements in how we analyze lattice data. I will discuss two improvements we have developed in chapters ?? and 6. Chapter ?? introduces an improvement to the gradient flow method for calculating the renormalized step scaling function. Chapter 6 introduces an improvement to the traditional MCRG method of calculating the bare step scaling function. Finally work in BSM physics on the lattice has pushed forward our knowledge of formulating lattice actions and understanding the nature of lattice artifacts. Many of the theories that have been studied are sufficiently

un-QCD like that lattice QCD intuitions are being upended with a new tribal knowledge.

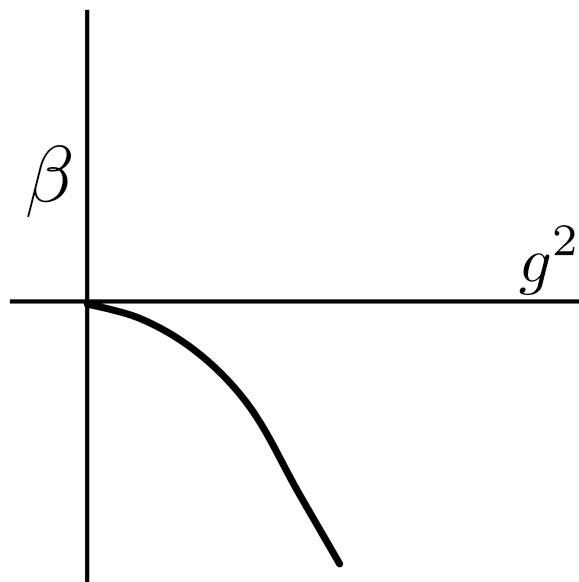


Figure 2.4: In a confining theory, such as QCD, the running of the coupling starts out negative. As g^2 grows the running becomes increasingly negative. Chiral symmetry is broken and the theory is described by confined bound states.

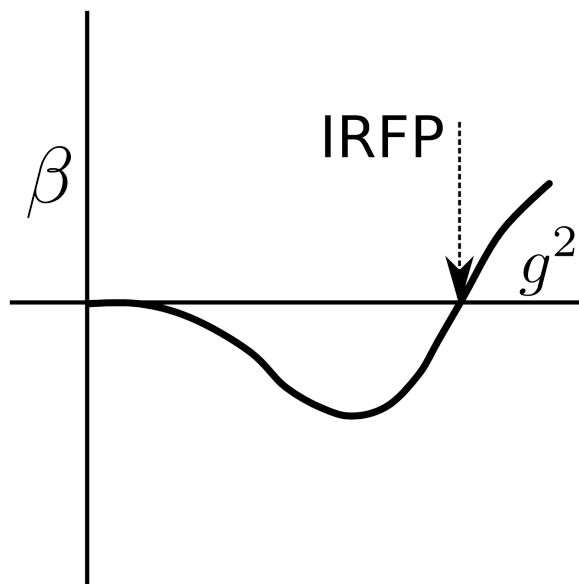


Figure 2.5: In a conformal field theory the β function starts out negative but turns around and crosses the $\beta = 0$ axis. At this crossing the theory has an infrared fixed point. The renormalized flow for the theory will run to the IR fixed point and the theory is conformal.

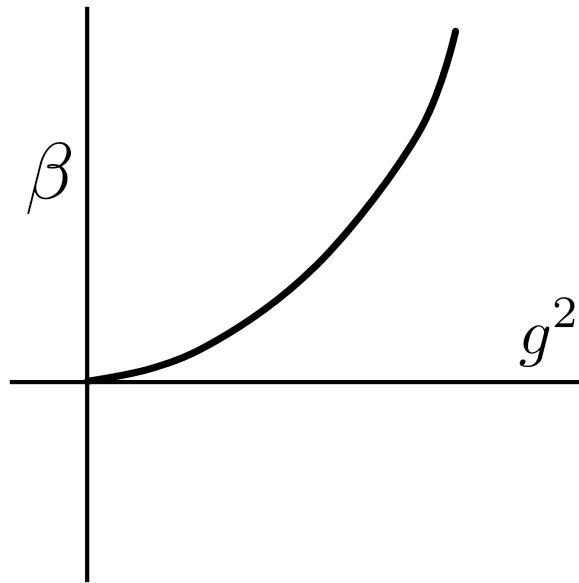


Figure 2.6: If the fermionic degrees of freedom overwhelm those of the bosons the beta function will start positive. This is the loss of asymptotic freedom. The beta function for such a theory resembles that of QED.

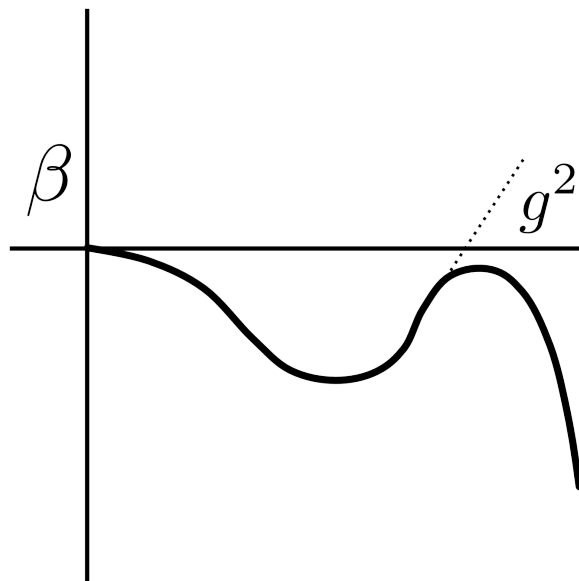


Figure 2.7: In a walking theory the beta function looks similar to the conformal picture of figure 2.5. The crucial difference is that before the β function develops a zero, shown by the dashed line, chiral symmetry is spontaneously broken driving the beta function back down. Near the would be IRFP where the dashed line has a zero, the coupling runs very slowly and is said to be walking.

Chapter 3

Lattice Field Theory

3.1 The Lattice

Lattice gauge theory is a method to regularize a quantum gauge theory nonperturbatively. To date it is the only controlled way to study strongly coupled systems from first principals. Lattice methods have been extremely successful in studying the strong nuclear force and many Lattice QCD results are now at the sub percent accuracy. As a mature field there are many excellent books on the subject that go above and beyond the level of rigor that I will pursue in the following discussion [38, 110, 29]. My aim is to provide some technical details relevant to the calculations that I perform in later chapters and give the reader a flavor of the subject.

The lattice regularizes the theory by discretizing Euclidean spacetime onto a regular grid of sites which are connected by links. The most common discretization is that of a four dimensional hypercube. For simplicity I will assume that all the sides of our hypercube are of length L (volume L^4), although this is not necessary and is in fact not true of many lattice calculations.

3.2 Gauge Fields

3.2.1 Pure Gauge

The simplest quantum field theory to put on the lattice is a pure Yang-Mills gauge theory. In group theory these are $SU(N_C)$ gauge theories without any pesky fermions. The most straight

forward discretization of a pure gauge action is in terms of the plaquette

$$P_{x,\mu\nu} = \text{Tr} \left[U_{x,\mu} U_{x+\hat{\mu},\nu} U_{x+\hat{\nu},\mu}^\dagger U_{x,\nu}^\dagger \right], \quad (3.1)$$

where $\mu \neq \nu$ and the trace over color and spin indices is implied. The plaquette is the smallest closed loop and therefore the smallest gauge invariant observable, corresponding to the links multiplied around a unit square of the lattice. The Hermitian conjugate, $P_{x,\mu\nu}^\dagger$, is the same loop with the order of the link products in the opposite direction.

To illustrate how the action is built out of plaquettes let's consider an abelian gauge theory for simplicity. For an abelian gauge theory

$$F_{\mu\nu} = \partial_\mu A_\nu - \partial_\nu A_\mu \quad (3.2)$$

$$\mathcal{L} = \frac{1}{4} F_{\mu\nu} F^{\mu\nu} \quad (3.3)$$

$$S = \int \mathcal{L} d^4x = \int \frac{1}{4} F_{\mu\nu} F^{\mu\nu} \quad (3.4)$$

Introducing the lattice spacing a we have link variables

$$U_{x,\mu} = e^{igaA_\mu(x)}. \quad (3.5)$$

Here g is the charge, a is the lattice spacing, and A_μ is the vector potential along the link. Note that the link variable is gauge invariant under the transformation

$$U_{x,\mu} \rightarrow \Omega_x U_{x,\mu} \Omega_{x+a\hat{\mu}}. \quad (3.6)$$

Although not necessary for the abelian case, considering $A_\mu(x)$ to have an underlying group structure will help us generalize this discussion at the end. Therefore using the identity $\exp(A)\exp(B) = \exp(A + B + \frac{1}{2}[A, B] + \dots)$ for matrices A and B , we can now construct the plaquette

$$\begin{aligned} P_{x,\mu\nu} = & e^{ig} \exp \left[aA_\mu(x) + aA_\nu(x + \hat{\mu}) - aA_\mu(x + \hat{\nu}) - aA_\nu(x) \right. \\ & - \frac{a^2}{2} [A_\mu(x), A_\nu(x + \hat{\mu})] - \frac{a^2}{2} [A_\mu(x + \hat{\nu}), A_\nu(x)] \\ & + \frac{a^2}{2} [A_\nu(x + \hat{\mu}), A_\mu(x + \hat{\nu})] + \frac{a^2}{2} [A_\mu(x), A_\nu(x)] \\ & \left. + \frac{a^2}{2} [A_\mu(x), A_\mu(x + \hat{\nu})] + \frac{a^2}{2} [A_\nu(x + \hat{\mu}), A_\nu(x)] + O(a^3) \right] \end{aligned} \quad (3.7)$$

Taylor expanding the shifted terms yields

$$A_\nu(x + \hat{\mu}) = A_\nu(x) + a\partial_\mu A_\nu(x) + O(a^2), \quad (3.8)$$

so that the plaquette becomes

$$\begin{aligned} P_{x,\mu\nu} &= e^{ig} \exp \left[a^2 (\partial_\mu A_\nu(x) - \partial_\nu A_\mu(x) + [A_\mu(x), A_\nu(x)]) + O(a^3) \right] \\ &= \exp \left[ig a^2 F_{\mu\nu}(x) + O(a^3) \right] \\ &= 1 + ig a^2 F_{\mu\nu}(x) - \frac{g^2 a^4}{2} F_{\mu\nu}(x) F^{\mu\nu}(x) + O(a^2) \end{aligned} \quad (3.9)$$

In the last step we Taylor expand about a . We can take the trace, note that $F_{\mu\nu}$ is traceless,

$$\text{Tr}[P_{x,\mu\nu}] = \text{Tr} \left[1 - \frac{g^2 a^4}{2} F_{\mu\nu}(x) F^{\mu\nu}(x) + O(a^2) \right]. \quad (3.10)$$

From this relationship it is clear that we can construct the following action from the plaquette

$$S_G(P_{x,\mu\nu}) = \frac{2}{g^2} \sum_x \sum_{\mu < \nu} (1 - P_{x,\mu\nu}). \quad (3.11)$$

As we take the continuum limit, $a \rightarrow 0$, we recover the continuum action

$$S \rightarrow \int d^4x \frac{1}{4} F_{\mu\nu} F^{\mu\nu}. \quad (3.12)$$

This can be trivially generalized to $SU(N_c)$ gauge theories

$$S_G(P_{x,\mu\nu}) = \beta \sum_x \sum_{\mu < \nu} \text{Re} \left[\text{Tr} \left[\mathbb{I} - \frac{1}{N_c} P_{x,\mu\nu} \right] \right], \quad (3.13)$$

where $\beta \equiv \frac{2N_c}{g^2}$. This is the Wilson gauge action [120].

There are several ways to improve upon the Wilson gauge action. The general idea is to add additional gauge invariant terms into the action with the goal of removing higher order lattice discretization errors [114, 119, 89]. There are many improved actions that have been developed in the past few decades. Although improved gauge actions in general are beyond the scope of this thesis, I will briefly comment on a negative adjoint term that we use in our action.

3.2.2 Negative Adjoint Term

In our studies we have added a negative adjoint term to the gauge part of the action. The purpose of this term is to avoid a well known spurious UV fixed point [17, 16, 19] in the fundamental adjoint plane show in 3.1. The mixed fundamental-adjoint action is

$$S_{AF} = \beta_F \sum_x \sum_{\mu < \nu} \text{Re} \left[\text{Tr} \left[\mathbb{I} - \frac{1}{N_c} P_{x,\mu\nu} \right] \right] + \beta_A \sum_x \sum_{\mu < \nu} \left[\text{Tr} \left[\mathbb{I} - \frac{1}{N_c^2} P_{x,\mu\nu}^\dagger P_{x,\mu\nu} \right] \right] \quad (3.14)$$

In our studies we set the ratio $\frac{\beta_A}{\beta_F} = -0.25$ which avoids a crossover that extends from the UV fixed point [64]. With our action the perturbative relation to the bare coupling is

$$\frac{6}{g^2} = \beta_F + 2\beta_A. \quad (3.15)$$

3.3 Fermionic Fields

3.3.1 Naive Fermions

Fermions prove to be a much thornier issue and great care must be taken with their discretization. Lets see what happens when we take a theory of one free fermion

$$S_f = \int d^4x \left(\bar{\psi}(x) \gamma^\mu \partial_\mu \psi(x) + m \bar{\psi}(x) \psi(x) \right), \quad (3.16)$$

and define the discretized lattice derivative to be the difference

$$\partial_\mu \psi(x) \rightarrow \Delta_\mu \psi_x = \frac{\psi_{x+\hat{\mu}} - \psi_{x-\hat{\mu}}}{2a}. \quad (3.17)$$

The fermionic fields ψ have the following gauge transformation:

$$\begin{aligned} \psi_x &\rightarrow \Omega(x) \psi_x \\ \bar{\psi}_x &\rightarrow \bar{\psi}_x \Omega^\dagger(x). \end{aligned} \quad (3.18)$$

We will now use the link variables $U_{x,\mu}$ to transport the ψ fields between neighboring sites

$$S_f = \frac{1}{2} \sum_{x,\mu} \bar{\psi}_x \gamma^\mu [U_{x,\mu} \psi_{x+\hat{\mu}} - U_{x-\hat{\mu}}^\dagger \psi_{x-\hat{\mu}}] + m \sum_x \bar{\psi}_x \psi_x. \quad (3.19)$$

Now we can construct the propagator in momentum space

$$S(p) = \frac{m - i \sum_{\mu} \gamma_{\mu} \sin(p_{\mu}a)/a}{m^2 + \sum_{\mu} \sin^2(p_{\mu}a)/a^2}, \quad (3.20)$$

where the momenta are restricted by the lattice discretization $p_{\mu} = \frac{2\pi i x_{\mu}}{N_{\mu}}$ for a lattice of N_{μ} sites in direction μ . For small momenta, $p_{\mu} \approx (0, 0, 0, 0)$ we can Taylor expand $\gamma^{\mu} \sin(p_{\mu}) \approx \not{p}$ and recover the continuum fermionic propagator $\frac{1}{\not{p} - m}$. Unfortunately we also recover the continuum propagator for fifteen other values of the four lattice momenta $p_{\mu} \approx (\pi, 0, 0, 0), p_{\mu} \approx (0, \pi, 0, 0), \dots, p_{\mu} \approx (\pi, \pi, \pi, \pi)$. Thus instead of describing one species of fermion, our theory instead describes 16 degenerate species of fermions. This is the fermion doubling problem. For each dimension d of the lattice we get 2^d species of fermions.

3.3.2 Wilson Fermions

Wilson solved the fermion doubling problem by adding a second derivative term to the action,

$$\Delta_{\mu}^2 \psi_x = \frac{\psi_{x+\hat{\mu}} - 2\psi_x + \psi_{x-\hat{\mu}}}{2a}. \quad (3.21)$$

Such a term does not affect our continuum extrapolation. The resulting action

$$S_f = -\frac{1}{2} \sum_{x,\mu} \bar{\psi}_x \left[(1 - \gamma^{\mu}) U_{x,\mu} \psi_{x+\hat{\mu}} + (1 + \gamma^{\mu}) U_{x-\hat{\mu}}^{\dagger} \right] + (4 + m) \sum_x \bar{\psi}_x \psi_x, \quad (3.22)$$

has the following free momentum space propagator

$$\begin{aligned} \frac{1}{a} S(p) &= \left[i\gamma^{\mu} \sin(p_{\mu}) + m + \sum_{\mu} (1 - \cos(p_{\mu})) \right]^{-1} \\ &= \frac{-i\gamma^{\mu} \sin(p_{\mu}) + m + \sum_{\mu} (1 - \cos(p_{\mu}))}{\sum_{\mu} \sin^2(p_{\mu}) + \left[m + \sum_{\mu} (1 - \cos(p_{\mu})) \right]^2}. \end{aligned} \quad (3.23)$$

Now all of the doubler modes with any component of momentum $p_{\mu} \approx \pi$ become massive. The mass doubler modes gain is proportional to the inverse of the lattice spacing. The caveat in this strategy is that the Wilson term explicitly breaks chiral symmetry on the lattice even at zero bare fermion mass.

Unfortunately there is no easy way around this problem. The Nielsen and Ninomiya no-go theorem states that you cannot place chiral fermions on a regular lattice of even dimension

without fermion doublers [94, 95, 96, 82]. Any solution to the doubler problem will have to sacrifice chiral symmetry, dimensionality, or the regularity of the lattice. There are several types of fermionic actions that trade off these pitfalls. Some actions are able to preserve a near exact chiral symmetry, generally by adding dimensionality, however these actions are extremely expensive and difficult to simulate at strong coupling. The only action I will address further is the staggered action because it is used in our study.

3.3.3 Staggered Fermions

Staggered fermions were introduced by Kogut and Susskind [83]. The staggered action mixes Dirac and space time indices of the naive fermionic action to reduce the number of doublers from 16 to 4. To accomplish this lets define new field variables $\psi(x)'$ and $\bar{\psi}(x)'$

$$\begin{aligned}\psi(n) &= \gamma_1^{n_1} \gamma_2^{n_2} \gamma_3^{n_3} \gamma_4^{n_4} \psi(n)' \\ \bar{\psi}(n) &= \bar{\psi}(n)' \gamma_4^{n_4} \gamma_3^{n_3} \gamma_2^{n_2} \gamma_1^{n_1}.\end{aligned}\tag{3.24}$$

The transformation matrix is products of γ_μ raised to the power of the corresponding component n_μ of the site index $n_\mu = (n_1, n_2, n_3, n_4)$. We can now use these fields to build the staggered action

$$S(\psi', \bar{\psi}') = a^4 \sum_n \bar{\psi}(n)' \mathbb{I} \left(\sum_{\mu=1}^4 \eta_\mu(x) \frac{\psi(n + \hat{\mu})' - \psi(n - \hat{\mu})'}{2a} + m \psi(n)' \right), \tag{3.25}$$

where η_μ are the staggered sign functions

$$\begin{aligned}n_1(n) &= 1 \\ n_2(n) &= (-1)^{n_1} \\ n_3(n) &= (-1)^{n_1+n_2} \\ n_4(n) &= (-1)^{n_1+n_2+n_3}.\end{aligned}\tag{3.26}$$

This action has four copies of itself, one for each Dirac component. We finalize the action by taking one of these copies and coupling it to the gauge fields

$$S(\chi, \bar{\chi}) = a^4 \sum_n \bar{\chi}(n) \left(\sum_{\mu=1}^4 \eta_\mu(x) \frac{U_\mu(n) \chi(n + \hat{\mu}) - U_\mu^\dagger(n - \hat{\mu}) \chi(n - \hat{\mu})}{2a} + m \chi(n) \right). \tag{3.27}$$

The fields χ and $\bar{\chi}$ are Grassmann fields with only color indices and lack Dirac structure. Since we have kept one of four identical copies of our naive fermionic action we have essentially kept 4 of the 16 doubler modes. This action can be interpreted as having a 2^4 dimensional hyper cube unit cell with four flavors of fermions spread across it. The four species of fermions that remain are called referred to as tastes. While taste symmetry is exact in the continuum, it is broken by lattice artifacts. Taste breaking can be especially problematic at strong coupling where the gauge fields are coarser. Smearing terms can be added to the lattice action to mitigate or remove lattice artifacts associated with taste breaking. I will discuss the smearing improvement that we use in our simulations in the next section.

In QCD studies, taste poses a unique problem when trying to study 2, or 2+1, or 2+1+1 flavors of fermions. The 4 degenerate flavors have to be transformed into 1 or 2 flavors. This is accomplished by a technique called rooting [2]. Although rooted calculations routinely reproduce correct results, there is still some controversy about the technique [30, 31, 84]. Fortunately in our studies we examine theories with multiples of 4 flavors thereby circumventing the rooting problem entirely.

The balance that staggered fermions strike makes them a good choice in many situations. One benefit of staggered fermions is that they preserve a $U(1)_A$ chiral symmetry. Wilson fermions explicitly break chiral symmetry while lattice actions that preserve chiral symmetry exactly are much more computationally expensive. Finally, unlike Wilson fermions, there is a protected Goldstone mode which protects fermion masses from additive renormalization.

3.3.4 Improved Fermions

In the simulations detailed in chapters 4-6 we use nHYP improved staggered fermions in our action. nHYP smearing was introduced in [70] and reduces the problem of lattice artifacts, specifically taste breaking, in staggered actions [62].

nHYP smearing is built out of three consecutive smearing steps that are restricted to the

hypercubes associated with the link being smeared. The first step:

$$\bar{\Gamma}_{n,\mu;\nu,\rho} = (1 - \alpha_3)U_{n,\mu} + \frac{\alpha_3}{2} \sum_{\pm\eta \neq \rho,\nu,\mu} U_{n,\eta} U_{n+\hat{\eta},\mu} U_{n+\hat{\mu},\eta}^\dagger \quad (3.28)$$

$$\bar{V} = \bar{\Gamma}(\bar{\Gamma}^\dagger \bar{\Gamma})^{-\frac{1}{2}}, \quad (3.29)$$

where $\bar{V} = \Gamma(\Gamma^\dagger \Gamma)^{\frac{1}{2}}$ is a $U(N)$ projection. The second step uses the links from the first smearing step and projection and performs another smearing and projection

$$\tilde{\Gamma}_{n,\mu;\nu} = (1 - \alpha_2)U_{n,\mu} + \frac{\alpha_2}{4} \sum_{\pm\rho \neq \nu,\mu} \bar{V}_{n,\rho;\nu,\mu} \bar{V}_{n+\hat{\rho},\mu;\rho,\nu} \bar{V}_{n+\hat{\mu},\rho;\nu,\mu}^\dagger \quad (3.30)$$

$$\tilde{V} = \tilde{\Gamma}(\tilde{\Gamma}^\dagger \tilde{\Gamma})^{-\frac{1}{2}}. \quad (3.31)$$

Finally the last step is a third iteration of the smearing step

$$\Gamma_{n,\mu} = (1 - \alpha_1)U_{n,\mu} + \frac{\alpha_1}{6} \sum_{\nu \neq \mu} \tilde{V}_{n,\nu;\mu} \tilde{V}_{n+\hat{\nu},\mu;\nu} \tilde{V}_{n+\hat{\mu},\nu;\mu}^\dagger \quad (3.32)$$

$$V = \Gamma(\Gamma^\dagger \Gamma)^{-\frac{1}{2}}. \quad (3.33)$$

Here the links $U_{n,\mu}$ are un-smeared and the nHYP smeared links are $V_{n,\mu}$. The parameters $(\alpha_1, \alpha_2, \alpha_3)$ are parameters which can be tuned to control the amount of smearing. The indices after the semi-colon are not included in the staple sum. The link products

$$\sqcap_{n,\mu} = \sum_{\nu \neq \mu} U_{n,\nu} U_{n+\hat{\nu},\mu} U_{n+\hat{\mu},\nu}^\dagger \quad (3.34)$$

is the staple sum of the un-smeared link $U_{n,\mu}$.

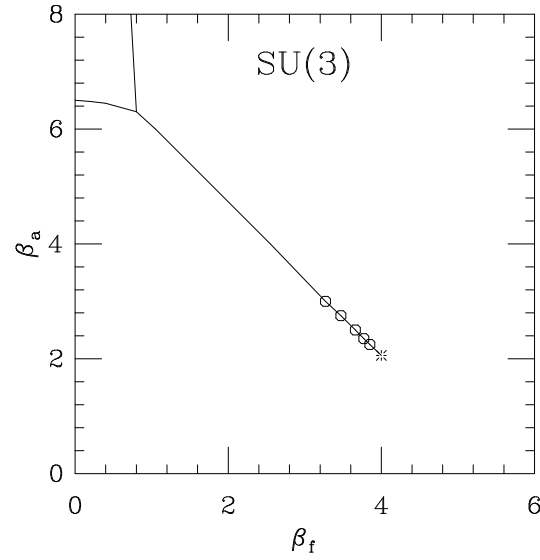


Figure 3.1: The phase diagram of SU(3) gauge theory in the fundamental-adjoint coupling plane. The solid lines are first order phase transitions. The phase transition ends in a second order critical point. Although the critical point lies above the β_F axis, the presence of the critical point influences the RG flow of actions where $\beta_A = 0$. Therefore we choose an action with $\beta_A < 0$ to avoid effects from the critical point. The phase diagram for confining SU(3) gauge theories with fermions is similar, only the location of the transitions changes.

Chapter 4

Monte Carlo Renormalization Group

4.1 Introduction

In recent years, many groups have initiated lattice investigations of strongly-coupled gauge-fermion systems beyond QCD. While the ultimate goal of these efforts is to explore potential new physics beyond the standard model, an essential step is to improve our theoretical understanding of the basic properties of these non-perturbative systems. In this chapter I study the renormalization group properties of SU(3) gauge theories with $N_f = 8$ and 12 nearly-massless fermions in the fundamental representation, through the Monte Carlo Renormalization Group (MCRG) two-lattice matching technique. This is one of several complementary analyses our group is involved with, two more of which (investigating Dirac eigenmode scaling and finite-temperature transitions) are discussed elsewhere [67, 107]. Recent references on SU(3) gauge theories with $N_f = 8$ and 12 include [48, 49, 6, 40, 86]; earlier works are reviewed in Ref. [56].

References [66, 65] study MCRG two-lattice matching for the 12-flavor system with nHYP-smearred staggered actions very similar to those we use here. Our gauge action includes both fundamental and adjoint plaquette terms, with coefficients related by $\beta_A = -0.25\beta_F$. The negative adjoint plaquette term lets us avoid a well-known spurious ultraviolet fixed point caused by lattice artifacts, and implies $\beta_F = 12/g^2$ at the perturbative level. In our fermion action, we use nHYP smearing with parameters (0.5, 0.5, 0.4), instead of the (0.75, 0.6, 0.3) used by Refs. [66, 65]. By changing the nHYP-smearing parameters in this way, we can access stronger couplings without encountering numerical problems. At such strong couplings, for both $N_f = 8$ and $N_f = 12$ we

observe a lattice phase in which the single-site shift symmetry (“ S^4 ”) of the staggered action is spontaneously broken (“ \mathcal{S}^4 ”) [28, 107].¹ In this work we only investigate couplings weak enough to avoid the \mathcal{S}^4 lattice phase.

In the next section, we review how the MCRG two-lattice matching technique determines the step-scaling function s_b in the bare parameter space. Although working entirely with bare parameters would be disadvantageous if our aim were to produce renormalized phenomenological predictions for comparison with experiment, our current explorations of the phase structures of the 8- and 12-flavor systems benefit from this fully non-perturbative RG approach, especially for relatively strong couplings. In Section 4.3 we present our results from the traditional MCRG two-lattice matching technique. While our 8-flavor s_b is significantly different from zero, for $N_f = 12$ we observe $s_b \lesssim 0$ for $\beta_F < 8$, indicating an infrared fixed point (IRFP).

We emphasize that while the existence of an IRFP is physical (scheme-independent), the coupling at which it is located depends on the choice of renormalization scheme. A limitation of traditional MCRG two-lattice matching is the need to optimize the RG blocking transformation separately for each lattice coupling β_F . As we explain below, this optimization forces us to probe a different renormalization scheme for each β_F , so that the bare step-scaling function we obtain is a composite of many different discrete β functions. Chapter 6 discusses an improvement to traditional MCRG that avoids this issue.

4.2 Method

In this section I will discuss the basic procedure for MCRG. MCRG is a real space renormalization group technique that can be used to study the critical properties of statistical systems. The idea was first proposed by Kadanoff. However it wasn’t until Wilson’s work that it became a well defined quantitative technique. The process of MCRG can be broken into two primary steps:

(1) Blocking

¹ Ref. [40] recently interpreted the \mathcal{S}^4 lattice phase in terms of relevant next-to-nearest neighbor interactions.

(2) Matching

Blocking integrates out the UV fluctuations, exposing the IR dynamics. Matching allows us to match pairs of couplings in such a way that we can measure the bare step scaling function. The steps scaling function is the discrete analog of the beta function in the continuum. In general it is possible to recover the beta function from the step scaling function. However, for our purposes this is not important. Our goal is to determine whether or not a particular theory is conformal in the IR. As long as strong enough couplings are studied, the IRFP will be visible in both the bare step scaling function and the renormalized beta function.

In the Wilsonian picture, couplings are characterized by their engineering dimension. In a four dimensional theory, a coupling is relevant if its engineering dimension is less than four. The coupling is marginal if its engineering dimension is equal to four and irrelevant if the engineering dimension is greater than four. The RG flow of the system is governed by fixed points and the renormalized trajectory. The systems we study have a repulsive UV fixed point. If the theory is conformal it will also have an attractive IR fixed point. The renormalized trajectory originates at the UV fixed point and flows in relevant directions towards the infrared. Flow in irrelevant directions will always be directed towards the renormalized trajectory. Finally flow along the renormalized trajectory will always flow into attractive fixed points. In theories with multiple fixed points, the renormalized trajectory connects the fixed points in coupling space. Flow from an arbitrary point in coupling space will always be towards the renormalized trajectory in irrelevant or marginal directions and along the renormalized trajectory in relevant directions. Generally the flow should be fastest in irrelevant directions.

4.2.1 Blocking

The first step in MCRG is to block the lattice. Blocking requires that the theory is defined on a regular lattice that has a discrete scaling symmetry. Discrete scaling symmetry, shown in figure 4.1 simply means that you can divide the sites of the lattice into regular blocks that tile the entire lattice and then replace those blocks with a single site. The new site can be located at a site from

the original lattice or anywhere in the block. The value at this site is typically computed as some function of the sites in the blocking group. The lattice has a discrete scaling symmetry if [18]

- (1) we perform this block replacement the same way to every block in the original lattice
- (2) we scale the lattice spacing of the lattice formed by the block variables: $a \rightarrow a' \equiv sa$
- (3) the lattice of block replaced observables is the same size as the original lattice

Here s is a scale factor which relates the original lattice to the blocked lattice.

The blocked lattice will have fewer sites than the original lattice. If the block variable has P sites from the original lattice and d dimensions then the relationship between the number of sites lost, the scale factor and the lattice dimension is $p = s^d$. Naturally the blocking process can be repeated in a recursive manner until blocking is no longer possible. As an example consider a 32^2 square lattice that is being blocked by a factor $s = 2$. Such a lattice can be blocked 4 times: $32^2 \rightarrow 16^2 \rightarrow 8^2 \rightarrow 4^2 \rightarrow 2^2$. Performing another blocking would result in a single point which doesn't have the same dimensionality as our initial $2 - D$ lattice. It is clear that some starting volumes are more flexible than others. A 14^2 lattice can only be blocked once, $14^2 \rightarrow 7^2$, since a 7^2 lattice can not be blocked with a scale factor of 2.

In a lattice simulation the quantities we are interested in are the link variables. The link variable $U_{n,\mu}$ encodes the gauge field at site n in all four $\hat{\mu}$ directions. A MCRG blocking step can be defined as a transformation that takes two links $U_{n,\mu}$ and $U_{n+\hat{\mu},\mu}$ and creates one link with twice the lattice spacing of the initial two links.

Employing a square blocking scheme as described above means the blocked lattice will have a factor of 2^N fewer links than the original lattice. In four dimensions this means that an observable such as a Wilson loop will have a factor of 16 less statistics on the blocked lattice compared to the original lattice. Fortunately, on an N dimensional hyper cubic lattice there are 2^N unique ways to perform this blocking. Therefore in four dimensions there are 16 unique blocking configurations. We can preserve statistics in our calculation by performing all of the unique blocking transformations and averaging all of the observables from those blocking transformations together. Additionally we

store the blocked links on a lattice of the original cardinality using offsets as shown in figure 4.2. This allows us to efficiently store a lattice blocked down to any allowable level and has the added benefit that only minor adjustments need to be made to our code to calculate blocked observables.

Lets consider how blocking will effect a lattice model with d dimensions and action $S(K_i)$. Here K_i are the set of all possible couplings in the action. In lattice simulations only a small number of these couplings are actually tuned. The system is characterized by one or more length scales i.e. the correlation length ξ . After blocking, the links have a lattice spacing scaled by s , $a' = sa$ such that the physical size of the box has not been changed. This removes the UV fluctuations below the length scale sa . Because the UV modes have been integrated out the blocked lattice is described by a new action S' with a new set of couplings $K_i^{n_b}$ where n_b is the number of blocking steps. As long as s is smaller than the lattice correlation length the infrared properties of the system are unaffected. Successive blocking steps will define a flow in coupling space:

$$K_i^{(0)} \rightarrow K_i^{(1)} \rightarrow K_i^{(2)} \rightarrow \dots \rightarrow K_i^{(n_b)}, \quad (4.1)$$

where K_i^0 denotes the couplings on the unblocked lattice.

Although the physical correlation length is unchanged the lattice correlation length after n_b blocking steps is scaled by

$$\xi^{n_b} = s^{-n_b} \xi^{(0)}. \quad (4.2)$$

Critical fixed points exist when $\xi = \infty$ and trivial fixed points exist when $\xi = 0$. The scaling behavior at critical fixed points are well understood theoretically. Here the linearized RG transformation predicts the scaling operators and the corresponding scaling dimension.

Figure 4.3 visualizes the renormalization group flow that I have described for a system with one fixed point and only one relevant coupling at the fixed point, K_0 . All other couplings are irrelevant and are collapsed onto the y-axis. The critical surface is shown in the figure is depicted at $K_0 = 0$. A simulation point in parameter space, P , is chosen. The flow lines approach the fixed point in the irrelevant directions and flow away from the fixed point in the relevant direction. After some number of RG steps the irrelevant operators have died out and the system is flowing

along the renormalized trajectory. The renormalized trajectory is uniquely determined by the RG transformation.

Finally, in our blocking step we also apply a nHYP smearing step. By changing the parameters in the smearing step we can control the renormalized trajectory that we will flow to under successive blocking steps. If we start with an infinite volume the choice of nHYP parameters would be of little consequence because we could block our system an infinite number of times until we reach the renormalized trajectory. Unfortunately in the real world we are not afforded that luxury. It is necessary to choose a smearing that guarantees you can reach the renormalized trajectory in a finite number of steps. We accomplish this by applying a nHYP smearing to the unblocked lattice. We choose the three parameters of the nHYP smearing to be $(\alpha, 0.2, 0.2)$ where we use several values of α , $0.3 \leq \alpha \leq 0.7$. After smearing the unblocked lattice, we multiply adjacent links to form a single blocked link.

4.2.2 Two-lattice matching procedures and the need for optimization

The goal of two lattice matching is to determine a series of lattice couplings, $\beta_0, \beta_1, \dots, \beta_n, \dots$, with lattice spacings that differ by a factor of s between consecutive points. That is $a(\beta_n) = a(\beta_{n-1})/s$. The scale change s is the same scale change as in the blocking discussion above. I denote the bare step scaling function as

$$s_b(\beta_n, s) = \beta_n - \beta_{n-1}. \quad (4.3)$$

The renormalized running coupling of theories governed by a Gaussian fixed point can be recovered with a bit of extra work. In weak coupling near the fixed point a renormalized coupling can be calculated. This coupling can be compared to a continuum regularization scheme. At stronger coupling a physical quantity or other quantity such as the Sommer parameter or the Wilson Flow is used to determine the lattice scale. Since the primary focus of this thesis is to determine the existence or lack thereof of an infrared fixed point in strong coupling I don't attempt to follow such a procedure. In the future finding the corresponding renormalized beta function may

be of interest.

Figure 4.4 demonstrates how two lattice matching is performed. In the figure we start with two points in coupling space P_1 and P_2 . Under successive RG transformations they flow along the lines indicated. If we can identify two sets of couplings that flow to the same point on the renormalized trajectory after the same number of blocking steps we know that the correlation lengths are identical. From equation 4.2 we know that if two sets of couplings flow to the same point on the renormalized trajectory, but one set of couplings required one less blocking step than the lattice correlation length differs from a factor of s . Thus we have accomplished our goal. We have identified two points in coupling space P_1 and P_2 with corresponding $\xi_2 = \xi_1/s$ after n and $(n - 1)$ blocking steps.

Demonstrating that the couplings are equal is equivalent to showing that the actions are identical, $S(P_1^{(n_b)}) = S(P_2^{(n_b-1)})$. Calculating the blocked action is extremely challenging. Fortunately we don't have to do this. We are saved because we don't need to know the exact form of the actions to demonstrate that they are equivalent. Instead we can show that expectation values of every operator measured on $P_1^{(n_b)}$ and $P_2^{(n_b-1)}$ are identical. Generating configuration ensembles of blocked lattices with the approximate Boltzman weight is trivial, we simply need to take an existing ensemble of configurations and then block them [113].

Putting everything together we can prescribe the following procedure for two lattice matching [63]:

- (1) Generate a configuration ensemble of size L^d with action $S(P_1)$. Block each configuration n_b times and measure a set of expectation values on the resulting $(L/s^{n_b})^d$ set.
- (2) Generate several configurations of size $(L/s)^d$ with action $S(P_2)$, where each P_2 is a set of trial couplings. Block each configuration $(n_b - 1)$ times and measure the same expectation values on the resulting $(L/s^{n_b})^d$ set. Compare the results with that obtained in step 1 and tune the coupling P_2 such that the expectation values agree. If they do we have determined a flow $P_1 \rightarrow P_2$ over scale s .

This method is powerful for several reasons. Since we compare measurements on the same lattice size, the finite volume corrections are small. Working on lattices larger than the correlation length of the system is not required. MCRG works in both the confined and deconfined phases. Since the matching is accomplished from measurements of local operators, their computation is cheap and the statistical errors are small. Finally because the flow lines are unique it is sufficient to match only one operator. The other operators should give the same prediction. If we fail to reach the renormalized trajectory it will be evident in that the other operators do not match. Therefore we can use the spread in the matching values of the different operators as a measure of systematic error in our matching.

The discussion so far has focused on fixed points with one relevant operator. If the fixed point has two relevant operators the matching process is identical, only now it is necessary to tune two operators to find pairs of points in a two dimensional coupling space. While this is possible, it is much easier to fix one of the relevant couplings to its fixed point value. By effectively removing one of the relevant couplings we can proceed as with the matching of the remaining coupling.

There is one final hurdle to overcome. On a finite volume we only have a finite number of blocking steps. If we are unlucky we might choose a block transformation, and therefore a renormalized trajectory, that is so far away from the bare parameters in couplings space that we never approach it. Accordingly after we have performed the maximum number of blocking steps the system may not reach the renormalized trajectory. This is a problem because matching requires both blocked ensembles to have reached the renormalized trajectory.

One solution is to use a bigger volume. Using a bigger volume means that we can perform more blocking steps, eventually we will arrive at the renormalized trajectory. This solution is clearly flawed. Lets assume that the largest volume you have is 32^4 and you can not reach the renormalized trajectory in four blocking steps. To achieve one more blocking step you need to jump to 64^4 lattices, if that is not enough then on to 128^4 lattices. We are now considering lattices that are larger than those used for large professional calculations! The computational cost of configuration generation scales roughly as $V^{5/4}$, in our doubling scenario generating larger volumes

will require at least $2^{5/4}$ more computational power.

Clearly we need another solution. As I mentioned above we apply a nHYP smearing step during our block transformation. This serves a crucial role, nHYP smearing has 3 parameters. Since the different block transformations correspond to different renormalized trajectories, tuning the nHYP parameters allows us to control the renormalized trajectory that we flow to. We use the first parameter of the nHYP smearing as an optimization parameter and fix the second and third. We identify the optimal blocking accordingly:

- (1) Consistent matching between the different operators: along the RT all expectation values should agree on the matched configuration sets. Any deviation is a measure that the RT has not been reached.
- (2) Consecutive blocking steps should give the same matching coupling.

4.2.3 Chirally Broken Theories

In a confining and chiral broken theory, such as QCD, the RG flow will away from the repulsive UV fixed point into the IR. If the fermions are massless the only relevant operator is the gauge coupling β . This is the β that we set in the action of our lattice simulation. Therefore the renormalized trajectory will flow along the β axis in coupling space from infinity to zero. RG flows will flow to the renormalized trajectory and then along it as shown in figure 4.5.

4.2.4 Conformal Theories

In a conformal system with massless fermions, the RG flow from points on the weak coupling side of the IRFP will also flow towards the renormalized trajectory in irrelevant directions. However unlike the QCD case the flow along the renormalized trajectory will terminate at the IRFP. At the IRFP β also becomes an irrelevant operator. Had we started from the strong coupling side of the IRFP we would observe something completely new. The flow would move towards the renormalized trajectory in the irrelevant directions but it would flow into the IRFP. The flow into the IRFP from

the strong coupling is ‘backwards’ from the flow in a chiral broken theory. The behavior of a conformal theory is shown in figure 4.6.

4.3 8 and 12 Flavor Results

In this section I employ the two lattice matching technique described earlier in this chapter to study SU(3) gauge theories with 8 and 12 chiral fermions in the fundamental representation. We proceed by repeatedly applying an nHYP RG blocking transformations with scale factor $s = 2$. The lattices used in this study have volumes $24^3 \times 48$, $12^3 \times 24$ and $6^3 \times 12$. On the largest $24^3 \times 48$ lattices that we use in this current study, we work with fermion masses $m = 0.0025$ to stay near the $m = 0$ critical surface. Under RG blocking with scale factor s , the fermion mass changes as $s^{1+\gamma_m}$ where γ_m is the mass anomalous dimension. Therefore we use $m = 0.01$ on $12^3 \times 24$ and $m = 0.02$ on $6^3 \times 12$ lattices. We have explicitly checked that these masses are small enough to introduce only negligible finite-mass effects, by generating lattices with $m = 0$ for some points and obtaining indistinguishable results.

Under RG blocking on the $m = 0$ critical surface, the system flows toward the renormalized trajectory in irrelevant directions, and along it in the relevant direction. By blocking the larger lattices (with β_F) n_b times and the smaller lattices (with β'_F) only $n_b - 1$ times, we obtain blocked systems with the same lattice volume. If these blocked systems have both flowed to the same point on the renormalized trajectory, then we can conclude that $\xi(\beta_F) = 2\xi(\beta'_F)$ on the unblocked systems, as desired.

We determine whether the blocked systems have flowed to the same point on the renormalized trajectory by matching several short-range gauge observables: the plaquette, all three six-link loops, and two planar eight-link loops. For a given β_F , each observable may predict a different $\Delta\beta_F \equiv \beta_F - \beta'_F$. The spread in these results is a systematic error that dominates our uncertainties. In principal more observables can be added, our choices are limited by the final volume of the blocked lattice.

In an IR-conformal system, the gauge coupling that is relevant at the perturbative gaussian

FP becomes irrelevant at the IRFP. The renormalized trajectory connects these two fixed points. When RG flows approach this renormalized trajectory, the two-lattice matching can be performed and interpreted the same way as in confining systems. In this region the gauge coupling flows to stronger couplings, $\Delta\beta_F > 0$ corresponding to a negative RG β function. The situation is less clear at stronger couplings where we might naïvely expect backward flow. If there is no ultraviolet FP in this region to drive the RG flow along a renormalized trajectory, the two-lattice matching might become meaningless. This issue affects every method that attempts to determine the flow of the gauge coupling in IR-conformal systems at strong coupling. In all published studies that report an IRFP, backward flow has only been observed in a very limited range of couplings in the immediate vicinity of the IRFP (cf. Ref. [56]).

Since we can block our lattices only a few times, we must optimize the two-lattice matching by requiring that consecutive RG blocking steps yield the same $\Delta\beta_F$. We identify the optimized $\Delta\beta_F$ with the bare step-scaling function s_b . The traditional technique reported in this chapter optimizes the RG blocking transformation (renormalization scheme). The new method we propose in Section 6 instead applies the Wilson flow to the lattice system prior to RG blocking, and optimizes the flow time t_f .

As in Refs. [66, 65], we use RG blocking transformations that include two sequential HYP smearings with parameters $(\alpha, 0.2, 0.2)$, and optimize α as shown in the left panel of Fig. 4.7. Qualitatively, this optimization finds the renormalization scheme for which the renormalized trajectory passes as close as possible to the lattice system with coupling β_F . Without optimization, residual flows in irrelevant directions can distort the results: this is the reason $\Delta\beta_F$ changes with α in Fig. 4.7, and also explains why increasing the number of blocking steps reduces this α -dependence.

We note that finite-volume effects are minimized by carrying out the optimization on blocked lattices of the same volume, which was reported by Ref. [66]. That is, we should compare $\Delta\beta_F$ from matching $12^3 \times 24$ blocked to $3^3 \times 6$ vs. $6^3 \times 12$ blocked to $3^3 \times 6$ with that from matching $24^3 \times 48$ blocked to $3^3 \times 6$ vs. $12^3 \times 24$ blocked to $3^3 \times 6$. As in Refs. [66, 65], we do not explore weak enough couplings to recover the two-loop perturbative predictions $s_b \approx 0.3$ for $N_f = 12$.

4.3.1 8 Flavors

In our study of SU(3) gauge theories with 8 flavors we calculated the step scaling function at four couplings corresponding to $\beta_F = 5.4, 5.6, 6, 7$. We were unable to perform matching at stronger coupling due to the appearance of the \mathcal{S}^4 phase at $\beta_f = 4.6$. At weaker couplings our results for the step scaling function are lower than those predicted by two loop perturbation theory. Perturbation theory predicts a value of $s_b \approx 0.6$. This implies that our simulations are at couplings too strong to perform comparison to perturbation theory. Figure 4.8 displays our final results. The error bars are indicative of the spread the step scaling function predicted by different observables.

It is clear that the step scaling function is far removed from zero. This indicates that the 8 flavor theory does not have an infrared fixed point. While the calculated step scaling function is not in explicit agreement with perturbative value it is somewhat close. The fact our value for the step scaling function is below the perturbative value indicates that we are likely in strong enough coupling that perturbation theory is not reliable. Furthermore interpolating between the our measured s_b and the perturbative value does not require a fixed point to develop. This picture is consistent with 8 flavors being chirally broken and confining over the range of couplings simulated.

4.3.2 12 Flavors

Our investigation of SU(3) gauge theory with 12 flavors reveals behavior very different from the 8 flavor results. We calculated the step scaling function at $\beta_F = 5, 6, 7, 8, 9, 10$. As with the 8 flavor case we were unable to perform matching at stronger couplings because of the \mathcal{S}^4 phase at $\beta = 2.7$. At the weakest couplings studied, we find that our results for the step scaling function are smaller than those predicted by the two loop perturbative calculation which predicts $s_b \approx 0.3$. As with the 8 flavor results, this implies that we are at strong enough coupling that perturbation theory is unreliable.

The results shown in 4.9 show that the step scaling function for $\beta_F = 8, 9, 10$ is positive while the step scaling function at $\beta_F = 5, 6, 7$ is negative or close to zero. This is consistent

with backwards flow at strong coupling and an infrared fixed point at $\beta_F < 8$. Unfortunately we are unable to cleanly identify where the fixed point is. I discuss this point in the next section. Independent of the location of the fixed point, our results are not consistent with a chirally broken theory.

4.3.3 The Case of the Wandering Fixed Point

Recall that our optimization of the RG blocking transformation means that we use a different renormalization scheme for each coupling β_F , so these bare step scaling functions are composites of several different discrete β functions. In an IR conformal system, each renormalization scheme has a fixed point because the existence of the fixed point is scheme independent. The location of the fixed point however is a scheme dependant property. Figure 4.10 demonstrates how this can happen and complicate locating the IRFP. Therefore it is reasonable to expect the IRFP to be located at a different point in coupling space for each β_F that we calculate s_b for. For example, with $N_f = 12$ at $5 \leq \beta_F \leq 6$, our optimization selects renormalization schemes with the fixed point near β_F , so that s_b is roughly consistent with zero over an extended range.

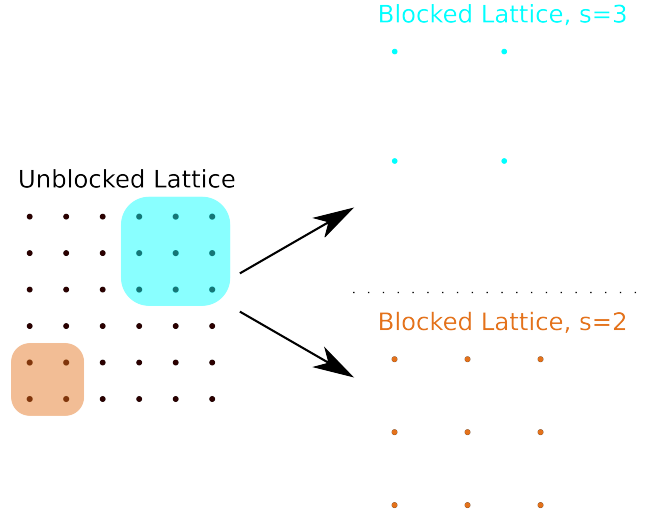


Figure 4.1: The original 6×6 lattice on the left possesses a discrete scaling symmetry of $s = 2$ and $s = 3$. The shaded orange square is a $s = 2$ block variable. The resulting orange 3×3 blocked lattice in the bottom right formed by replacing each block variable with a single site in the upper left corner of the block. The cyan shaded region shows a $b = 3$ block variable. Performing a block transformation that replaces each block with a point in the upper left of the block produces the 2×2 blocked lattice shown in the upper right.

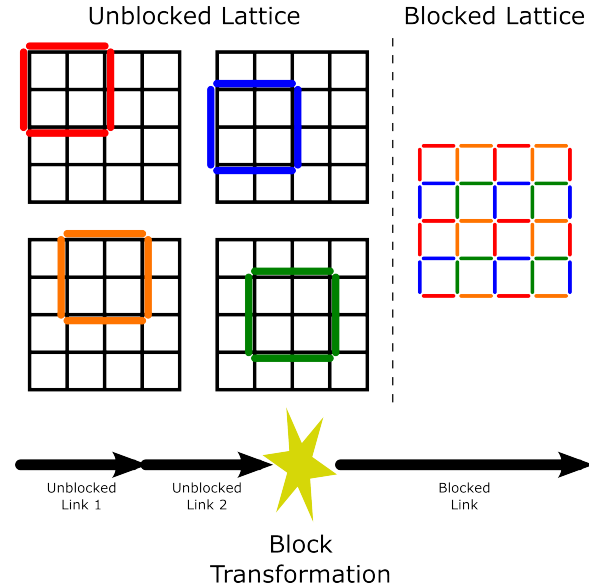


Figure 4.2: This figure shows how links are blocked on the lattice. Two adjacent links in the same direction are block transformed to form one link of twice the lattice spacing. We perform all possible block transformation shown on the left as the red, blue, green, and orange block tilings of the unblocked lattice. We then store the links of the blocked lattice as shown on the right hand side of the figure.

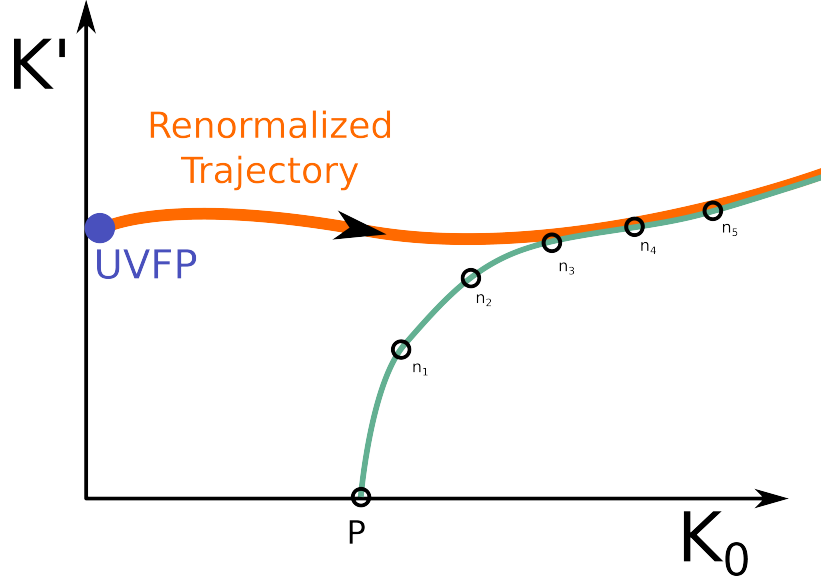


Figure 4.3: Here I show coupling space for a system with one relevant direction K_0 . All irrelevant directions are collected in K' . We simulate at a point P in parameter space. As we block the system the effective couplings will change. In the diagram here the couplings reach the renormalized trajectory after 3 blocking steps. Further blocking steps move the couplings along the renormalized trajectory.

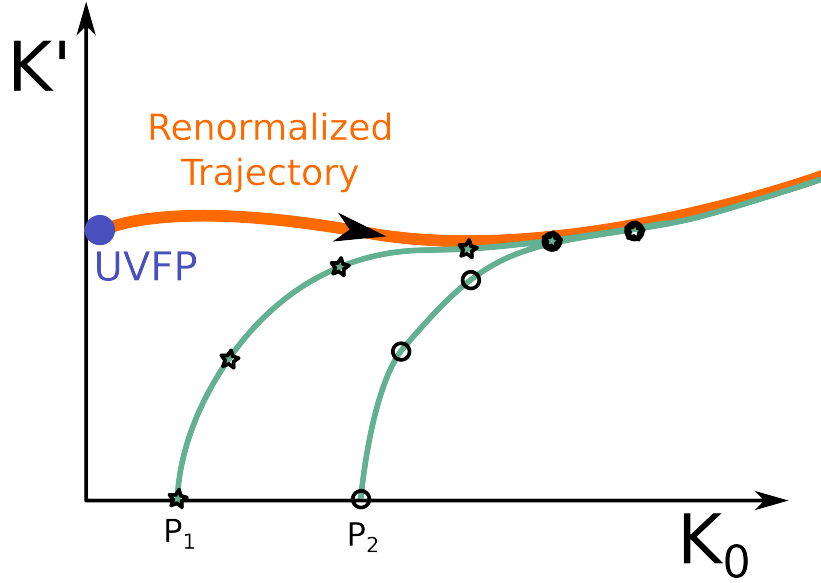


Figure 4.4: For matching we pick two points in coupling space P_1 and P_2 . After 3 blocking steps, shown as circles the effective action of P_2 has reached the renormalized trajectory. P_1 requires 4 blocking steps shown as stars but reaches the same point on the renormalized trajectory. Because P_2 took one less blocking step its correlation length ξ is a factor of s smaller than the correlation length of the first ensemble that started at P_1 . By choosing pairs of points in coupling space that block to the same point on the renormalized trajectory in coupling space, we can construct a discrete step scaling function.

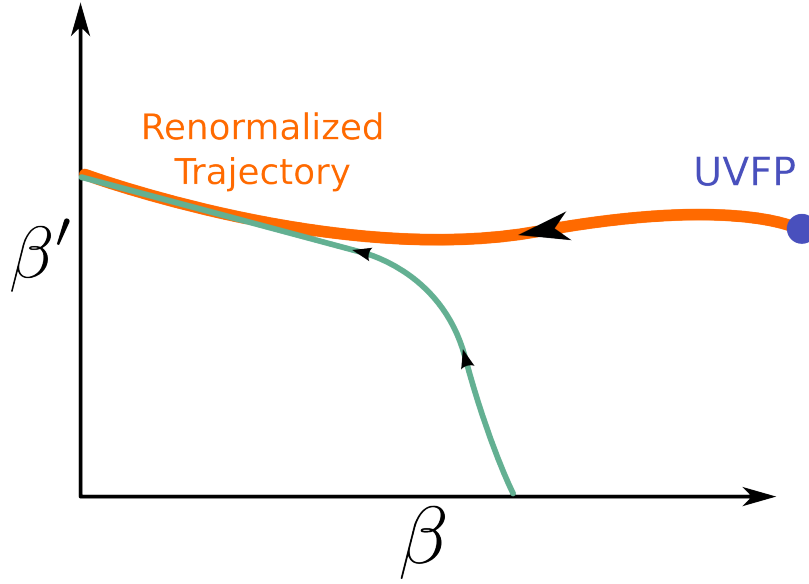


Figure 4.5: The RG flow of a confining theory on the $m = 0$ critical surface. β is the relevant gauge coupling, β' are irrelevant couplings.

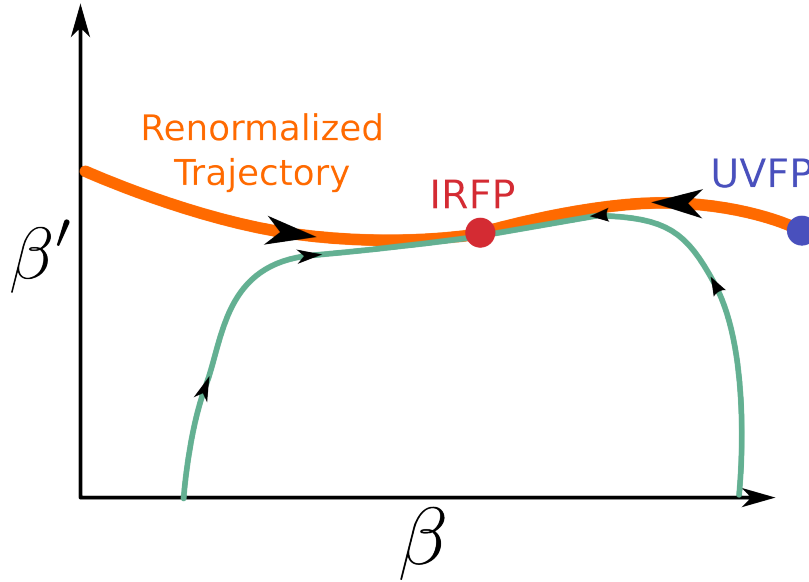


Figure 4.6: The RG flow of a conformal theory on the $m = 0$ critical surface. β is the relevant gauge coupling, β' are irrelevant couplings.

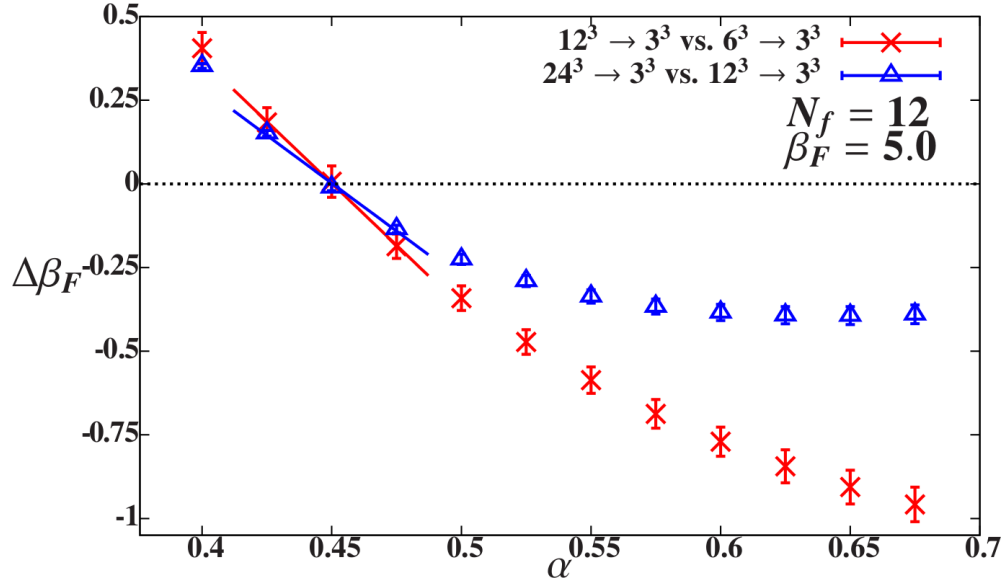


Figure 4.7: Optimization of the HYP-smearing parameter α in the RG blocking transformation, for $\beta_F = 5.0$. The uncertainties on the data points are dominated by averaging over the different observables as described in the text.

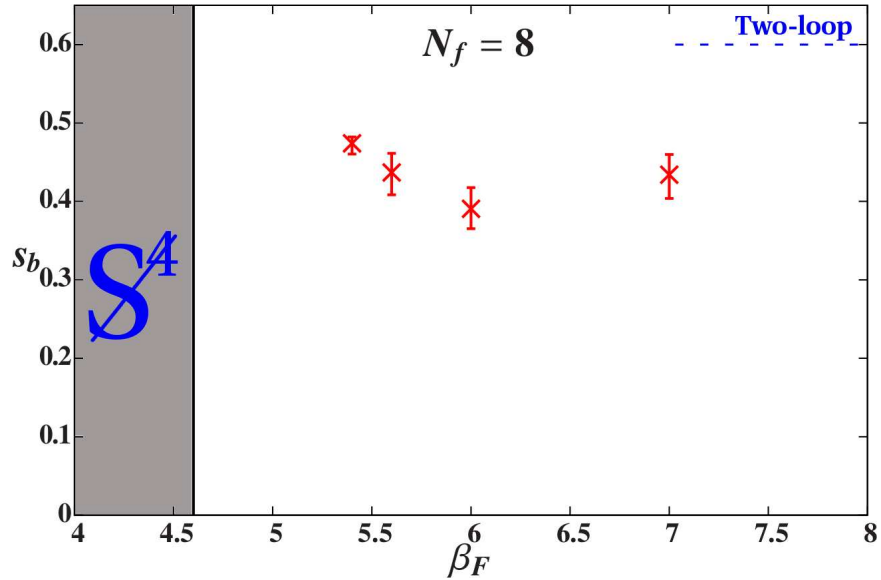


Figure 4.8: Results for the bare step-scaling function s_b from traditional MCRG two-lattice matching with $24^3 \times 48$, $12^3 \times 24$ and $6^3 \times 12$ lattice volumes for $N_f = 8$. The blue dashed lines are perturbative predictions for asymptotically weak coupling.

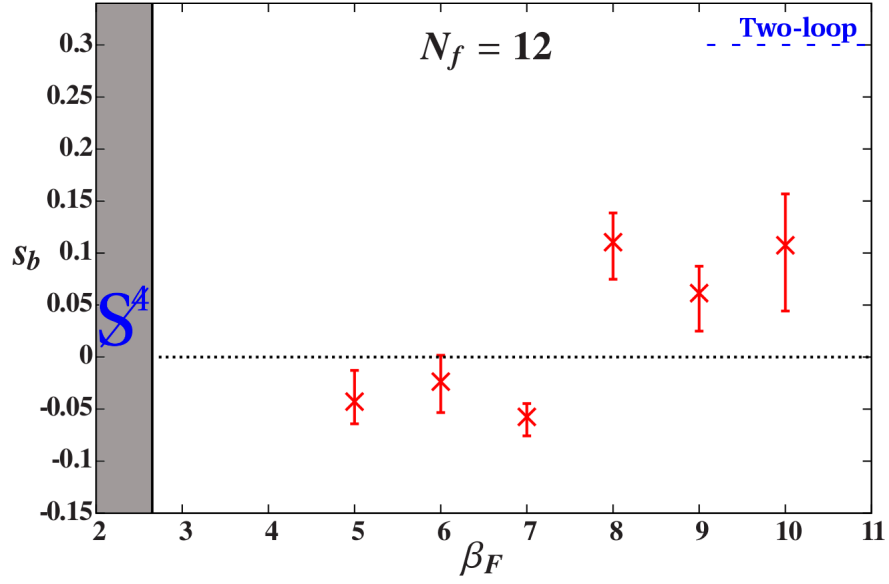


Figure 4.9: Results for the bare step-scaling function s_b from traditional MCRG two-lattice matching with $24^3 \times 48$, $12^3 \times 24$ and $6^3 \times 12$ lattice volumes for $N_f = 12$. The blue dashed lines are perturbative predictions for asymptotically weak coupling.

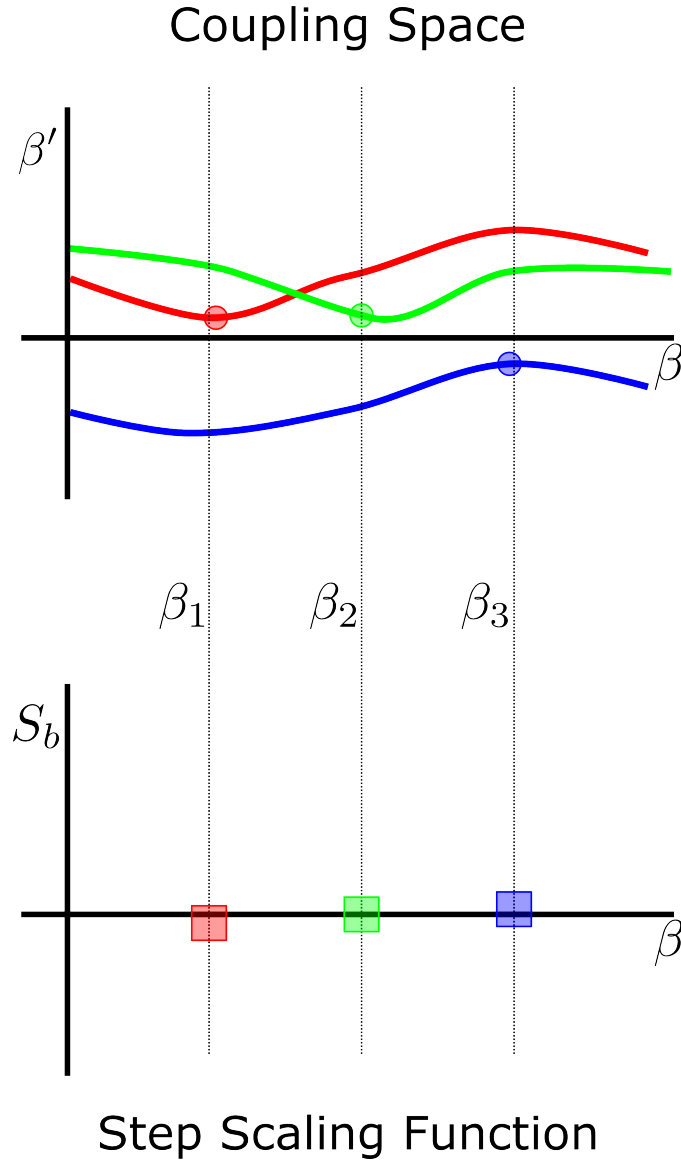


Figure 4.10: An illustration of how optimizing the block transformation can result in difficulties locating an IRFP, β is the relevant gauge coupling and β' are irrelevant couplings. The upper figure shows the renormalized trajectory in red, green, and blue found by optimizing the RG transformation at β_1, β_2 , and β_3 respectively. The location of the IRFP changes in each renormalized trajectory, in this picture the IRFP is moved to the coupling we perform MCRG at. The resulting s_b is consistent with zero across a wide range of couplings.

Chapter 5

Wilson Flow

5.1 Introduction

Asymptotically-free $SU(N)$ gauge theories coupled to N_f massless fundamental fermions are conformal in the infrared if N_f is sufficiently large, $N_f \geq N_f^{(c)}$. Their renormalization group (RG) β functions possess a non-trivial infrared fixed point (IRFP) where the gauge coupling is an irrelevant operator. Although this IRFP can be studied perturbatively for large N_f near the value at which asymptotic freedom is lost [22, 14], as N_f decreases the fixed point becomes strongly coupled. Systems around $N_f \approx N_f^{(c)}$ are particularly interesting strongly-coupled quantum field theories, with non-perturbative conformal or near-conformal dynamics. Their most exciting phenomenological application is the possibility of a light composite Higgs boson from dynamical electroweak symmetry breaking [47, 91, 8, 52, 3]. Due to the strongly-coupled nature of these systems, lattice gauge theory calculations are a crucial non-perturbative tool with which to investigate them from first principles. Many lattice studies of potentially IR-conformal theories have been carried out in recent years (cf. the recent reviews [93, 57] and references therein). While direct analysis of the RG β function may appear an obvious way to determine whether or not a given system flows to a conformal fixed point in the infrared, in practice this is a difficult question to address with lattice techniques. In particular, extrapolation to the infinite-volume continuum limit is an essential part of such calculations.

In the case of $SU(3)$ gauge theory with $N_f = 12$ fundamental fermions, several lattice groups have investigated the step scaling function, the discretized form of the β function. To date, these

studies either did not reach a definite conclusion [64, 86] or may be criticized for not properly taking the infinite-volume continuum limit [9, 10, 64, 66, 76, 100]. At the same time, complementary numerical investigations have been carried out, considering for example the spectrum, or bulk and finite-temperature phase transitions [39, 46, 7, 33, 28, 27, 48, 49, 6, 4, 80]. The different groups performing these studies have not yet reached consensus regarding the infrared behavior of the 12-flavor system.

Our own $N_f = 12$ results favor the existence of a conformal IRFP, which we observe in Monte Carlo RG studies [66, 100]. Our zero- and finite-temperature studies of the lattice phase diagram show a bulk transition consistent with conformal dynamics [107, 69]. From the Dirac eigenvalue spectrum [27], and from finite-size scaling of mesonic observables [24], we obtain consistent predictions for a relatively small fermion mass anomalous dimension: $\gamma_m^* = 0.32(3)$ and $0.235(15)$, respectively. While this conclusion, if correct, would render the 12-flavor system unsuitable for composite Higgs phenomenology, we consider $N_f = 12$ to remain an important case to study. Considerable time and effort has already been invested to obtain high-quality lattice data for the 12-flavor system. Until different methods of analyzing and interpreting these data can be reconciled – or the causes of any remaining disagreements can be clarified – it will not be clear which approaches are most reliable and most efficient to use in other contexts.

The recent development of new running coupling schemes based on the gradient flow [88, 87, 51, 50, 53] provides a promising opportunity to make progress. In this work we investigate step scaling using the gradient flow running coupling.¹ We begin by introducing a non-perturbative improvement to this technique, which increases our control over the continuum extrapolation by reducing the leading-order cut-off effects. While this improvement is phenomenological in the sense that we have not derived it systematically through a full improvement program, it is generally applicable to any lattice gauge theory of interest and in our tests removes all $\mathcal{O}(a^2)$ cut-off effects. We illustrate it first for 4-flavor $SU(3)$ gauge theory, a system where the running coupling has

¹ We are aware of two other ongoing investigations of the $N_f = 12$ gradient flow step scaling function, by the authors of Ref. [51] and Ref. [86].

previously been studied with both Wilson [115] and staggered [97, 51, 50] fermions. We then turn to $N_f = 12$, where we show that the infinite-volume continuum limit is well defined and predicts an IRFP. In both the 4- and 12-flavor systems, our improvement can remove all observable $\mathcal{O}(a^2)$ effects, despite the dramatically different IR dynamics. We conclude with some comments on other systems where improved gradient flow step scaling may profitably be applied.

5.2 Improving the Gradient Flow Step Scaling

The gradient flow is a continuous invertible smearing transformation that systematically removes short-distance lattice cut-off effects [88, 87]. At flow time $t = a^2 t_{\text{lat}}$ it can be used to define a renormalized coupling at scale $\mu = 1/\sqrt{8t}$

$$g_{\text{GF}}^2(\mu = 1/\sqrt{8t}) = \frac{1}{\mathcal{N}} \langle t^2 E(t) \rangle, \quad (5.1)$$

where “ a ” is the lattice spacing, t_{lat} is dimensionless, and the energy density $E(t) = -\frac{1}{2} \text{ReTr} [G_{\mu\nu}(t) G^{\mu\nu}(t)]$ is calculated at flow time t with an appropriate lattice operator. We evolve the gradient flow with the Wilson plaquette term and use the usual “clover” or “symmetric” definition of $G_{\mu\nu}(t)$. The normalization \mathcal{N} is set such that $g_{\text{GF}}^2(\mu)$ agrees with the continuum $\overline{\text{MS}}$ coupling at tree level.

If the flow time is fixed relative to the lattice size, $\sqrt{8t} = cL$ with c constant, the scale of the corresponding coupling $g_c^2(L)$ is set by the lattice size. Like the well-known Schrödinger functional (SF) coupling, $g_c^2(L)$ can be used to compute a step scaling function [51, 50, 53]. The greater flexibility of the gradient flow running coupling is a significant advantage over the more traditional SF coupling. A single measurement of the gradient flow will provide g_c^2 for a range of c . In our study we obtain g_c^2 for all $0 \leq c \leq 0.5$ separated by $\delta t_{\text{lat}} = 0.01$. Each choice of c corresponds to a different renormalization scheme, which can be explored simultaneously on the same set of configurations [53].

The normalization factor \mathcal{N} in finite volume has been calculated for anti-periodic boundary conditions (BCs) in refs. [51, 50], and for SF BCs in Ref. [53]. In this work we use anti-periodic

BCs, for which

$$\frac{1}{\mathcal{N}} = \frac{128\pi^2}{3(N^2 - 1)(1 + \delta(c))} \quad \delta(c) = \vartheta^4 \left(e^{-1/c^2} \right) - 1 - \frac{c^4 \pi^2}{3}, \quad (5.2)$$

where $\vartheta(x) = \sum_{n=-\infty}^{\infty} x^{n^2}$ is the Jacobi elliptic function. For $0 \leq c \leq 0.3$ the finite-volume correction $\delta(c)$ computed in Ref. [51] is small, $|\delta(c)| \leq 0.03$. As explained in refs. [51, 50], the RG β function of g_{GF}^2 is two-loop universal with SF BCs, but only one-loop universal with anti-periodic BCs.

At non-zero lattice spacing g_{GF}^2 has cut-off corrections. These corrections could be $\mathcal{O}(a)$ for unimproved actions, and even $\mathcal{O}(a)$ -improved actions could have large $\mathcal{O}(a^2[\log a]^n)$ -type corrections [12, 13]. In existing numerical studies of staggered or $\mathcal{O}(a)$ -improved Wilson fermions the leading lattice corrections appear to be $\mathcal{O}(a^2)$ [53, 111],

$$g_{\text{GF}}^2(\mu; a) = g_{\text{GF}}^2(\mu; a = 0) + a^2 \mathcal{C} + \mathcal{O}(a^4[\log a]^n, a^4). \quad (5.3)$$

It is possible to remove, or at least greatly reduce, the $\mathcal{O}(a^2)$ corrections in eq. 5.3 by defining

$$\tilde{g}_{\text{GF}}^2(\mu; a) = \frac{1}{\mathcal{N}} \langle t^2 E(t + \tau_0 a^2) \rangle, \quad (5.4)$$

where $\tau_0 \ll t/a^2$ is a small shift in the flow time. In the continuum limit $\tau_0 a^2 \rightarrow 0$ and $\tilde{g}_{\text{GF}}^2(\mu) = g_{\text{GF}}^2(\mu)$.

There are several possible interpretations of the t -shift in eq. 5.4. The gradient flow is an invertible smearing transformation, so one can consider τ_0 as an initial flow that does not change the IR properties of the system but leads to a new action. The gradient flow coupling \tilde{g}_{GF}^2 in eq. 5.4 is calculated for this new action. Alternatively one can consider the replacement of $\langle t^2 E(t) \rangle$ with $\langle t^2 E(t + \tau_0 a^2) \rangle$ as an improved operator for the energy density. In either case the t -shift changes the $\mathcal{O}(a^2)$ term of $g_{\text{GF}}^2(\mu; a)$. If we expand $\tilde{g}_{\text{GF}}^2(\mu)$ in $\tau_0 a^2$,

$$\tilde{g}_{\text{GF}}^2(\mu; a) = \frac{1}{\mathcal{N}} \langle t^2 E(t) \rangle + \frac{a^2 \tau_0}{\mathcal{N}} \left\langle t^2 \frac{\partial E(t)}{\partial t} \right\rangle, \quad (5.5)$$

and choose τ_0 such that the second term in eq. 5.5 cancels the $a^2 \mathcal{C}$ term in eq. 5.3, we remove the leading lattice artifacts

$$\tilde{g}_{\text{opt}}^2(\mu; a) = g_{\text{GF}}^2(\mu; a = 0) + \mathcal{O}(a^4[\log a]^n, a^4). \quad (5.6)$$

Full $\mathcal{O}(a^2)$ improvement through a systematic improvement program would require adding terms to improve the flow equation, the action, the boundary conditions, and the energy density operator $\langle t^2 E(t) \rangle$ [111]. Since our proposed improvement involves only a single parameter τ_0 , this τ_0 itself must depend on other parameters, most importantly on $\tilde{g}_{\text{GF}}^2(\mu)$ and on the bare coupling through the lattice spacing dependence of the term $\langle t^2 \frac{\partial E(t)}{\partial t} \rangle$ in eq. 5.5. Optimizing τ_0 both in the renormalized and bare couplings could remove the predictive power of the method. Fortunately, as we will see in the next section, our numerical tests indicate that it is sufficient to choose τ_0 to be a constant or only weakly $\tilde{g}_{\text{GF}}^2(\mu)$ dependent to remove most $\mathcal{O}(a^2)$ lattice artifacts.

Since the gradient flow is evaluated through numerical integration, the replacement $g_{\text{GF}}^2 \rightarrow \tilde{g}_{\text{GF}}^2$ can be done by a simple shift of t without incurring any additional computational cost. The optimal t -shift τ_{opt} can be identified by a simple procedure when the gradient flow is used for scale setting, which we will consider in a future publication. In this chapter we concentrate on the step scaling function and find the τ_{opt} that removes the $\mathcal{O}(a^2)$ terms of the discrete β function corresponding to scale change s ,

$$\beta_{\text{lat}}(g_c^2; s; a) = \frac{\tilde{g}_c^2(L; a) - \tilde{g}_c^2(sL; a)}{\log(s^2)}. \quad (5.7)$$

5.3 4 Flavor Test

We illustrate the t -shift improvement with the $N_f = 4$ SU(3) system. This theory was recently studied by refs. [51, 50] using gradient flow step scaling with staggered fermions. The 4-flavor SF running coupling was previously considered in Ref. [115] using $\mathcal{O}(a)$ -improved Wilson fermions, and in Ref. [97] using staggered fermions. In our calculations we use nHYP-smeared [71, 70] staggered fermions and a gauge action that includes an adjoint plaquette term in order to move farther away from a well-known spurious fixed point in the adjoint–fundamental plaquette plane [28]. As in Ref. [51] we impose anti-periodic BCs in all four directions, which allows us to carry out computations with exactly vanishing fermion mass, $m = 0$. For the discrete β function we consider the scale change $s = 3/2$ and compare lattice volumes $12^4 \rightarrow 18^4$, $16^4 \rightarrow 24^4$ and $20^4 \rightarrow 30^4$.

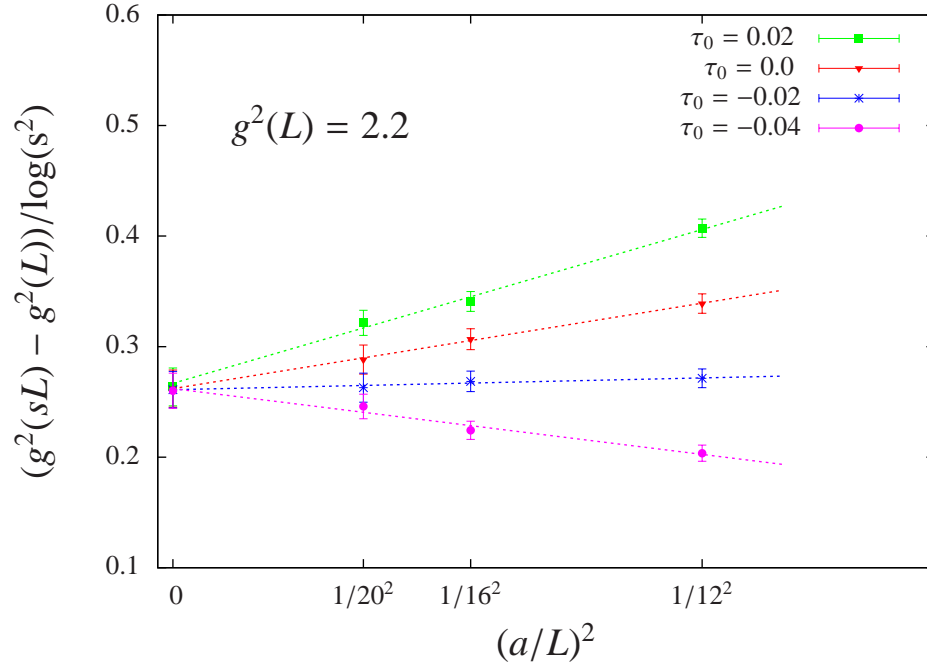


Figure 5.1: Continuum extrapolations of the discrete β_{lat} function of the $N_f = 4$ system at $\tilde{g}_c^2(L) = 2.2$ with several different values of the t -shift coefficient τ_0 . The dotted lines are independent linear fits at each τ_0 , which predict a consistent continuum value.

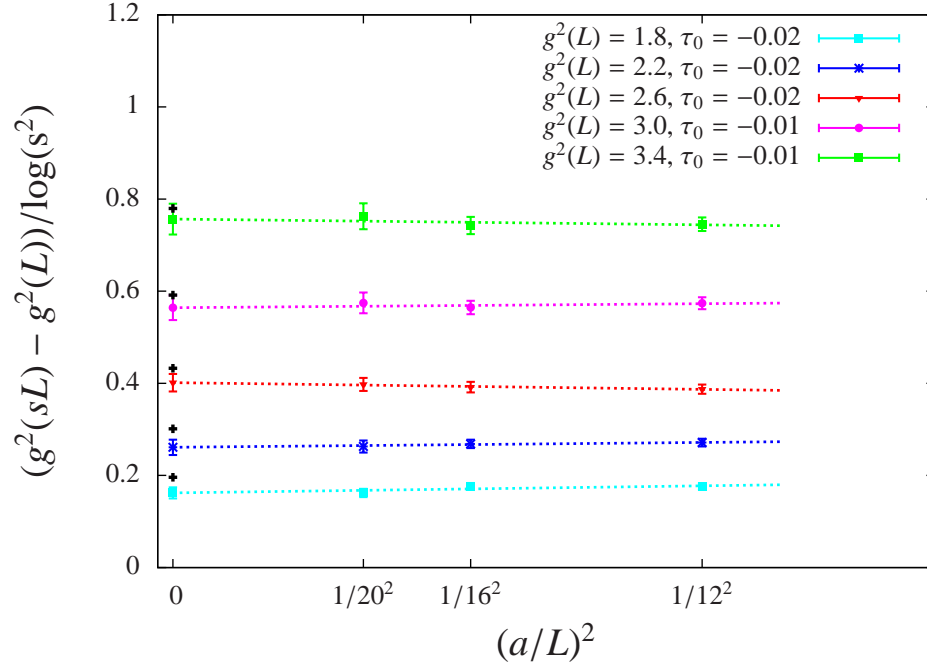


Figure 5.2: Continuum extrapolations of the discrete β_{lat} function of the $N_f = 4$ system for several different $\tilde{g}_c^2(L)$ values. For $\tilde{g}_c^2(L) = 1.8, 2.2$ and 2.6 $\tau_0 = -0.02$ is near-optimal, while the larger couplings $\tilde{g}_c^2(L) = 3.0$ and 3.4 require $\tau_0 = -0.01$ to remove most $\mathcal{O}(a^2)$ effects. The colored points at $(a/L)^2 = 0$ are the continuum extrapolated results, while the black crosses at $(a/L)^2 = 0$ show the corresponding two-loop perturbative predictions.

We accumulated 500–600 measurements of the gradient flow coupling, with each measurement separated by 10 molecular dynamics time units (MDTU), at 7–8 values of the bare gauge coupling on each volume. We consider the $c = 0.25$ scheme, as opposed to $c = 0.3$ used in Ref. [51], because smaller c gives better statistics at the expense of larger lattice artifacts. As discussed above, we aim to reduce these lattice artifacts through the non-perturbative improvement we have introduced. We follow the fitting procedure described in Ref. [115].

Full details of this study will be presented in Ref. [25]. Here we provide a representative illustration of the t -shift optimization. Figure 5.1 shows the dependence of the discrete β function on $(a/L)^2$ when $\tilde{g}_c^2(L) = 2.2$ with several values of the t -shift parameter τ_0 . The red triangles correspond to no improvement, $\tau_0 = 0$. The data are consistent with linear dependence on a^2 and extrapolate to 0.262(17), about 2σ below the two-loop perturbative value of 0.301. The slope of the extrapolation is already rather small, $b = 11(3)$. By adding a small shift this slope can be increased or decreased. With $\tau_0 = -0.02$ no $\mathcal{O}(a^2)$ effects can be observed – the corresponding slope is $b = 1.5(3.1)$ – and we identify this value as near the optimal τ_{opt} . The data at different τ_0 extrapolate to the same continuum value, even when the slope b is larger than that for $\tau_0 = 0$. This is consistent with the expectation that the t -shift changes the $\mathcal{O}(a^2)$ behavior of the system but does not affect the continuum limit. Since our action produces relatively small $\mathcal{O}(a^2)$ corrections even without improvement, the t -shift optimization has little effect on the continuum extrapolation, though the consistency between different values of τ_0 is reassuring.

It is interesting that the cut-off effects in our unimproved results, characterized by the slope b of the red triangles in Fig. 5.1, are more than three times smaller than those shown in fig. 4 of Ref. [51]. This difference grows to about a factor of four when we consider the larger $c = 0.3$ used in that study, suggesting that the t -shift optimization could have a more pronounced effect with the action used in Ref. [51]. The cause of the reduced lattice artifacts with our action is not obvious. Both our action and that used by Ref. [51] are based on smeared staggered fermions, though we use different smearing schemes. The different smearing might have an effect, as might the inclusion of the adjoint plaquette term in our gauge action. This question is worth investigating in the future.

In principle τ_{opt} could be different at different g_c^2 couplings but in practice we found little variation. Figure 5.2 shows near-optimal continuum extrapolations of the discrete β function at several values of $\tilde{g}_c^2(L)$. At each $\tilde{g}_c^2(L)$ the continuum extrapolated result is consistent within $\sim 2\sigma$ with the two-loop perturbative prediction, denoted by a black cross in Fig. 5.2. Comparable consistency with perturbation theory was found in previous studies [115, 97, 51, 50].

5.4 12 Flavor Results

We use the same lattice action with $N_f = 12$ as with $N_f = 4$ and consider six different volumes: 12^4 , 16^4 , 18^4 , 24^4 , 32^4 and 36^4 . This range of volumes allows us to carry out step scaling analyses with scale changes $s = 4/3$, $3/2$ and 2 . As for $N_f = 4$ we performed simulations in the $m = 0$ chiral limit with anti-periodic BCs in all four directions. Depending on the volume and bare coupling β_F we accumulated 300–1000 measurements of the gradient flow coupling g_c^2 for $0 \leq c \leq 0.5$, with 10 MDTU separating subsequent measurements. Here we will consider only $c = 0.2$. Full details of our ensembles and measurements, studies of their auto-correlations, and additional analyses for $c = 0.25$ and 0.3 will appear in Ref. [25]. The choice of $c = 0.2$ minimizes the statistical errors, and we find the IRFP in this scheme to be at a weaker coupling than for larger c , which is numerically easier to reach. The typical trade-off for these smaller statistical errors would be larger cut-off effects, but as discussed in previous sections these cut-off effects can be reduced by our non-perturbative improvement.

Figure 5.3 shows the running coupling $g_c^2(L)$ as the function of the bare gauge coupling β_F for different volumes. The interpolating curves are from fits similar to those in Ref. [115]. The curves from different volumes cross in the range $6.0 \leq \beta_F \leq 6.5$. The crossing from lattices with linear size L and sL defines the finite-volume IRFP coupling $g_\star^2(L; s)$:

$$g_c^2(L) = g_c^2(sL) \implies g_\star^2(L; s) = g_c^2(L). \quad (5.8)$$

If the IRFP exists in the continuum limit then the extrapolation

$$\lim_{(a/L)^2 \rightarrow 0} g_\star^2(L; s) \equiv g_\star^2 \quad (5.9)$$

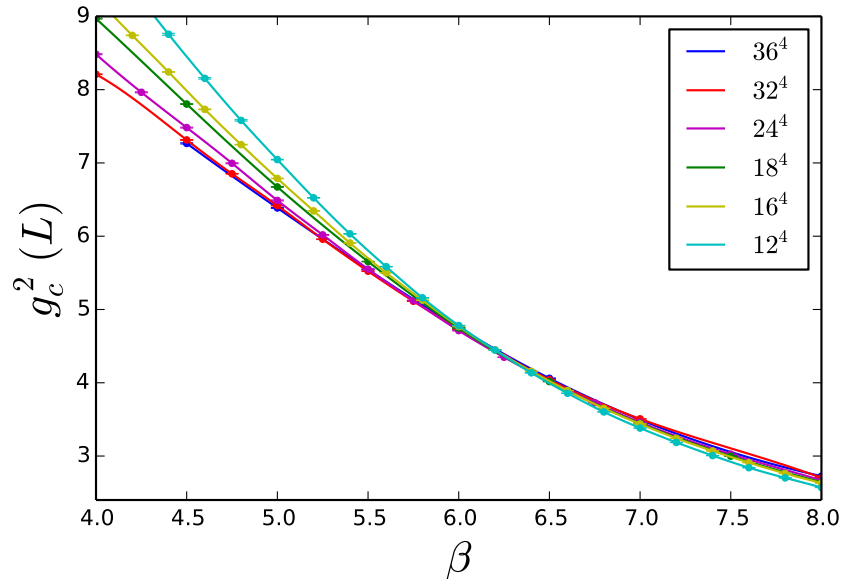


Figure 5.3: The $N_f = 12$ running coupling $g_c^2(L)$ versus the bare coupling β_F on several volumes, for $c = 0.2$. Crossings between results from different volumes predict the finite volume IRFP coupling $g_*^2(L)$ in this scheme.

has to be finite and independent of s .² Figure 5.4 illustrates the continuum extrapolation of $g_\star^2(L)$ with scale change $s = 2$ for various choices of the t -shift parameter τ_0 . The red triangles correspond to no shift, $\tau_0 = 0$. Their $(a/L)^2 \rightarrow 0$ continuum extrapolation has a negative slope, and the leading lattice cut-off effects are removed with a positive t -shift, $\tau_{\text{opt}} \approx 0.04$. A joint linear extrapolation of the $\tau_0 = 0, 0.02, 0.04$ and 0.06 results, constrained to have the same continuum limit at $(a/L)^2 = 0$, predicts $g_\star^2 = 6.21(25)$. However, these results all come from the same measurements, and are therefore quite correlated. While it is an important consistency check that the continuum limit does not change with τ_{opt} , just as for $N_f = 4$, the uncertainty in the continuum-extrapolated g_\star^2 from this joint fit is not reliable.

Instead, we should consider only the results with the near-optimal $\tau_{\text{opt}} \approx 0.04$. As we show in Fig. 5.5, $\tau_{\text{opt}} \approx 0.04$ is also near-optimal for scale changes $s = 3/2$ and $4/3$. None of these results have any observable $\mathcal{O}(a^2)$ effect, making the extrapolation to the continuum very stable. Each scale change predicts a continuum IRFP for $N_f = 12$. The three sets of results in Fig. 5.5 come from matching different volumes, making a joint fit legitimate. This continuum extrapolation predicts that the IR fixed point is located at renormalized coupling $g_\star^2 = 6.18(20)$ in the $c = 0.2$ scheme.

5.5 Summary

We have considered step scaling based on the gradient flow renormalized coupling, introducing a non-perturbative $\mathcal{O}(a^2)$ improvement that removes, or at least greatly reduces, leading-order cut-off effects. This phenomenological improvement increases our control over the extrapolation to the infinite-volume continuum limit, as we demonstrated first for the case of SU(3) gauge theory with $N_f = 4$ massless staggered fermions. Turning to $N_f = 12$, we found that the continuum limit was well defined and predicted an infrared fixed point even without improvement. Applying our proposed improvement reinforced this conclusion by removing all observable $\mathcal{O}(a^2)$ effects. For the finite-volume gradient flow renormalization scheme defined by $c = 0.2$, we find the continuum

² We thank D. N6gr6adi for useful discussions of the continuum limit.

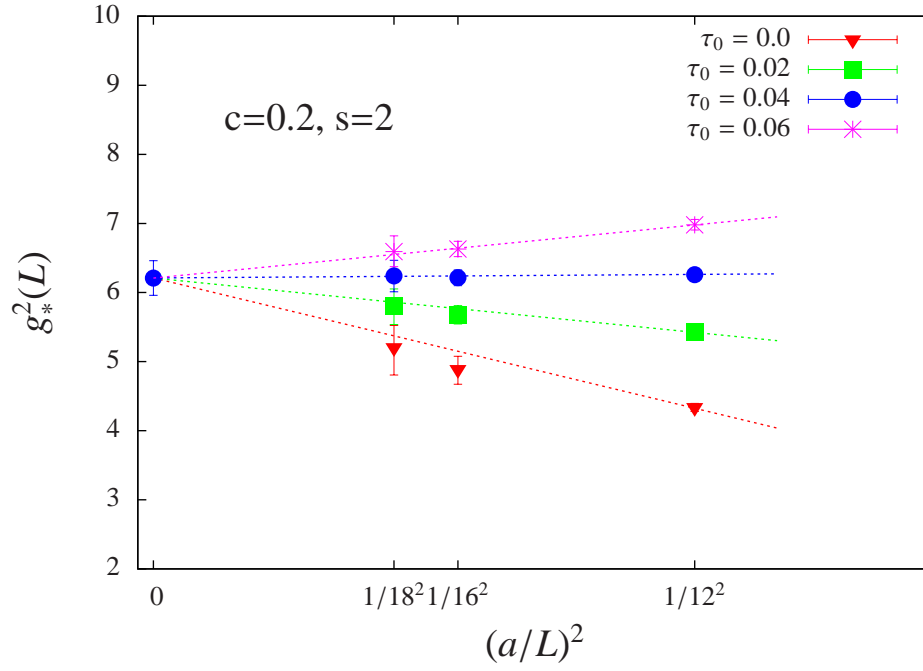


Figure 5.4: Continuum extrapolations of the 12-flavor finite volume IRFP $g_*^2(L)$, with several different t -shift coefficients τ_0 for fixed scale change $s = 2$. The dotted lines are a joint linear fit constrained to have the same $(a/L)^2 = 0$ intercept, which gives $g_*^2 = 6.21(25)$.

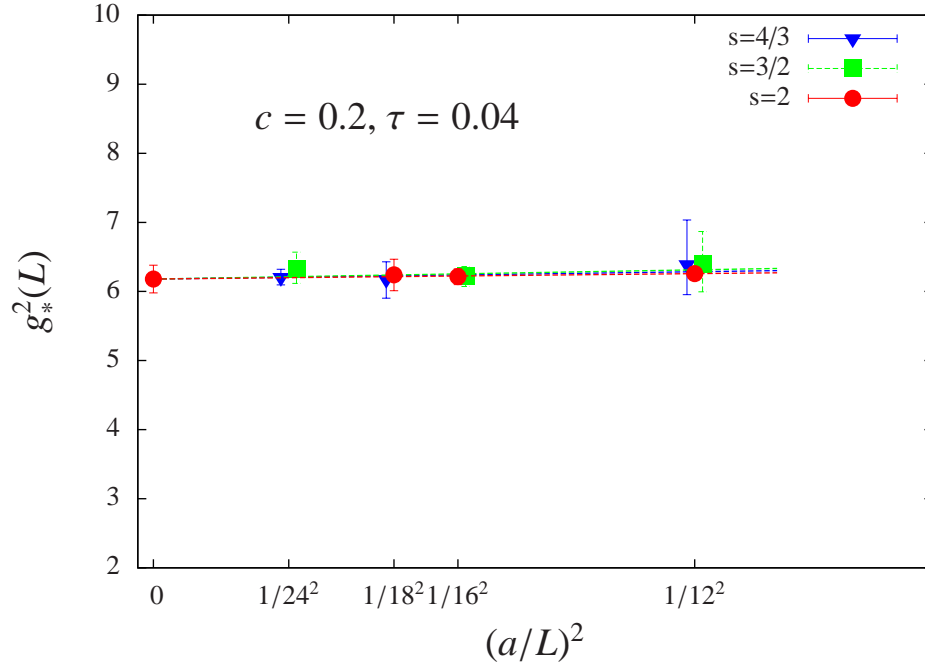


Figure 5.5: Continuum extrapolations of the 12-flavor finite volume IRFP $g_*^2(L)$, with several different scale changes for the near-optimal improvement coefficient $\tau_{\text{opt}} \approx 0.04$. The $s = 4/3$ and $3/2$ data points are horizontally displaced for greater clarity. The dashed lines are a joint linear fit constrained to have the same $(a/L)^2 = 0$ intercept, which gives $g_*^2 = 6.18(20)$.

conformal fixed point to be located at $g_\star^2 = 6.18(20)$.

The 12-flavor system has been under investigation for some time, and other groups have studied its step scaling function [86, 9, 10, 76]. However, this work is the first to observe an IRFP in the infinite-volume continuum limit. There are likely several factors contributing to this progress. While we did not invest more computer time than other groups, we have employed a well-designed lattice action. The adjoint plaquette term in our gauge action moves us farther away from a well-known spurious fixed point, while nHYP smearing allows us to simulate at relatively strong couplings. The gradient flow coupling itself appears to be a significant improvement over other schemes,³ and our non-perturbative improvement also contributes to obtaining more reliable continuum extrapolations.

Our non-perturbative improvement is general and easy to use in other systems. It does not rely on the lattice action or fermion discretization, though we suspect that the improvement may not be effective if there are $\mathcal{O}(a)$ artifacts, e.g. for unimproved Wilson fermions. Since $\mathcal{O}(a)$ -improved lattice actions are standard, this does not appear to be a practical limitation. We look forward to seeing our proposal applied both to QCD and to other conformal or near-conformal systems.

³ C.-J. D. Lin has told us about dramatic improvements in auto-correlations when using the gradient flow coupling compared to the twisted Polyakov loop coupling of Ref. [86].

Chapter 6

Wilson Flow MCRG

6.1 Introduction

For the past several years many lattice groups have been involved in studying strongly-coupled near-conformal gauge–fermion systems. Some of these models may be candidates for new physics beyond the standard model, while others are simply interesting non-perturbative quantum field theories. Because the dynamics of these lattice systems are unfamiliar, it is important to study them with several complementary techniques. Not only does this allow consistency checks, it can also provide information about the most efficient and reliable methods to investigate near-conformal lattice theories.

Monte Carlo Renormalization Group (MCRG) two-lattice matching is one of several analysis tools that we are using to investigate SU(3) gauge theories with many massless fermion flavors. This technique predicts the step-scaling function s_t in the bare parameter space. In a previous work [99] we proposed an improved MCRG method that exploits the Wilson flow to obtain a bare step-scaling function that corresponds to a unique discrete β function. We briefly review our Wilson-flow-optimized MCRG (WMCRG) procedure in Sections 6.2. It is important to note that we are investigating a potential infrared fixed point (IRFP) where the coupling is irrelevant: its running slows and eventually stops. This is challenging to distinguish from a near-conformal system where the gauge coupling runs slowly but does not flow to an IRFP. The observation of a backward flow that survives extrapolation to the infinite-volume limit could provide a clean signal. In Section 6.3 we report WMCRG results for SU(3) gauge theory with $N_f = 12$ flavors of massless

fermions in the fundamental representation.

This 12-flavor model has been studied by many groups, including Refs. [10, 39, 46, 7, 66, 33, 28, 80, 86, 6, 48, 49, 76, 27, 5, 69, 68, 26]. Using new ensembles of 12-flavor gauge configurations generated with exactly massless fermions, our improved WMCRG technique predicts a conformal IRFP where the step-scaling function vanishes. As with every method, it is essential to study the systematic effects. For WMCRG the most important systematic effects are due to the finite volume and limited number of blocking steps. While we are not able to carry out a rigorous infinite-volume extrapolation, the observed zero of the bare step-scaling function is present for all investigated lattice volumes and renormalization schemes, and agrees with the earlier MCRG results of Ref. [66]. The results of our complementary $N_f = 12$ investigations of finite-temperature phase transitions [107, 69], the Dirac eigenmode number [27, 26], and finite-size scaling [68] are also consistent with the existence of an infrared fixed point and IR conformality.

6.2 Wilson Flow Optimized MCRG

As an alternative to optimizing the RG blocking transformation, and thus changing the renormalization scheme at each coupling β_F , here we propose to use the Wilson flow to move the lattice system as close as possible to the renormalized trajectory of a fixed renormalization scheme.

The Wilson flow is a continuous smearing transformation [92] that can be related to the $\overline{\text{MS}}$ running coupling in perturbation theory [87]. Refs. [51, 50] recently used the Wilson flow to compute a renormalized step-scaling function in a way similar to Schrödinger functional methods. While this approach appears very promising, it is based on perturbative relations that are only fully reliable at weak coupling. Here we do not use this perturbative connection, instead applying the Wilson flow as a continuous smearing that removes UV fluctuations. The Wilson flow moves the system along a surface of constant lattice scale in the infinite-dimensional action-space; it is not a renormalization group transformation and does not change the IR properties of the system.

Our goal is to use a one-parameter Wilson flow transformation to move the lattice system as close as possible to the renormalized trajectory of our fixed RG blocking transformation. This

is shown in figure 6.1. We proceed by carrying out two-lattice matching after applying the Wilson flow for a flow time t_f on all lattice volumes. (The Wilson flow is run only on the unblocked lattices, not in between RG blocking steps.) As above, since we can block our lattices only a few times, we must optimize t_f by requiring that consecutive RG blocking steps yield the same $\Delta\beta_F$, as shown in 6.2. As for traditional MCRG, increasing the number of blocking steps reduces the dependence on the optimization parameter; in the limit $n_b \rightarrow \infty$, our results would be independent of t_f .

With Wilson-flowed MCRG we can efficiently determine bare step-scaling functions that correspond to unique RG β functions. The uniqueness of the β function is a result of using a fixed block transformation. In this work our block transformation consists of nHYP smearing the unblocked lattice and then multiplying adjacent links in the same direction.

The ability to study a unique β function opens up interesting directions for future studies. By comparing different β functions around the perturbative gaussian FP, we can study scaling violations in the lattice system. In IR-conformal systems, we can investigate the scheme-dependence of the β function near the IRFP, an issue explored in perturbation theory by Ref. [104].

6.3 12 Flavor Results

Our WMCRG results for the 12-flavor system are obtained on gauge configurations generated with exactly massless fermions. Our lattice action uses nHYP-smearred staggered fermions as described in Ref. [28], and to run with $m = 0$ we employ anti-periodic boundary conditions in all four directions. All of our analyses are carried out at couplings weak enough to avoid the unusual strong-coupling “ \mathcal{S}^4 ” phase discussed by Refs. [28, 69].

We perform three-lattice matching with volumes 6^4 – 12^4 – 24^4 and 8^4 – 16^4 – 32^4 . Three-lattice matching is based on two sequential two-lattice matching steps, to minimize finite-volume effects [66]. Both two-lattice matching steps are carried out on the same final volume V_f . We denote the number of blocking steps on the largest volume by n_b , and tune the length of the initial Wilson flow by requiring that the last two blocking steps predict the same step-scaling function. Using the 8^4 – 16^4 – 32^4 data we determine the bare step-scaling function for $n_b = 3$ and $V_f = 4^4$ as

well as $n_b = 4$ and $V_f = 2^4$, while the 6^4 – 12^4 – 24^4 data set is blocked to a final volume $V_f = 3^4$ ($n_b = 3$). This allows us to explore the effects of both the final volume and the number of blocking steps. We investigate three renormalization schemes by changing the HYP smearing parameters in our blocking transformation [99]: scheme 1 uses smearing parameters (0.6, 0.2, 0.2), scheme 2 uses (0.6, 0.3, 0.2) and scheme 3 uses (0.65, 0.3, 0.2).

Figs. 6.3, 6.4 and 6.5 present representative results for 12 flavors. All of the bare step-scaling functions clearly show $s_b = 0$, signalling an infrared fixed point, for every n_b , V_f and renormalization scheme. Appropriately for an IR-conformal system, the location of the fixed point is scheme dependent. We observe that the fixed point moves to stronger coupling as the HYP smearing parameters in the RG blocking transformation increase.

When we block our 8^4 , 16^4 and 32^4 lattices down to a final volume $V_f = 2^4$ (corresponding to $n_b = 4$), the observables become very noisy, making matching more difficult. The problem grows worse as the HYP smearing parameters increase, and our current statistics do not allow reliable three-lattice matching for $V_f = 2^4$ in schemes 2 and 3. To resolve this issue, we are accumulating more statistics in existing 32^4 runs, and generating additional 32^4 ensembles at more values of the gauge coupling β_F . These additional data will also improve our results for scheme 1, which we show in Fig. 6.5. Different volumes and n_b do not produce identical results in scheme 1, suggesting that the corresponding systematic effects are still non-negligible. We can estimate finite-volume effects by comparing $n_b = 3$ with $V_f = 3^4$ and $V_f = 4^4$. Systematic effects due to n_b can be estimated from $n_b = 4$ and $V_f = 2^4$, but this is difficult due to the noise in the 2^4 data. Even treating the spread in the results shown in Fig. 6.5 as a systematic uncertainty, we still obtain a clear zero in the bare step-scaling function, indicating an IR fixed point.

In this chapter we have shown how the Wilson-flow-optimized MCRG two-lattice matching procedure proposed in Ref. [99] improves upon traditional lattice renormalization group techniques. By optimizing the flow time for a fixed RG blocking transformation, WMCRG predicts a bare step-scaling function s_b that corresponds to a unique discrete β function. Applying WMCRG to new 12-flavor ensembles generated with exactly massless fermions, we observe an infrared fixed point

in s_b . The fixed point is present for all investigated lattice volumes, number of blocking steps and renormalization schemes, even after accounting for systematic effects indicated by Fig. 6.5. This result reinforces the IR-conformal interpretation of our complementary $N_f = 12$ studies of phase transitions [107, 69], the Dirac eigenmode number [27, 26], and finite-size scaling [68].

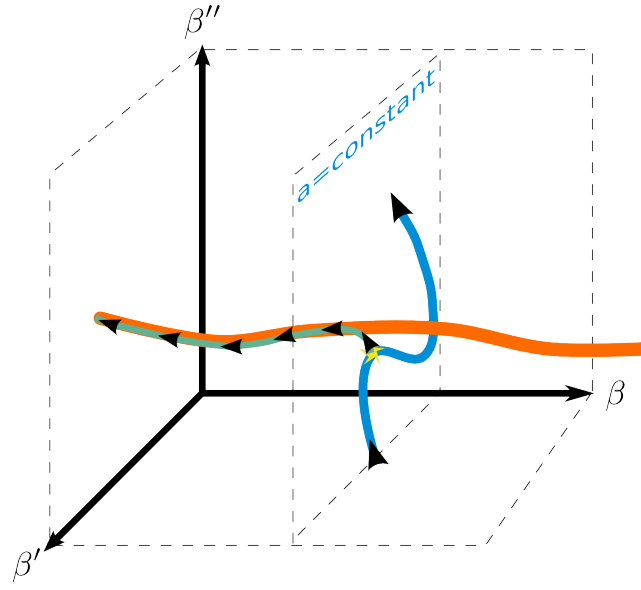


Figure 6.1: In Wilson Flow MCRG we use the Wilson Flow (blue) to approach the renormalized trajectory. An optimization step similar to that used in MCRG allows us to locate the flow time that gets us closest to the renormalized trajectory (orange). We then block our lattice using a fixed block transformation (green).

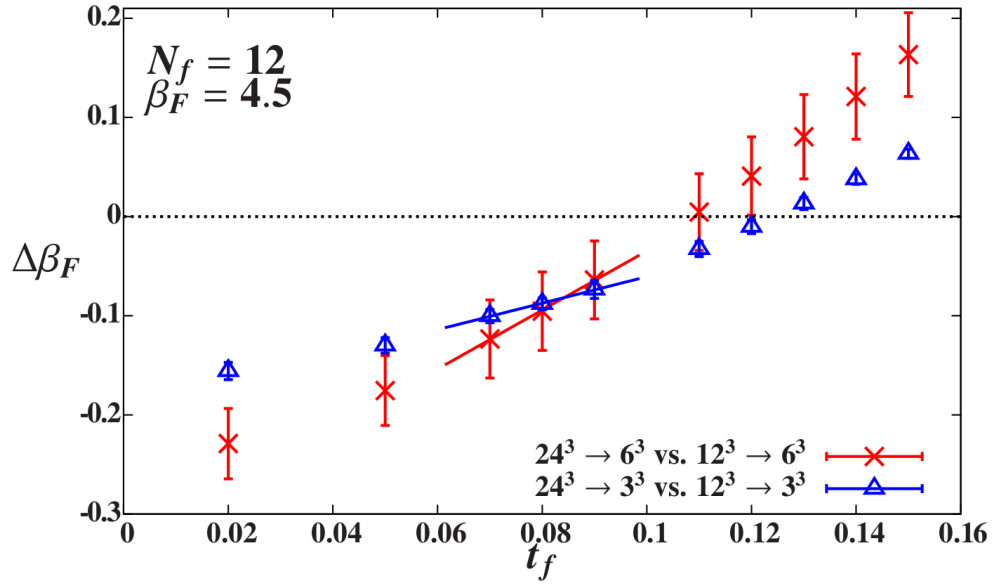


Figure 6.2: Optimization of the Wilson flow time t_f with fixed $\alpha = 0.5$, for $\beta_F = 4.5$. The uncertainties on the data points are dominated by averaging over the different observables.

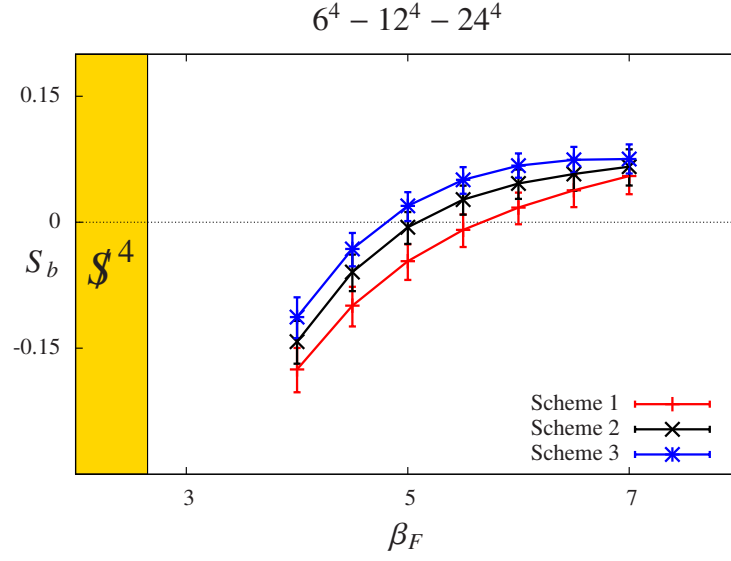


Figure 6.3: The bare step-scaling function s_b predicted by three-lattice matching with 6^4 , 12^4 and 24^4 lattices blocked down to 3^4 , comparing three different renormalization schemes. The error bars come from the standard deviation of predictions using the different observables.

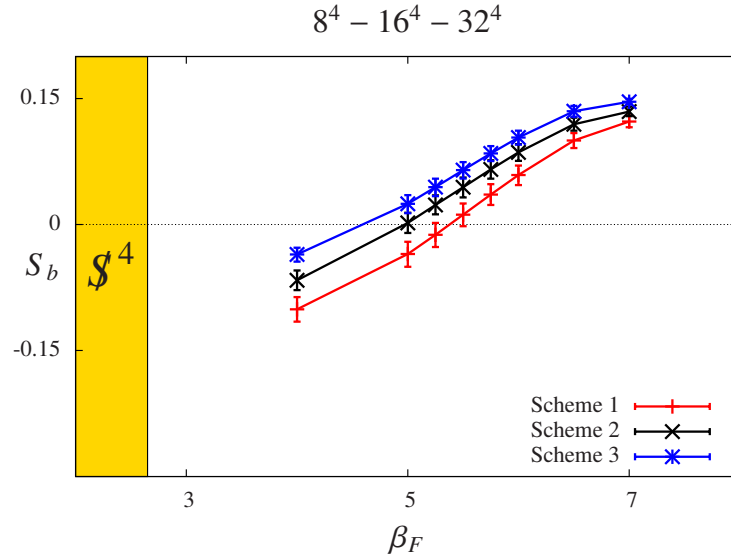


Figure 6.4: As in Fig. 6.3, the bare step-scaling function s_b for three different renormalization schemes from three-lattice matching, now using 8^4 , 16^4 and 32^4 lattices blocked down to 4^4 .

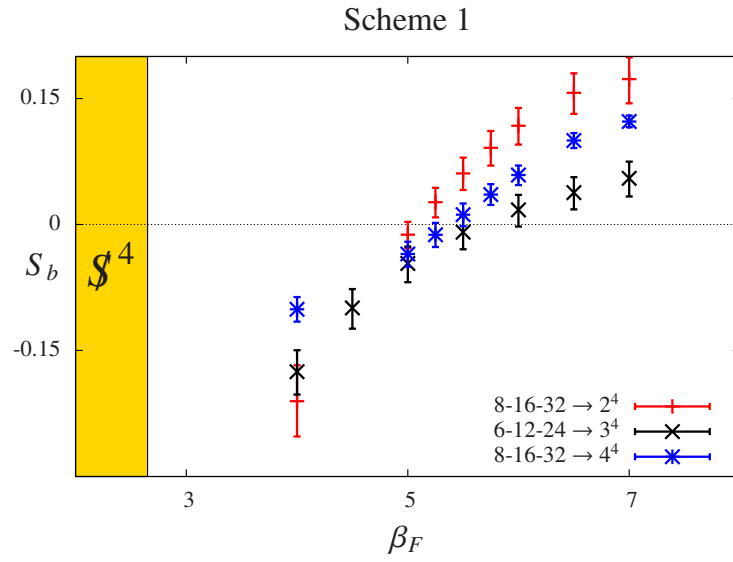


Figure 6.5: The bare step-scaling function s_b for scheme 1, comparing three-lattice matching using different volumes: 6^4 , 12^4 and 24^4 lattices blocked down to 3^4 (black \times s) as well as 8^4 , 16^4 and 32^4 lattices blocked down to 4^4 (blue bursts) and 2^4 (red crosses).

Chapter 7

Conclusion

Truly it is an exciting time in particle physics. For almost two decades since the discovery of the top quark no new fundamental particles were discovered in collider experiments. After a long wait the Higgs particle has been discovered, completing the standard model. Although the standard model is complete we know it must be an effective theory and our knowledge of the Higgs mass puts constraints on beyond standard model physics.

Many proposals for beyond standard model physics, including technicolor, are strongly coupled theories and thus inherently nonperturbative. Since the lattice offers the only controlled means of studying non perturbative field theories in a controlled manner, it is natural that strongly coupled beyond standard model is an active area of lattice research. Studying BSM physics on the lattices has created many new challenges for the lattice community to solve. Unlike QCD, we don't know the answer to most questions beforehand, and there are no experimental results to compare lattice results with. Furthermore just because a technique is successful when studying QCD does not mean that it will work just as well in systems that are very different from QCD. Accordingly it is important to approach each theory we study carefully and with an open mind. Only when several methods converge on the same result can that result be trusted. Additionally since there is more than one way to put a continuum theory on the lattice, it is important to understand the effect of the lattice action and lattice artifacts.

The lattice search for viable technicolor theories has focused on exploring gauge theories with $SU(N_C)$ colors and N_f fermions in some representation R . Changing these parameters generates

theories with dramatically different behavior. Theories with a small number of fermionic flavors in a lower representation behave similarly to QCD. If more fermionic degrees of freedom are added, the theory develops an infrared fixed point. If enough degrees of freedom are added the theory will lose asymptotic freedom. The location of the conformal window in the parameter space N_C, N_f , and R is fundamentally a question of strong dynamics. Perturbative and quasi perturbative calculations exist for the bounds of the conformal window but they can only serve as a guide to locate interesting theories for numerical studies.

Ultimately we are interested in the behavior of theories that may exhibit a slowly running coupling. This behavior, also called walking, may exist just below the conformal window. In a walking theory, γ_m must be $\mathcal{O}(1)$ for the theory to be phenomenologically successful.. Walking is widely believed to be necessary for any technicolor model to fit current experimental bounds on flavor changing neutral currents. To date a viable walking theory has not yet been found.

Many groups, including our own have explored the conformal window with several goals in mind. First understand the extent of the conformal window. Second improve lattice techniques in conformal systems. Third explore the bottom of the conformal window for walking behavior.

In this thesis I have discussed three methods that can be used to understand the β function. These methods are general and work for confining and conformal systems, they have the benefit that distinguishing between the two is straight forward. I first discussed MCRG and presented results for $SU(3)$ gauge theory with $N_f = 8$ and $N_f = 12$. Our results were consistent with the 12 flavor theory exhibiting an IRFP. The 8 flavor theory did not show a fixed point and appeared to be chirally broken and confining.

Next I introduced the gradient flow step scaling and showed results for $SU(3)$ gauge theory with $N_f = 4$ and $N_f=12$. The 4 flavor results were simply to show that our improvement works in a theory we know to be chirally broken. They also serve as a contrast to the 12 flavor results which clearly indicate an IRFP. Another group has recently applied this technique, with our improvement, to $SU(3)$ gauge theory with 8 flavors [72] their results show that while the theory runs slower than perturbation theory predicts, they can not locate an IRFP.

Finally I introduced an improvement to MCRG that uses the Wilson flow as an optimization. We call this technique Wilson Flow MCRG or WMCRG. WMCRG is like MCRG in that we are calculating the discrete step scaling function. By using the Wilson Flow as an optimization we are able to probe a single renormalized trajectory. Results from WMCRG for $SU(3)$ gauge theory of 12 flavors of fermions in the fundamental representation shows very clear evidence of a fixed point.

One of the benefits of these three techniques is that they are computationally inexpensive and do not require specific lattice dimensions or special boundary conditions. That is the lattices that we use for these step scaling studies are able to be used in other studies as well. Our group has been involved with several other analysis. We studied finite temperature phase transitions [107], finding a bulk phase transition in the 12 flavor theory and no such transition in the 8 flavor theory. We also used the Dirac eigenmodes to calculate γ_m over several length scales [26, 27]. Results from both of these studies are consistent with an IR conformal 12 flavor theory. For several years the IR physics of $SU(3)$ gauge theory with 12 flavors of fermions in the fundamental representation was a topic of debate. Several early papers drew different conclusions from their analysis. A consensus has emerged that the theory is indeed conformal. This consensus is a result of improvements in our understanding of quantum field theories as well as implements in lattice techniques used to study those quantum field theories.

In contrast to the 12 flavor theory, the results for 8 flavors show no sign of an IRFP. The fact that 8 flavors is chirally broken is not controversial. However, peculiar behavior in the eigenmode study seems to suggest that the running of the coupling is slow over a wide range of energy scales. These results seem to be confirmed by recent studies using our improved Wilson Flow step scaling function. It is still not clear if the running is simply slower than expected or is the long sought for walking scenario. Clearly $SU(3)$ gauge theory with 8 flavors of fermions in the fundamental representation is an interesting theory that should be studied further.

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