

COMPUTING THE ENERGY SCALE Λ_{YM} OF PURE
YANG-MILLS THEORY FROM A NOVEL LATTICE QCD
CODE WITH GRADIENT FLOW

by

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Abstract

The behavior of Quantum Chromodynamics, the quantum field theory that describes the strong interaction and the dynamics of quarks and gluons, at energies lower than a scale $\Lambda_{QCD} \approx 200 - 300$ MeV is not fully understood. The complexity of the problem stems from the exponential increase of the renormalized coupling constant towards low energies, which prevents a perturbative approach. Lattice QCD is a numerical method that has proven to be largely successful in describing the dynamics of strongly interacting quantum fields set on a discretized euclidean space time grid.

We developed a framework to perform lattice calculations of an SU(3) Yang-Mills theory with no dynamical fermion flavors. This theory, even if unphysical, presents most of the features of QCD and is generally a good approximation of it. In order to study divergent quantities and lattice discretization effects at low energies the recently proposed Gradient Flow method [1] has been used and smoothing properties of the flow on gauge fields at non zero flow time are presented.

The procedure for the discretization of the field theory on a four-dimensional lattice is discussed as well as the main algorithms needed for the calculations. Most notably the Metropolis algorithm, which is a Markov Chain Monte Carlo method, that can be used to generate statistical ensembles of lattice gauge field configurations. Some of the more technical details of the structure of the program that has been developed are also shown.

The implementation of the Gradient Flow method is proven to give results for simple topological properties that are consistent with ones found in literature, in particular we report for the topological susceptibility the value of $\chi^{1/4} = 186.9(4.9)$ MeV. These results, together with the analysis of the integrated autocorrelation time of the output data and the performance scaling properties of the program, suggest that the implementation is valid.

A method to estimate for the scale Λ_{YM} of fermionless Yang-Mills theory starting from the lattice results for the dimensionless quantity $t_f^2 \langle E \rangle$ has been developed. The challenge of finding an unbiased region in the energy spectrum in which the continuum limit extrapolation of $t_f^2 \langle E \rangle$ as a function of the energy scale $q = 1/\sqrt{8t_f}$ matches the perturbative expansion of the same quantity was solved by systematically fitting to all energy ranges within a window. The final result, taken as the weighted median of the fit parameters using the fit quality as weights, was found to be in agreement, for all loop-order corrections considered, with the one found in literature [2] using an entirely different method.

Chapter 1

Introduction

The theory that best describes the subatomic world as we know it is the Standard Model (SM) of particle physics. The interaction of quantum fields, the building blocks of the universe, are described within an extremely elegant mathematical framework. The particles, as they are most commonly intended, are nothing but ripples, local excitations, of the fields.

The strong interaction is one of the components of the Standard Model and arguably it is the most complicated of all. Quantum Chromodynamics (QCD) is the quantum field theory that describes it; the particle fields it considers are the quarks and the gauge fields (the mediators of the force) are the gluons. The reason why it is so difficult to study is because the “strength” of the force, given by the coupling constant, changes rapidly at different energies. The high-energy limit is relatively simple to describe: perturbative expansions in powers of the coupling constant can be performed, leading to advanced analytical results, for example the calculation of cross sections for collision processes. These computations are the key ingredient for understanding the phenomenology of particle accelerators.

Studying the low-energy limit of the strong interaction is a completely different problem. In a larger picture, the whole branch of physics called nuclear structure is fundamentally investigating the nature of the residual strong interactions between nucleons, protons and neutrons. These studies can be approached experimentally, by using data collected at Rare Isotope Beams facilities, to construct effective potentials between nucleons. These results are then used in many-body quantum mechanical simulations to get better and better insight on the underlying theory. Other approaches start from Effective Field Theories to model from “first principles” the nuclear interactions.

One of the biggest challenges in the theory of strong interactions is to link the description given by nuclear theory to the underlying Quantum Chromodynamics fundamental theory. The main problem is that the coupling constant of QCD becomes large at low energies. This makes perturbative calculations impossible to be carried out. Numerical methods are one of the possible solutions to the problem.

Lattice QCD is the most successful description of the strong forces at low energies that we know today. The key idea is to discretize spacetime in a four-dimensional grid of points. Quarks are set on sites of this lattice and physical properties are extracted by considering discretized paths

between different lattice sites. From the time it was first developed by K.G.Wilson, Lattice QCD has evolved dramatically, largely due to the increase of computational power over the years. Some of the most interesting calculations are still not technically feasible with the computational resources of today, this makes Lattice QCD a prominent field in the context of High Performance Computing (HPC).

1.1 Goals and Motivation

As this work is intended to be a Computational Physics thesis, the final goals of the project are dual. As far as computational science is concerned, the aim was to develop a computer program that implemented the rather recently proposed Gradient Flow Method [1] for Lattice QCD. This has been done with the idea of laying the foundations for a possible future more complete framework for performing lattice calculations. This larger program, however, is clearly too much to develop in the time frame of a master thesis, so the first building blocks have been implemented in this work.

The resulting lattice simulation program has been used to analyze a physical problem related to the scale Λ_{YM} of Pure Yang-Mills Theory. By that we mean the quantum field theory based on the SU(3) gauge symmetry group, the same of QCD, but with no dynamical fermion flavors. It is an interesting first approximation of the full theory, that however requires a simpler numerical implementation.

The specific problem we investigated regarded the matching between lattice calculations and analytical results from perturbation theory in the context of the Gradient Flow Method. We propose an estimate of the scale of Pure Yang-Mills theory starting from the lattice computation of the dimensionless quantity $t_f^2 \langle E \rangle$, which is related to the action of the gauge fields as they are evolved by the gradient flow equation. The same object has a known perturbative expansion at an energy scale dependent from the fictitious dimension t_f , the flow time, that we try to fit to the data extracted from our numerical simulation to estimate Λ_{YM} .

Other methods to compute the scale of Pure Yang-Mills theory have been previously studied in literature. The method we propose could be a simpler alternative to those. Future work is planned to extend the results of this thesis to the case of QCD and its scale Λ_{QCD} , using gauge field configurations generated with more sophisticated Lattice QCD codes, to check the validity of the method.

1.2 Our Contribution

This thesis, as every other scientific work, has its foundation in many other previously published work. It is then worth to state clearly what the author has contributed to.

The main contribution is the development of a completely novel Lattice code: starting from the very basics of SU(3) matrix arithmetics; to the sampling of the space of gauge field configuration via the Metropolis algorithm and finally to the implementation of a numerical integrator for the gradient flow equation.

There exist several other professional Lattice QCD code bases that are well-tested and highly optimized. The aim of this work is not to compete with these packages, but rather to gain some insight into the key methods and strategies as well to understand the challenges that the problem poses. The development process resulted in approximately 10000 lines of code written in C++ and this required most of the time spent on the thesis.

The second contribution can be summarized in the analysis of the feasibility of the estimation of Λ_{YM} from data collected with the lattice implementation of the gradient flow. This new method, which appears to produce consistent results, could be the subject of future studies.

1.3 Structure of the Thesis

This work is structured into three parts. In part I we will discuss the general features of QCD, chapter 2, to give the reader some of the basics of this complicated and very diverse theory. Elements of the discretization process needed to develop Lattice QCD and some basic definitions will be provided in chapter 3, while in chapter 4 more advanced topic such as the gradient flow and the reference scale fixing in Lattice QCD will be discussed.

In part II we introduce the numerical implementation we developed by presenting the main algorithms that have been used and the general structure of the code, chapter 5. In the following chapter the tests that have been performed to set up the algorithms and to check the performance scaling are presented.

Finally, in part III we first give an overview of the observables that have been computed on the ensembles of gauge field configurations that we have generated, chapter 7. Then in chapter 8 we present the procedure that has been used to estimate the scale Λ_{YM} and the final results that have been obtained.

Part I

Overview of QCD and Lattice Field Theories

Chapter 2

A Primer on QCD in the Continuum

The Standard Model of particle physics (SM) is the theory of fundamental particles and their interactions. Three of the four known fundamental forces are described by it, the exception is gravity, so that all phenomena regarding the electromagnetic, weak and strong forces are included within the theory.

The SM describes quantized fields defined on all space-time whose excitations are commonly identified with particles. There are two categories of fields, depending on the spin statistic they can be: fermions, quarks or leptons in the SM, which are the constituents of matter; or bosons, which mediate the interactions between particles. A summary of the fundamental particles of the SM, grouped by their properties and with some basic data, is shown in fig. 2.1

From a group theory point of view, the SM is the composition of three different local gauge symmetry groups, each associated with a fundamental force:

$$\text{SU}(3)_C \times \underbrace{\text{SU}(2)_L \times \text{U}(1)_Y}_{\text{broken to } \text{SU}(2)_W \times \text{U}(1)_Q} . \quad (2.1)$$

The $\text{U}(1)_Q$ symmetry group is associated with the electromagnetic interaction and $\text{SU}(2)_W$ is the weak interaction. They both derive from the symmetry $\text{SU}(2)_L \times \text{U}(1)_Y$, which is spontaneously broken through the Higgs Mechanism, .

Quantum Chromodynamics, commonly referred to as QCD, is the quantum field theory, contained in the SM, that describes the behavior of strongly interacting matter, that is quarks and gluons. It is a non-abelian gauge theory based on an $\text{SU}(3)$ symmetry group. The associated quantum number is called “color charge” which pictorially can assume the values of *red* (r), *green* (g), *blue* (b), *anti-red* (\bar{r}), *anti-green* (\bar{g}) or *anti-blue* (\bar{b}). In this chapter, we will discuss the QCD Lagrangian density, its properties and some of the major results of the theory. Some intermediate knowledge of Quantum Field Theory is assumed, derivations and proofs mainly follow the reasoning found in [4].

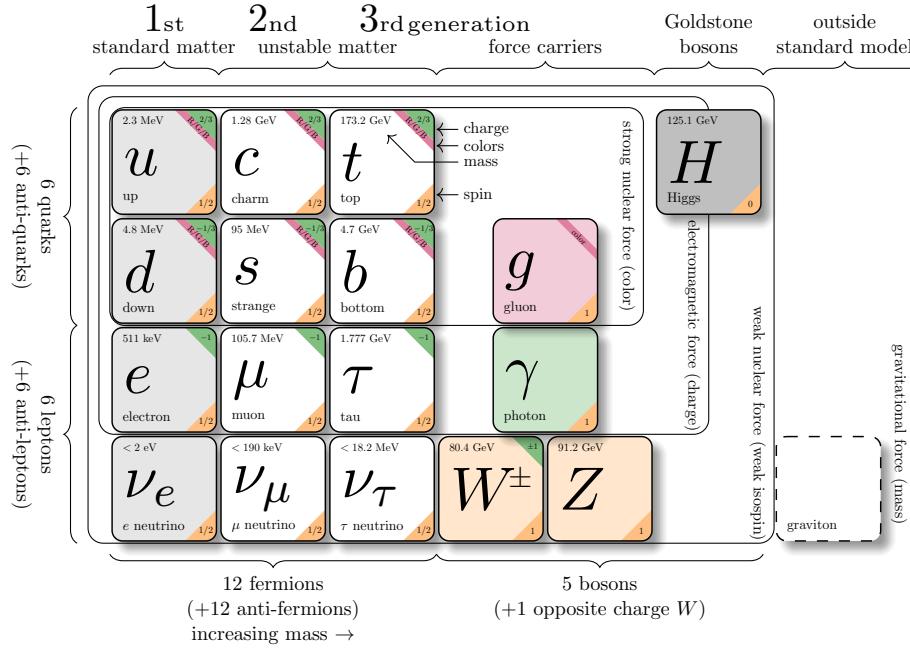


Figure 2.1: Summary of the particles in the Standard Model [3]

2.1 The QCD Lagrangian

In Quantum Field Theory (QFT) the characterizing equation of a theory is its Lagrangian density because it contains all the information about the fields that are involved, their properties and most importantly, their interactions. QCD is a non-abelian gauge theory based on the SU(3) local gauge symmetry group. The simplest Lagrangian is the Yang-Mills Lagrangian:

$$\mathcal{L}_{QCD} = \sum_{f=1}^{N_f} \bar{\psi}_f (i\gamma^\mu D_\mu - m_f) \psi_f - \frac{1}{4} (G_{\mu\nu}^a)^2. \quad (2.2)$$

Here ψ_f and $\bar{\psi}_f$ represent the complex-valued fermion fields of flavor f , with mass m_f . These are associated with the quark fields and come in six flavors: u (up), d (down), s (strange), c (charm), b (bottom) and t (top). The second element in the Lagrangian is the Gluon Field Strength Tensor, $G_{\mu\nu}^a$. The symbols μ and ν are Lorentz indices while a refers to the indices of the generators of the gauge group, SU(3) in this case. The Dirac's matrices γ^μ are also introduced. Note that Einstein's summing convention on repeated indices is implicit, for example when taking the square of the field strength tensor three sums are applied. The definition of $G_{\mu\nu}^a$ is:

$$G_{\mu\nu}^a = \frac{i}{g_0} [D_\mu, D_\nu] = \partial_\mu A_\nu^a - \partial_\nu A_\mu^a + g_0 f^{abc} A_\mu^b A_\nu^c. \quad (2.3)$$

In this equation A_μ^a is the gluon field that carries a Lorentz index and a group generator index, g_0 is the bare coupling constant of the strong interaction and the f^{abc} are the structure constants of SU(3), which satisfy:

$$[t^a, t^b] = i f^{abc} t^c \quad (2.4)$$

with t^a being the generators of the algebra $\mathfrak{su}(3)$. The covariant derivative D_μ is defined to be:

$$D_\mu = \partial_\mu - ig_0 t^a A_\mu^a. \quad (2.5)$$

With the information contained in the Lagrangian density, the behavior and the interactions of all fields are set.

2.1.1 Feynman Rules of QCD

A key element that is needed to perform perturbative calculations in a quantum field theory are Feynman rules. These are a set of equations and rules that represent the propagation of fields and the interaction vertices of the theory. For the case of QCD, being a non-abelian gauge theory, some vertices represent interactions between gauge bosons only, as opposed to Quantum Electrodynamics, QED, that forbids photon-photon interactions.

To begin with, we need to write out all of the terms of the Lagrangian density individually. We assume only one quark flavor as no term in the Lagrangian can change this quantum number,

$$\begin{aligned} \mathcal{L}_{QCD} &= \bar{\psi}(i\gamma^\mu D_\mu - m)\psi - \frac{1}{4}(G_{\mu\nu}^a)^2 \\ &= \bar{\psi}(i\gamma^\mu \partial_\mu + g_0 \gamma^\mu t^a A_\mu^a - m)\psi \end{aligned} \quad (2.6)$$

$$\begin{aligned} &\quad - \frac{1}{4} \left(\partial_\mu A_\nu^a - \partial_\nu A_\mu^a + g_0 f^{abc} A_\mu^b A_\nu^c \right) \left(\partial_\mu A_\nu^a - \partial_\nu A_\mu^a + g_0 f^{ade} A_\mu^d A_\nu^e \right) \\ &= \bar{\psi}(i\gamma^\mu \partial_\mu - m)\psi \\ &\quad - \frac{1}{4} (\partial_\mu A_\nu^a - \partial_\nu A_\mu^a)^2 \\ &\quad - g_0 f^{abc} (\partial_\mu A_\nu^a) A^{b\mu} A^{c\nu} \\ &\quad - \frac{1}{4} g_0^2 f^{abc} f^{ade} A_\mu^b A_\nu^c A^{d\mu} A^{e\nu} \\ &\quad + g_0 \gamma^\mu t^a A_\mu^a \bar{\psi}\psi. \end{aligned} \quad (2.7)$$

It is convention to represent the terms of the Lagrangian in a pictorial way, using what are called Feynmann diagrams. In the usual notation, fermions are drawn by a solid arrow, whose orientation is set based on whether it represents a particle field ψ or an anti-particle field $\bar{\psi}$. Gluons are drawn as springs and present no orientation. The terms in the Lagrangian that contain only two fields (the first two in eq. (2.7)) define the propagators, as they connect two points in space. All the other terms represent vertices of interaction between the fields. The first thing to do is to define the quark and gluon propagators, which can be obtained from the first and second terms respectively of the last equality:

$$\begin{aligned} i \xrightarrow[k]{\quad\quad\quad} j &= \frac{i(\gamma^\mu k_\mu + m)}{k^2 - m + i\epsilon} \delta_{ij} \\ a_\mu \xrightarrow[k]{\quad\quad\quad} b_\nu &= -\frac{i}{k^2 + i\epsilon} \delta_{ab} g^{\mu\nu} \end{aligned}$$

with $g^{\mu\nu}$ being the metric tensor of Minkovskian space-time. The arrow above the fields define the momentum k direction. In the fermion propagator the indices i, j stand for the color. In the gluon propagator a and b represent the indices for the generators of the $\mathfrak{su}(3)$ algebra while μ, ν are Lorentz indices. For propagators these pairs of indices must all match between the start and end point, this explains the introduction of the two Kronecker deltas $\delta_{i,j}$ and $\delta_{a,b}$. The last three terms in eq. (2.7) represent the interaction vertices of QCD. In particular the third term is a three gluon vertex:

$$= g_0 f^{abc} [g^{\mu\nu}(p-q)^\rho + g^{\nu\rho}(q-k)^\mu + g^{\rho\mu}(k-p)^\nu].$$

Then the four-gluon vertex:

$$= -ig_0^2 [f^{abe} f^{acd} (g^{\mu\nu} g^{\sigma\rho} - g^{\mu\rho} g^{\nu\sigma}) + f^{abd} f^{ace} (g^{\mu\nu} g^{\sigma\rho} - g^{\mu\sigma} g^{\nu\rho}) + f^{abc} f^{aed} (g^{\mu\sigma} g^{\nu\rho} - g^{\mu\rho} g^{\nu\sigma})].$$

Finally from the last term in eq. (2.7) we have the gluon-quark interaction vertex:

$$= ig_0 t^a \gamma^\mu.$$

Note that the Lagrangian we considered, and the resulting Feynman rules, are over-simplified: the “full” Lagrangian contains terms from Faddeev-Popov ghosts, that are introduced to fix the problems that arise when fixing the gauge; and counter-terms from the renormalization procedure. Nevertheless, interesting qualitative features can be inferred from the Feynman rules: the fact that gluons interact with each other through the three- and four-gluon vertices; the non

flavor-changing interaction between gluons and quarks, which decouples completely different quark flavors; the “color-changing nature” of the quark-gluon interaction, given by the t^a matrix in the vertex term that shows how the color state of a fermion is changed by the absorption or emission of a gluon.

2.1.2 Gauge Symmetry of the Lagrangian

The concept of gauge invariance is crucial in the construction of a discretized lattice theory from the continuum one, it is then worth considering how it is introduced. For the QCD Lagrangian to be physical, it must be locally gauge invariant, not only on a global scale. Let’s consider a quark field $\psi(x)$, which truly is a triplet of fermion fields each with different color quantum numbers:

$$\psi(x) = \begin{pmatrix} \psi_r(x) \\ \psi_b(x) \\ \psi_g(x) \end{pmatrix}. \quad (2.8)$$

A local gauge transformation in the internal space of SU(3) is a unitary transformation of this three-component vector, or in simpler terms: a rotation in color space. Defining $\Omega(x)$ one such local transformation applied to $\psi(x)$ one has:

$$\psi(x) \rightarrow \psi'(x) = \Omega(x)\psi(x), \quad (2.9)$$

and for the Dirac adjoint:

$$\bar{\psi}(x) \rightarrow \bar{\psi}'(x) = \bar{\psi}(x)\Omega^\dagger(x). \quad (2.10)$$

The transformation can be parametrized in terms of some functions $\alpha^a(x)$, one for each generator of the group:

$$\Omega(x) = \exp[i\alpha^a(x)t^a]. \quad (2.11)$$

This internal space rotation if applied to the simple Dirac free-field Lagrangian would generate an additional term:

$$\begin{aligned} \mathcal{L}_{Dirac} &= \bar{\psi}(i\gamma^\mu\partial_\mu - m)\psi \rightarrow \mathcal{L}'_{Dirac} = \bar{\psi}\Omega^\dagger(i\gamma^\mu\partial_\mu - m)(\Omega\psi) \\ &= \bar{\psi}'(i\gamma^\mu\partial_\mu - m)\psi' + i\bar{\psi}'\gamma^\mu\psi(\partial_\mu\Omega). \end{aligned} \quad (2.12)$$

The extra term implies that the Dirac Lagrangian is not gauge invariant. In order to fix this problem a new field $A_\mu(x)$, the gauge field, is introduced. It is convention to include it into the definition of the covariant derivative, so that ∂_μ becomes $D_\mu = \partial_\mu - ig_0 A_\mu(x)$, as previously stated. In geometrical terms, D_μ represents the derivative along the tangent vectors of the manifold on which the field is defined. The field $A_\mu(x)$ is algebra-valued and, to be consistent with the previous section one needs to project it on a basis of generators of the group: $A_\mu(x) = A_\mu^a(x)t^a$ (remember the implicit sum).

The transformation rule for $A_\mu(x)$ is set in order to cancel the extra term in eq. (2.12) exactly. We impose:

$$D_\mu\psi(x) \rightarrow (D_\mu\psi(x))' = (\partial_\mu - ig_0 A'_\mu(x))\Omega(x)\psi(x) = \Omega(x)(D_\mu\psi(x)), \quad (2.13)$$

which fixes the transformation for $A_\mu(x)$ to:

$$A_\mu(x) \rightarrow A'_\mu(x) = \Omega(x) \left[A_\mu(x) + \frac{i}{g_0} \partial_\mu \right] \Omega^\dagger(x). \quad (2.14)$$

Computing the derivative acting on Ω^\dagger is not trivial, as the matrices in the exponent of eq. (2.11) do not necessarily commute. For an infinitesimal transformation we can expand the matrix Ω in powers of α as $\Omega(x) = \exp[i\alpha^a(x)t^a] \approx 1 + i\alpha^a(x)t^a$ and get the following for the gauge field transformation, for a single component:

$$A_\mu^a(x) \rightarrow A'_\mu^a(x) = A_\mu^a(x) + \frac{i}{g_0} \partial_\mu \alpha^a(x) + f^{abc} A_\mu^b(x) \alpha^c(x) \quad (2.15)$$

which is the last element needed. One can now look at all the possible gauge invariant objects that can be constructed with the fields ψ and A of dimension 4, the same of the Lagrangian. Apart from the one already present in the Dirac Lagrangian with the covariant derivative, there are only two additional terms that are gauge invariant and of dimension 4 and they both can be taken by considering the gauge field tensor:

$$G_{\mu\nu}^a \equiv \frac{i}{g_0} [D_\mu, D_\nu] = \partial_\mu A_\nu^a - \partial_\nu A_\mu^a + g_0 f^{abc} A_\mu^b A_\nu^c, \quad (2.16)$$

which is clearly a gauge invariant object, since it is the commutator of covariant derivatives, of dimension 2. One can now construct the gauge field kinetic term, the well known $-\frac{1}{4}(G_{\mu\nu}^a)^2$ and reconstruct eq. (2.2). However, there is an additional term, not included in the QCD Lagrangian, that is the “dual term”, or “theta term” which will be interesting further in the work. It is defined as $\theta G_{\mu\nu}^a \tilde{G}^{a\mu\nu}$ where $\tilde{G}^{a\mu\nu} = \epsilon^{\mu\nu\rho\sigma} G_{a\rho\sigma}$ is the dual of the field tensor and $\epsilon_{\mu\nu\rho\sigma}$ the anti-symmetric Levi-Civita tensor of rank 4. With it the Lagrangian becomes:

$$\mathcal{L}_{QCD} = -\frac{1}{4}(G_{\mu\nu}^a)^2 + \bar{\psi}(i\cancel{D} - m)\psi + \theta G_{\mu\nu}^a \tilde{G}^{a\mu\nu}. \quad (2.17)$$

The theta term is usually neglected because there is no experimental evidence of it, but in principle it cannot be excluded. It is the simplest CP violating term that can be added to the QCD Lagrangian and for this is of particular interest in the study of CP phenomena, like the nucleon electric dipole moment (EDM)[5].

2.2 General Properties of QCD

Quantum Chromodynamics exhibits a set of features as a theory that is common to all non-abelian gauge theories. We have already seen one, that is the direct interaction of the gauge bosons, something that is not allowed in abelian theories as QED. Other interesting properties emerge when trying to renormalize the theory and are in general linked to the fixing of the scale, which leads to the concept of running coupling. In the particular case of QCD, the coupling constant at in a low-energy regime leads to Confinement, while in the high-energy limit Asymptotic Freedom emerges.

2.2.1 Running Coupling

The Renormalization Group Equation (RGE) defines the rate at which the renormalized coupling g varies as the renormalization scale μ changes, through the beta function of the coupling constant, $\beta(g)$, defined as:

$$\beta(g) = \frac{d}{d\log(\mu)} g(\mu), \quad (2.18)$$

For a generic non-abelian theory $SU(N)$ one can expand the β function in orders of g as:

$$\beta(g) = b_0 g^3 + b_1 g^5 + b_2 g^7 + \dots \quad (2.19)$$

One can then integrate up to arbitrary order eq. (2.18) and get an expression for $g(\mu)$. The coefficients b_i are obtained from computing the contribution of higher and higher order diagrams to the coupling. The value of b_0 , from 1-loop corrections, is:

$$b_0 = -\frac{1}{(4\pi)^2} \left(\frac{11}{3}N - \frac{2}{3}N_f \right). \quad (2.20)$$

Here N is the dimension of the gauge group $SU(N)$ and N_f the number of flavors of the fermions. The minus sign in front implies that any non-abelian theory with a sufficiently small number of fermions, less than $\frac{11}{2}N$ (that is ≤ 16 for QCD), is “asymptotically free”, meaning that at high energy the coupling vanishes and particles don’t feel any interaction.

The RGE is usually expressed in terms of the analog of the fine structure-constant for the strong force, $\alpha_s = g^2/4\pi$. The first order solution is given by plugging eq. (2.20) into eq. (2.18) and integrating. In terms of α_s at a momentum scale μ we get:

$$\alpha_s(\mu) = \frac{\alpha_s^0}{1 + \frac{b_0 \alpha_s^0}{4\pi} \log(\mu^2/M^2)}, \quad (2.21)$$

here α_s^0 is the value of the coupling at an energy M , which is set by the integration. The typical choice for μ when measuring the running coupling is the mass of the Z boson, an experimentally very well known value. This mass a is sufficiently large energy scale for perturbation theory to be applied safely to QCD.

The coupling depends on the arbitrarily chosen renormalization point M , a convenient choice is then to rewrite the equation defining a scale Λ that satisfies:

$$1 = \frac{b_0 \alpha_s^0}{4\pi} \log(M^2/\Lambda^2). \quad (2.22)$$

This simplifies eq. (2.21) to its well-known form, correct up to one loop corrections:

$$\alpha_s(\mu) = \frac{4\pi}{b_0 \log(\mu^2/\Lambda^2)}. \quad (2.23)$$

In fig. 2.2 a higher order approximation of $\alpha_s(\mu)$ is plotted against some selected experimental values.

In section 4.4.1 we will show a more precise result, correct up to four-loop corrections, for the running coupling $\alpha_s \mu$. The scale, often referred to as Λ_{QCD} , is the energy at which the

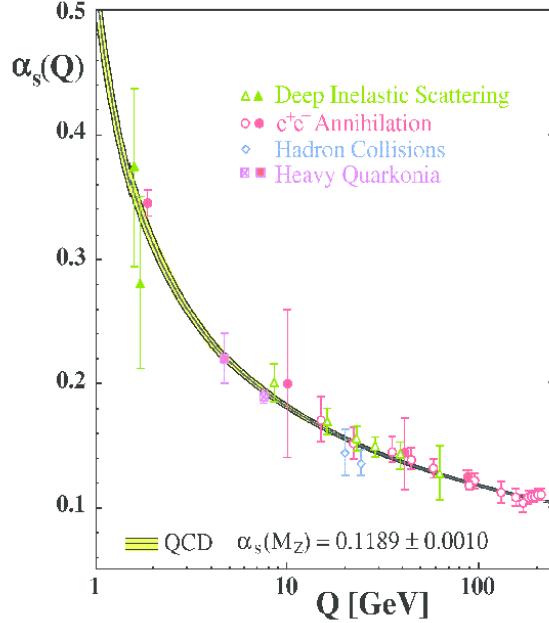


Figure 2.2: The strong coupling as a function of the energy scale. The data points are experimental results; the black solid line and yellow bands represent the QCD prediction using the reported value of the coupling at the mass of the Z boson (α_s^0 in our notation). Image taken from [6]

interactions become strong and perturbation theory becomes invalid. We can see that for μ approaching the value of Λ eq. (2.23) diverges. Experimental values [7] suggest that $\Lambda_{QCD} \approx 200 - 300$ MeV, meaning that perturbation theory can be applied from momenta roughly above the 1 GeV scale, where $\alpha_s \approx 0.4$.

One of the main goals of this thesis is to find a simple and not numerically expensive way to estimate the scale parameter Λ from lattice calculations. In this work the scale of pure Yang-Mills theory, which we denote as Λ_{YM} is studied. This is the case of no fermion flavors, but in future work the determination of Λ_{QCD} with 2 or $2 + 1$ dynamical fermions will also be studied.

Asymptotic Freedom

Equation (2.23) is a clear indication that QCD at high energies has a small coupling. Asymptotic freedom is the property of gauge theories, QCD is usually the example of it, that causes the interactions between the fields to become weaker as the energy scale increases, this is because of the inverse log dependence of the coupling on the energy. Grand Unification Theories (GUT), for example, are based on the fact that there exists an energy scale at which the strong force has a coupling equivalent to the one of the electroweak interaction.

The discovery of asymptotic freedom by Gross, Wilczek and Politzer [8, 9], was used, in the late Seventies when the fundamental theory was still debated, as an indication that QCD is indeed the correct theory.

Asymptotically free theories can be analyzed perturbatively at sufficiently large energies and are believed to be consistent up to any energy scale.

Confinement

In general, color charged particles cannot be isolated, so that quarks and gluons are not detectable alone, but always in the form of hadrons: colorless objects formed of multiple quarks. Hadrons can be of two types: mesons (quark-antiquark) and baryons (three quarks or three anti-quarks). Glueballs, combinations of gluons such that the total is colorless, are also in principle allowed, but have not been observed yet. Confinement is the phenomenon for which it is not possible to isolate a color charge, single quarks or gluons, from a hadron without producing other new hadrons. Single colored particles within very small time scales undergo hadronization, the process of spawning new quarks or anti-quarks from the vacuum to balance the total color charge and produce colorless matter. This is the reason for the fundamentally different nature of the strong force compared to the other forces, no perturbative expansion can be made at low energies. The exact proof of how confinement is introduced in non-abelian gauge theories like QCD is yet not known. Qualitatively, however, it can be explained by the fact that the gauge bosons of the theory, the gluons, carry color charge just like the matter fields, the quarks.

At low energies from eq. (2.23) we can infer that the coupling constant increases exponentially, so particles that have momenta comparable to the scale Λ_{QCD} interact very strongly with the gauge field and any other particle. This implies, for example, that the force between two particles does not vanish for long distances, but instead it increases. The usual picture is to consider the gluons being exchanged by two quarks at rest. These gluons would form flux-tubes between the sources that, if stretched by separating the quarks, eventually store enough energy to make a quark-anti-quark pair energetically favorable, as illustrated in the following figure.

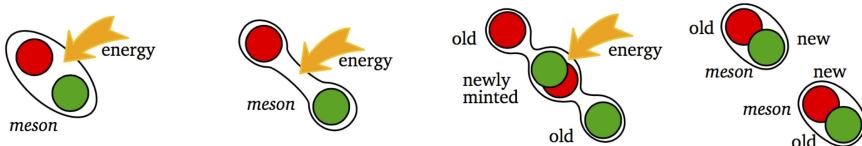


Figure 2.3: Representation of confinement using flux-tubes . As two color sources are pulled apart, the energy stored in the color field between the sources increases so much that a $q\bar{q}$ pair is formed from the vacuum energy.

One could also look at the inverse of the energy scale, which is a distance in natural units, of roughly $\Lambda_{QCD}^{-1} \approx 1$ fm, that is approximately the size of light hadrons, a further proof that quark sources cannot be torn apart easily.

The strong low-energy interaction is the cause for the large discrepancy between the mass of the baryons and the mass of their constituent quarks. Considering a proton for example, its constituent quark (two up-quarks and one down-quark) have masses that sum up to 10 MeV. However, the total mass of the proton is 938 MeV, this means that the 99% of its mass is given by the binding energy of quarks and gluons.

2.3 Methods and Regimes of Chromodynamics

Given the very different behaviors of the strong force at different energy scales, QCD needs to be dealt with in various ways depending on the scale of interest. At high energies, perturbation theory can be applied safely, but at low ones no expansion in the coupling constant can be made.

Perturbative QCD

At high energies, like the scale of the large particle accelerators that are currently available, the QCD coupling constant is sufficiently small to allow analytical calculations of Feynman diagrams to be meaningful. Hadronization is neglected as the time scale is small enough to consider it a post-collision process. For example, at a scale where the strong force is comparable to the electroweak interaction, the cross section of $e^+e^- \rightarrow \text{hadrons}$ compared to that of $e^+e^- \rightarrow \mu^+\mu^-$ can be used to measure the number of quark flavors that are active below a certain energy. In fig. 2.4 the perturbative QCD (pQCD) ratio of the hadronic to muonic cross section for the electron-positron annihilation is shown compared to experimental data. A step is visible around 3 – 4 GeV and it is caused by the triggering of the bottom quark sector in the process.

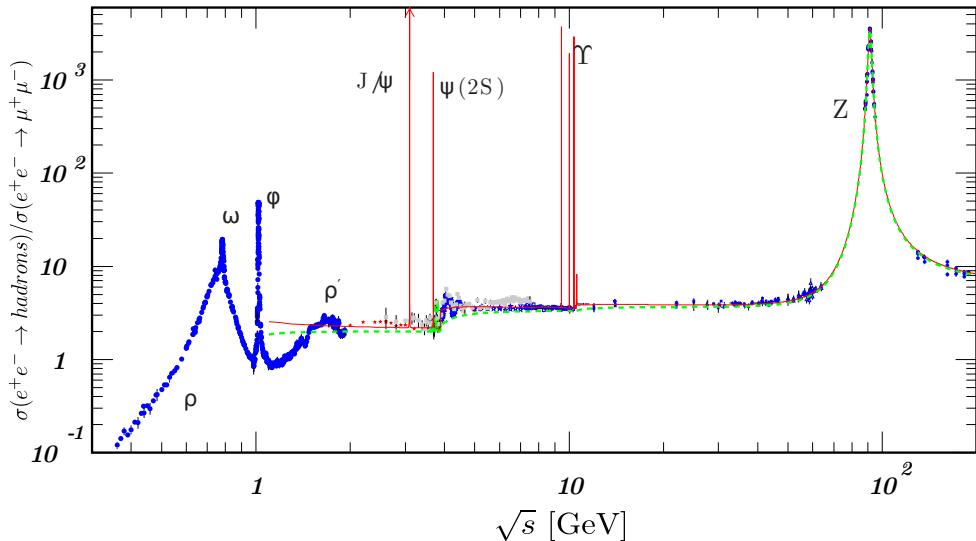


Figure 2.4: Ratio of the experimental cross section of $e^+e^- \rightarrow \text{hadrons}$ and $e^+e^- \rightarrow \mu^+\mu^-$ as collected by the Particle Data Group [10]. Naive quark model is the dashed green line and pQCD results are also shown as a solid red line. pQCD appears to agree well to experiment in non resonant regions of the energy spectrum.

Lattice Methods

To deal with the non-perturbative sector of the strong interaction the most widespread approach is to use numerical simulations, in particular, lattice methods, it is common to refer to it as Lattice QCD. The main idea, which will be expressed more in detail in Chapter 3, is to discretize spacetime and evaluate the quantum field only at fixed sites on a hyper-cubic lattice. Physical quantities are then computed stochastically on ensembles of such gauge fields. This approach, however, is very expensive from a computational point of view and so far no calculation at physical quark masses with a sufficiently small lattice spacing, in principle closer to the continuum theory, have been performed.

Effective Field Theories

An interesting problem is to link QCD directly with nuclear forces, that are long-range remnants of the strong interaction at a hadron level. The problem is of high interest because there is yet no fundamental theory of nuclear interactions from first principles. The most common approach is to define an Effective Field Theory (EFT), starting from a low energy approximation of chromodynamics, that preserves most of the symmetries of the underlying theory.

Chiral EFT (χ EFT) is one of the most popular of such theories; it starts from considering nucleons as a fundamental SU(2) group in isospin and describes the interactions between them through the exchange of pions [11]. It promotes particles that are not fundamental in QCD to the basic blocks of a low-energy effective Lagrangian, with nucleons as the fermion fields and pions as Nambu-Goldstone bosons [12] of the theory. Its Lagrangian is constructed in a systematic manner considering all possible interaction vertices between hadrons [13]:

$$\mathcal{L}_{\chi EFT} = \mathcal{L}_{\pi\pi} + \mathcal{L}_{\pi N} + \mathcal{L}_{NN} + \text{(three hadron terms)} + \dots \quad (2.24)$$

where $\mathcal{L}_{\pi\pi}$ is the term describing the dynamics between pions, \mathcal{L}_{NN} represents the term for interactions between two nucleons and $\mathcal{L}_{\pi N}$ the one between one pion and one nucleon and so on. Each term is further expanded in powers of Q/Λ_χ where Q is the pion momentum and Λ_χ is the energy scale at which the theory breaks down because of the pions having an energies comparable to the mass of the nucleons. One constructs order by order all possible terms:

$$\begin{aligned} \mathcal{L}_{NN} &= \mathcal{L}_{NN}^{(0)} + \mathcal{L}_{NN}^{(2)} + \mathcal{L}_{NN}^{(4)} + \dots \\ \mathcal{L}_{\pi N} &= \mathcal{L}_{\pi N}^{(1)} + \mathcal{L}_{\pi N}^{(2)} + \mathcal{L}_{\pi N}^{(3)} + \mathcal{L}_{\pi N}^{(4)} + \dots \\ \mathcal{L}_{\pi\pi} &= \mathcal{L}_{\pi\pi}^{(2)} + \mathcal{L}_{\pi\pi}^{(4)} + \dots \end{aligned} \quad (2.25)$$

one then computes all possible diagrams order by order and truncates the expansion when needed. This is mostly useful for computing reduced matrix elements for nuclear many-body calculations [14].

One interesting remark, as seen in fig. 2.5, is that the approach of χ EFT spontaneously generates interaction matrix elements for three or more nucleons.

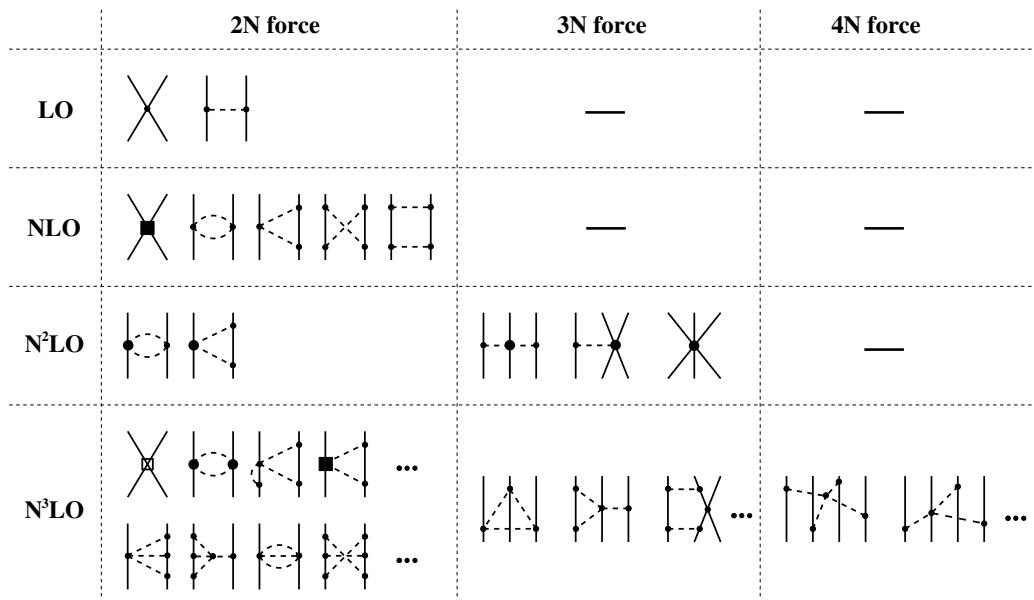


Figure 2.5: Diagrams of the leading order terms in χEFT [13], organized by truncation error. The solid lines represent nucleons, dashed lines are for pions. Different types of vertices are shown with different symbols.

Chapter 3

Lattice Field Theories and Lattice QCD

Lattice QCD is one of the main Lattice Field Theories. It deals with the strong force at low energies, handling non-perturbative calculations in a numerical way. The idea of discretizing spacetime in a lattice and perform calculations of field theories was proposed by Wilson in 1974 [15] as an alternative method to study the confinement of quarks. It has proven to be the most systematic approach to non-perturbative theories like QCD. In this chapter we will describe briefly how this discretization procedure is performed and some of the main computational strategies that are involved.

3.1 Discretizing Field Theories

The starting point for Lattice QCD is Feynman's path-integral formalism, but expressed in Euclidean spacetime, through a Wick rotation. An observable of some field ϕ is then given by:

$$\langle O[\phi] \rangle = \frac{1}{Z[\phi]} \int \mathcal{D}[\phi] O[\phi] e^{-S_E[\phi]}, \quad (3.1)$$

and the euclidean correlator of two quantities $O_1[\phi]$ and $O_2[\phi]$ between times t_1 and t_2 is:

$$\langle O_1[\phi](t_1)O_2[\phi](t_2) \rangle = \frac{1}{Z[\phi]} \int \mathcal{D}[\phi] O_1[\phi(t_1)]O_2[\phi(t_2)] e^{-S[\phi]}, \quad (3.2)$$

where the partition function $Z[\phi]$ is defined as:

$$Z = \int \mathcal{D}[\phi] e^{-S_E[\phi]}, \quad (3.3)$$

and $S_E[\phi]$ is the euclidean action of the field:

$$S_E[\phi] = \int d^4x \mathcal{L}[\phi], \quad (3.4)$$

Evaluating path-integrals is not possible in general with analytical tools so, in order to allow numerical computations, the Euclidean spacetime is discretized on a lattice, fixing the value of the field at only specific and equally spaced position in spacetime.

3.1.1 The Quantum Mechanical Example

To better understand the discretization procedure for quantum field theories, it is useful to look at the simpler case of a quantum mechanical system, for example of a single particle in a potential V . The Hamiltonian of the system is simply:

$$\hat{H} = \frac{\hat{p}^2}{2m} + \hat{V}, \quad (3.5)$$

where we used natural units ($\hbar = c = 1$). In path integral formalism the propagator between a state $|x_i\rangle$ at t_i to another state $|x_f\rangle$ at t_f can be expressed as:

$$\langle x_f | e^{-\hat{H}(t_f - t_i)} | x_i \rangle = \int \mathcal{D}x(t) e^{-S_E[x]}, \quad (3.6)$$

with $S_E[x]$ being the euclidean action:

$$S_E[x(t)] = \int_{t_i}^{t_f} dt \mathcal{L}[x(t)] = \int_{t_i}^{t_f} dt \frac{m}{2} \dot{x}(t)^2 + V[x(t)]. \quad (3.7)$$

Note that in the path integral in eq. (3.6) the exponent contains no imaginary unit. This is because we are using the euclidean action, related to the classical action by a Wick rotation $it \rightarrow t$.

To compute the propagator numerically one can chose to discretize the time dimension into slices t_0, t_1, \dots, t_N of distance a one from another. A path then becomes a vector of positions at fixed time points:

$$x = \{x(t_0), x(t_1), \dots, x(t_N)\}. \quad (3.8)$$

The integral is then changed from a continuous definition to an $N - 2$ dimensional integral (the boundary points are fixed):

$$\int \mathcal{D}x(t) e^{-S_E[x](t)} \rightarrow \frac{1}{A} \int dx_1 dx_2 \dots dx_{N-1} e^{-S_{lat}[x]}, \quad (3.9)$$

where A is some normalization. The classical action is now only computed on the fixed points of the discretized time coordinate. In particular we can approximate the integral between t_j and t_{j+1} , using the finite difference discretization for the derivative of $x(t)$ and the average potential, as:

$$\int_{t_j}^{t_{j+1}} dt \mathcal{L}[x(t)] \approx a \left[\frac{m}{2} \left(\frac{x_{j+1} - x_j}{a} \right)^2 + \frac{1}{2} (V[x_{j+1}] - V[x_j]) \right]. \quad (3.10)$$

With this approximation the action of one path, or “configuration”, is the sum of the approximated integrals from:

$$S_{lat}[x] \approx \sum_{i=0}^{N-1} \left[\frac{m}{2a} (x_{j+1} - x_j)^2 + aV(x_i) \right]. \quad (3.11)$$

This discretization procedure for a simple quantum mechanical system (whose space coordinate can be seen as a one-dimensional scalar field) allows to compute the path integrals in eqs. (3.1) and (3.2) using, in the case for the expectation value for example:

$$\begin{aligned}\langle O[x] \rangle &= \frac{1}{Z[x]} \int \mathcal{D}[x] O[x] e^{-S_E[x]} = \frac{1}{Z[x]} \int dx_1 dx_2 \dots dx_{N-1} O[x] e^{-S_{lat}[x]} \\ &= \int dx_1 dx_2 \dots dx_{N-1} P[x] O[x],\end{aligned}\quad (3.12)$$

where we have identified $P[x] = \frac{1}{Z[x]} e^{-S_{lat}[x]}$ as the probability distribution function of the observable O . One can then compute the expectation value numerically by choosing random configurations according the specified PDF, that depends on the action and compute the expectation value as the average of the observable over the chosen configurations.

3.2 Discretization of QCD on the Lattice

In the case of QCD the discretization procedure is more complicated than the one shown in the previous section for a one dimensional harmonic oscillator, because there are two types of field at play. Moreover the gluon gauge field A and the N_f fermionic quark fields ψ are defined in four-dimensional spacetime. In order to perform numerical calculations on the lattice, it is necessary to discretize the Euclidean action for both fields on a hyper-cubic lattice Λ :

$$\Lambda = \{n = (n_x, n_y, n_z, n_t)\} \text{ where } n_x, n_y, n_z = 1, 2, \dots, N; \quad n_t = 1, 2, \dots, N_T. \quad (3.13)$$

The choice of the lattice spacing, usually denoted a , is arbitrary, but most often it is chosen to be equal for all dimensions, space and time alike. Our discretized fields are then constrained to have values only on the points an instead of on a continuum spacetime x^μ :

$$\phi(x) \xrightarrow{\text{discretization}} \phi(an). \quad (3.14)$$

In this section the discretization is performed following a structure that resembles section 2.1.2, that is to start from the fermion fields and imposing gauge invariance.

3.2.1 Naïve Discretization of Fermions

The main ideas for the discretization procedure of fermions that we present are taken from [16], quoting the main intermediate steps that lead to the formulation of Lattice QCD. The reviews [17, 18] have also been useful references.

The starting point is the fermionic euclidean action in the continuum:

$$S_F[\psi, \bar{\psi}] = \int d^4x \bar{\psi}(x) (\gamma_\mu \partial^\mu + m) \psi(x), \quad (3.15)$$

now we discretize the euclidean spacetime on a lattice Λ of spacing a , each point will be denoted with n . The partial derivative can be turned into the central finite difference between

neighboring points along the direction of the derivative:

$$\partial_\mu \psi(x) \rightarrow \frac{\psi(n + \hat{\mu}) - \psi(n - \hat{\mu})}{2a}. \quad (3.16)$$

The discretized fermion action is then the sum of the lagrangian density over all lattice sites multiplied by the unit volume a^4 :

$$S_F[\psi, \bar{\psi}] = a^4 \sum_{n \in \Lambda} \bar{\psi}(n) \left[\sum_{\mu=1}^4 \gamma_\mu \frac{\psi(n + \hat{\mu}) - \psi(n - \hat{\mu})}{2a} + m\psi(n) \right]. \quad (3.17)$$

As we did in section section 2.1.2, we try to apply a local gauge transformation $\Omega(n)$ to the field. It is simple to show that some terms of the derivative in the action are not gauge invariant:

$$\bar{\psi}(n)\psi(n \pm \hat{\mu}) \rightarrow \bar{\psi}(n)\Omega^\dagger(n)\Omega(n \pm \hat{\mu})\psi(n \pm \hat{\mu}). \quad (3.18)$$

The problem is in the way the derivative was discretized and the solution is the introduction of an additional field that has the correct transformation laws. This field must transform under a local gauge transformation in a way that connects the values of the fermion field at two different lattice sites. Because of this property we denote the auxiliary field as “link variables”, later the matching with the gauge field will be introduced, and by convention it is written as $U(n, n + \hat{\mu}) = U_\mu(n)$. Furthermore, it must depend on the orientation along the μ direction, from eq. (3.18):

$$U(n, n - \hat{\mu}) \equiv U_{-\mu}(n) = U_\mu(n - \hat{\mu})^\dagger. \quad (3.19)$$

In particular, we want it to transform as:

$$\begin{aligned} \bar{\psi}(n)U_\mu(n)\psi(n + \hat{\mu}) &\rightarrow \bar{\psi}(n)\Omega^\dagger(n)U'_\mu(n)\Omega(n + \hat{\mu})\psi(n \pm \hat{\mu}) \\ \bar{\psi}(n)U_{-\mu}(n)\psi(n - \hat{\mu}) &\rightarrow \bar{\psi}(n)\Omega^\dagger(n)U'_{-\mu}(n)\Omega(n - \hat{\mu})\psi(n \pm \hat{\mu}), \end{aligned} \quad (3.20)$$

from which we infer that:

$$\begin{aligned} U_\mu(n) &\rightarrow U'_\mu(n) = \Omega(n)U_\mu(n)\Omega^\dagger(n + \hat{\mu}) \\ U_{-\mu}(n) &\rightarrow U'_{-\mu}(n) = \Omega(n)U_{-\mu}(n)\Omega^\dagger(n - \hat{\mu}). \end{aligned} \quad (3.21)$$

the field $U_\mu(n)$ is an SU(3) valued field, that can be intuitively thought of as a set of oriented links between the sites of the lattice. A visual representation is given in fig. 3.1,

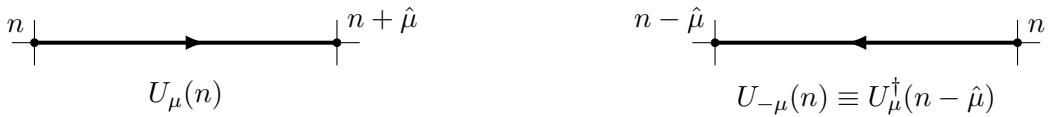


Figure 3.1: Schematic representation of the link variables $U_\mu(n)$ and $U_{-\mu}(n)$ on the lattice.

With this results we can write a gauge invariant lattice fermion action as:

$$S_F[\psi, \bar{\psi}] = a^4 \sum_{n \in \Lambda} \bar{\psi}(n) \left[\sum_{\mu=1}^4 \gamma_\mu \frac{U_\mu(n)\psi(n + \hat{\mu}) - U_{-\mu}(n)\psi(n - \hat{\mu})}{2a} + m\psi(n) \right]. \quad (3.22)$$

3.2.2 The Gauge Transporter and the Wilson Loop

For a more formal definition of the link variables we can look at the continuum gauge transporter $G(x, y)$, which is defined as the path-ordered product, denoted with \mathcal{P} , of a gauge field $A_\mu(x)$ along some curve \mathcal{C} between two points x and y in spacetime:

$$G(x, y) = \mathcal{P} \exp \left(ig_0 \int_{\mathcal{C}} A_\mu(x') dx'^\mu \right). \quad (3.23)$$

An important thing to note now is that gauge transporter belongs to the gauge group, $SU(3)$ for QCD, and not to the algebra, $\mathfrak{su}(3)$. Up to order $\mathcal{O}(a)$ the integral on the straight line connecting two points on the lattice can be approximated with $aA_\mu(n)$ for small lattice spacings. For convenience the gauge field is scaled by a factor $1/g_0$, giving:

$$\tilde{A}_\mu(x) = \frac{1}{g_0} A_\mu(x), \quad (3.24)$$

but for notation's sake we'll drop the tilde from now on. One is then left with:

$$\begin{aligned} G(n, n + \hat{\mu}) \equiv U_\mu(n) &= \exp(iA_\mu(n)) = \mathbb{1} + iaA_\mu(n) + \mathcal{O}(a^2) \\ G(n, n - \hat{\mu}) \equiv U_{-\mu}(n) &= \exp(-iaA_\mu(n)) = \mathbb{1} - iaA_\mu(n - \hat{\mu}) + \mathcal{O}(a^2), \end{aligned} \quad (3.25)$$

where in the second equation, for the negative direction, the approximation is taken on the arrival point instead on the initial one. With this definition is easy to see the relation we just stated earlier on the orientation change:

$$U(n, n - \hat{\mu}) = U_{-\mu}(n) = U_\mu^\dagger(n). \quad (3.26)$$

One can now define a Wilson loop as a gauge transporter that has the same start and final point. In the continuum case it can be shown, see for example ref. [4], that the trace of such an object is gauge invariant. On the lattice the minimal Wilson loop is just a square on a fixed plane $\mu\nu$, and we call this a Plaquette:

$$\begin{aligned} P_{\mu\nu}(n) &= U_\mu(n)U_\nu(n + \hat{\mu})U_{-\mu}(n + \hat{\mu} + \hat{\nu})U_{-\nu}(n + \hat{\nu}) \\ &= U_\mu(n)U_\nu(n + \hat{\mu})U_\mu^\dagger(n + \hat{\nu})U_\nu^\dagger(n), \end{aligned} \quad (3.27)$$

a pictorial representation of this equation is given in fig. 3.2.

3.2.3 The Wilson Gauge Action

In the continuum theory the Wilson loop is related to the square of the field strength tensor, the same term that enters the Lagrangian. The plaquette is therefore a good starting point for the creation of a discretized version of the Euclidean action for the gluon field.

Recalling the Baker-Campbell-Hausdorff (BCH) formula for the product of exponentials of matrices:

$$\exp(A) \exp(B) = \exp \left(A + B + \frac{1}{2}[A, B] + \dots \right), \quad (3.28)$$

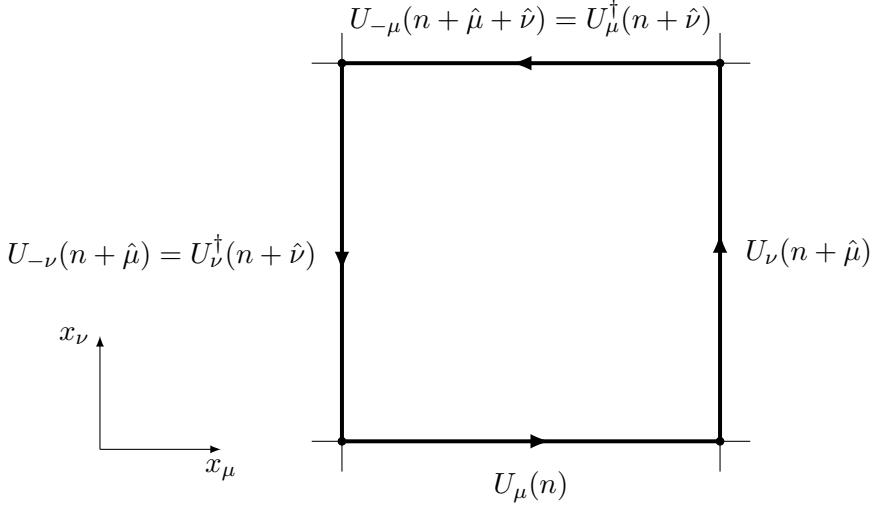


Figure 3.2: The Plaquette, as defined on the lattice, is the oriented product of the link variables of a minimal square in the plane $\mu\nu$.

and by inserting eq. (3.25) repeatedly into the definition of the plaquette one has:

$$\begin{aligned} P_{\mu\nu}(n) &= \exp \left[iaA_{\mu}(n) + iaA_{\nu}(n + \hat{\mu}) - iaA_{\mu}(n + \hat{\nu}) - iaA_{\nu}^{\dagger}(n) - \frac{a^2}{2}[A_{\mu}(n), A_{\mu}(n + \hat{\mu})] \right. \\ &\quad - \frac{a^2}{2}[A_{\nu}(n + \hat{\mu}), A_{\nu}(n)] + \frac{a^2}{2}[A_{\mu}(n), A_{\nu}(n)] + \frac{a^2}{2}[A_{\nu}(n + \hat{\mu}), A_{\mu}(n + \hat{\nu})] \\ &\quad \left. + \frac{a^2}{2}[A_{\mu}(n), A_{\mu}(n + \hat{\nu})] + \frac{a^2}{2}[A_{\nu}(n + \hat{\mu}), A_{\nu}(n)] + \mathcal{O}(a^3) \right]. \end{aligned} \quad (3.29)$$

Now the terms that are shifted from the site n can be expanded as:

$$A_{\mu}(n + \hat{\nu}) = A_{\mu}(n) + a\partial_{\nu}A_{\mu}(n) + \mathcal{O}(a^2) \quad (3.30)$$

and with this substitution most terms cancel and one gets:

$$\begin{aligned} P_{\mu\nu}(n) &= \exp [ia^2(\partial_{\mu}A_{\nu}(n) - \partial_{\nu}A_{\mu}(n) + i[A_{\mu}(n), A_{\nu}(n)]) + \mathcal{O}(a^3)] \\ &= \exp [ia^2G_{\mu\nu}(n) + \mathcal{O}(a^3)], \end{aligned} \quad (3.31)$$

where we recovered the gauge field strength tensor using $G_{\mu\nu} = \partial_{\mu}A_{\nu} - \partial_{\nu}A_{\mu} + i[A_{\mu}, A_{\nu}]$. The relation found in eq. (3.31) can be used to build the euclidean lattice action term for the gluons. In particular, to be able to recover the continuum action, we would like a term of the form:

$$S_G[U] = \frac{a^4}{2g^2} \sum_{n \in \Lambda} \sum_{\mu\nu} \text{Tr}(G_{\mu\nu}(n)^2) \xrightarrow{a \rightarrow 0} \frac{1}{4g^2} \int d^4x G_{\mu\nu}(x)^2 = S_G[A]. \quad (3.32)$$

Note that the $1/g_0^2$ term, which could seem strange since it is not in the known continuum Lagrangian, is given by the scaling of the field in eq. (3.24). Up to order $\mathcal{O}(a^2)$ this can be obtained, in terms of the link variables, by:

$$S_G[U] = \frac{2}{g^2} \sum_{n \in \Lambda} \sum_{\mu < \nu} \text{ReTr}(\mathbb{1} - P_{\mu\nu}(n)). \quad (3.33)$$

Higher order corrections to this action can be computed analytically by considering higher orders in the BCH expansion in eq. (3.29) and in the exponential expansion in eq. (3.31).

3.2.4 Lattice Fermions

The discretization of fermions as discussed in section 3.2.1 is incomplete as it leads to unphysical results. It becomes evident if one considers the Fourier transform of the propagator. Let's first rewrite eq. (3.22) in a more compact way, introducing the lattice Dirac propagator $M_{xy}[U]$:

$$S_F[\psi, \bar{\psi}] = a^4 \sum_{n \in \Lambda} \bar{\psi}(x) M_{xy}[U] \psi(y), \quad (3.34)$$

with

$$M_{xy}[U] = \sum_{\mu=1}^4 \gamma_\mu \frac{U_\mu(x) \delta_{x,(y-\hat{\mu})} - U_{-\mu}^\dagger(x) \delta_{x,(y+\hat{\mu})}}{2a} + m \delta_{x,y}. \quad (3.35)$$

In momentum space the propagator becomes:

$$\tilde{M}_{pq} = \delta(p - q) \tilde{M}(p) \quad \text{where} \quad \tilde{M}(p) = m \mathbb{1} + \frac{i}{a} \sum_{\mu=1}^4 \gamma_\mu \sin(p_\mu a). \quad (3.36)$$

In order to calculate the inverse of the propagator in real space we need to invert the one in momentum space and perform an inverse Fourier transform. However, the inverse of the propagator in Fourier space has multiple poles, for example in the case of massless fermions:

$$\tilde{M}(p)^{-1} \Big|_{m=0} = \frac{-ia \sum_\mu \gamma_\mu \sin(p_\mu a)}{\sum_\mu \sin^2(p_\mu a)}. \quad (3.37)$$

The problem vanishes for $a \rightarrow 0$, the continuum case, returning just one fermion type. On the contrary, for non-vanishing lattice spacings multiple fermions, one for each pole, emerge. This is known as the “doubling problem” and the extra fermions are called doublers. The solution, proposed by Wilson [15], is to modify the propagator adding a term that makes the poles in the inverse of the Fourier transformed propagator vanish. In momentum space the Wilson propagator is defined as:

$$\tilde{M}^W(p) = m \mathbb{1} + \frac{i}{a} \sum_{\mu=1}^4 \gamma_\mu \sin(p_\mu a) + \mathbb{1} \frac{1}{a} \sum_{\mu=1}^4 (1 - \cos(p_\mu a)). \quad (3.38)$$

Now, by applying the inverse Fourier transform, one gets for coordinate space:

$$M_{xy}^W[U] = \frac{1}{2a} \sum_{\mu=\pm 1}^{\pm 4} (\mathbb{1} - \gamma_\mu) U_\mu(x) \delta_{x,(y-\hat{\mu})} + \left(m + \frac{4}{a} \right) \delta_{x,y}, \quad (3.39)$$

where the shorthand notation $\gamma_{-\mu} = -\gamma_\mu$ has been introduced.

The final form of the Wilson Fermion Action is:

$$S_F[\psi, \bar{\psi}] = a^4 \sum_{n \in \Lambda} \bar{\psi}(x) M_{xy}^W[U] \psi(y). \quad (3.40)$$

Now that all the needed information about the action and the fields is set, mainly through equations eq. (3.33) and eq. (3.40), the picture of how to discretize QCD from the continuum Minkovskian spacetime to an euclidean spacetime lattice is complete.

3.3 Path Integrals on the Lattice

To express expectation values and correlators on the lattice we employ the path integral formalism. The partition function, as we have seen earlier, is the path integral of the fields over the whole space of the exponential of the action. For the case of QCD the fields are U , ψ and $\bar{\psi}$:

$$Z = \int \mathcal{D}\psi \mathcal{D}\bar{\psi} \mathcal{D}U e^{-S_E[\psi, \bar{\psi}, U]}, \quad (3.41)$$

with the action being the sum of the gluonic and fermionic parts:

$$S_E[\psi, \bar{\psi}, U] = S_G[U] + S_F[\psi, \bar{\psi}, U] = S_G[U] + \bar{\psi} M[U] \psi. \quad (3.42)$$

The immediate simplification is to integrate out the fermion fields. As in the continuum case one can perform the integration of the Grassmann-valued fields.

Grassmann numbers are defined to be anticommutative, meaning that for any two Grassmann numbers a, b the relation $ab = -ba$ holds. A useful general property of these numbers regards the evaluation of multidimensional Gaussian integrals. For set of Grassmann numbers θ_i , their complex conjugates θ_i^* and a Hermitean matrix K one can define the following integral:

$$\begin{aligned} \int \mathcal{D}\Theta^* \mathcal{D}\Theta e^{-\Theta^* K \Theta} &= \left(\prod_i \int d\theta_i^* d\theta_i \right) e^{-\theta_i^* K_{ij} \theta_j} = \left(\prod_i \int d\theta_i^* d\theta_i \right) e^{-\sum_i \theta_i^* k_i \theta_i} \\ &= \prod_i k_i = \det K, \end{aligned} \quad (3.43)$$

A more detailed proof of this relation can be found in [4]. This result is very different from what one would get in the case of real numbers, where the analog integral that of is a multi-variable gaussian, with known solution $(2\pi)^n / \det(K)$.

With the result of eq. (3.43) we can simplify greatly the partition function:

$$Z = \int \mathcal{D}\psi \mathcal{D}\bar{\psi} \mathcal{D}U e^{-S[\psi, \bar{\psi}, U]} = \int \mathcal{D}U e^{-S_G[U]} \det M[U] \quad (3.44)$$

In a similar fashion as in statistical mechanics, the expectation value of an observable on the lattice can be computed, recalling the general form eq. (3.1), as:

$$\langle O \rangle = \frac{1}{Z} \int \mathcal{D}U O[\psi, \bar{\psi}, U] e^{-S_G[U]} \det M[U]. \quad (3.45)$$

To compute fermionic correlators however, another relation of Grassmann integrals is needed:

$$\begin{aligned} \int \mathcal{D}\Theta^* \mathcal{D}\Theta \theta_a^* \theta_b e^{-\Theta^* K \Theta} &= \left(\prod_i \int d\theta_i^* d\theta_i \right) \theta_a^* \theta_b e^{-\theta_i^* K_{ij} \theta_j} = \left(\prod_i \int d\theta_i^* d\theta_i \right) \theta_a^* \theta_b e^{-\sum_i \theta_i^* k_i \theta_i} \\ &= (\det K)(K^{-1})_{ab}, \end{aligned} \quad (3.46)$$

which allows us to write for the correlator of a fermion from a site m , the to a site n on the lattice:

$$\langle O_n(\psi)O_m(\bar{\psi}) \rangle = \frac{1}{Z} \int \mathcal{D}U O_n[\psi, U]O_m[\bar{\psi}, U]e^{-S_G[U]}(\det M[U])(M[U]^{-1})_{mn}. \quad (3.47)$$

This last result is crucial for the numerical evaluation of fermionic observables. For example, for a given “source” $\bar{\psi}(x)$ and a “sink” $\psi(y)$ the correlator between the two can be computed via path integrals, but this requires inverting the fermion propagator matrix.

The above expressions for the partition function and the observables cannot be evaluated or simplified analytically any further, so the usual approach is to approximate the path integral numerically. The main idea is to create an ensemble of field configurations using the probability distribution function:

$$P[U] = \frac{e^{-S_G[U]} \det M}{Z}. \quad (3.48)$$

Given an ensemble $\mathcal{U} = \{U_1, U_2, \dots, U_N\}$ of configurations sampled with the PDF above, the integral in eq. (3.45) can be approximated as:

$$\langle O \rangle = \int \mathcal{D}U P[U]O[\psi, \bar{\psi}, U] \approx \frac{1}{N} \sum_{i=1}^N O(\psi, \bar{\psi}, U_i), \quad (3.49)$$

so that the expectation value of O is the average of the observable measure on the configurations of \mathcal{U} . The interesting, but difficult, problem that we have to face now is the choice of the ensemble. The solution is to use Monte Carlo methods, in particular to sample the space of the configurations using a Markov chain. In chapter 5 a more detailed description of this problem, and a sketch of the implementation that has been developed to address it, will be presented.

3.3.1 Pure Gauge Field Theory

Computing full QCD on the lattice is computationally expensive, mainly due to the integration of fermions via the determinant. From a numerical point of view, the determinant needs to be computed at every step of the Markov chain that is used to evaluate the path-integral. This operation affects the time cost of sampling the configuration space dramatically. A first approach is to neglect the determinant completely, considering it constant. This, on a deeper level, effectively means removing dynamical fermions, freezing them to the lattice sites. This approximation is usually referred to as “quenched QCD”, or QQCD.

The properties of this theory, that is then reduced to a simple Yang-Mills theory with zero fermion flavors, are still interesting to study and have historically played a very important role, being the only accessible simulation until sufficient computing power became available. For example, from eq. (2.20) one can see that the theory presents confinement and asymptotic freedom.

3.3.2 Observables

On the lattice, given the transformation eq. (3.21), any product of link variables that starts and ends at the same site, a closed loop, is gauge invariant. The average values of these objects over

the whole lattice can be linked to physical observables, for example the field tensor. In a more general form any observable $L[U]$ of the type:

$$L[U] = \text{Tr} \left[\prod_{(n,\mu) \in \mathcal{P}} U_\mu(n) \right], \quad (3.50)$$

where \mathcal{P} is a closed path of links on the lattice, is a gauge invariant object.

Plaquette

The simplest observable, which we have already encountered upon defining the Wilson action in eq. (3.33), is the plaquette. This is the minimal closed loop on the lattice it is related to the action. For each lattice site there are 6 possible plaquettes to be computed, given all the combinations of euclidean indices. One can consider the average value of the plaquette on a lattice field configuration as an observable related to the value of the action of the field, and consequently to the coupling constant. A proper definition of the observable, computed on a lattice Λ is:

$$P[U] = \frac{1}{6|\Lambda|} \sum_{n \in \Lambda} \sum_{\mu < \nu} P_{\mu\nu}(n), \quad (3.51)$$

where $P_{\mu\nu}(n)$ is the one defined in eq. (3.27) and $|\Lambda|$ is the notation that we introduce for the total number of lattice sites.

Energy Density

The energy of the field is proportional to the square of the field tensor, so the energy density is in particular:

$$E[U] = -\frac{1}{4|\Lambda|} \sum_{n \in \Lambda} \sum_{\mu < \nu} \text{ReTr}(G_{\mu\nu}(n) G_{\mu\nu}(n)). \quad (3.52)$$

In order to estimate this quantity, one has to compute the field tensor at every lattice site, square it and sum over the whole space. It is also usually normalized by the lattice volume (the number of sites) to get the density. The simplest definition of the field tensor is $G_{\mu\nu}^{(\text{plaq})}(n) = \mathbb{1} - P_{\mu\nu}(n)$ but this is not very accurate. A more symmetric definition can be obtained by the “clover leaf” field tensor, that is the average of all the plaquettes of a same plane $\mu\nu$ that start from a given lattice site taken all with the same orientation. To better understand the clover term, in figure fig. 3.3 a representation is given.

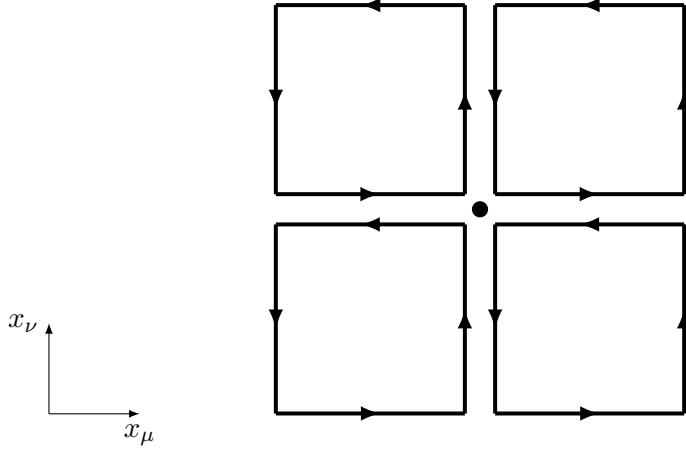


Figure 3.3: Schematic representation of the symmetric definition of the clover term $G_{\mu\nu}^{(clover)}(n)$ on the lattice. The field strength tensor for a plane $\mu\nu$ at a lattice site is given by the average of all the for plaquettes that start from a site n , all taken with the same orientation.

In terms of link variables this is equal to:

$$G_{\mu\nu}^{(clover)}(n) = \frac{1}{4} \left[U_\mu(n)U_\nu(n + \hat{\mu})U_\mu^\dagger(n + \hat{\nu})U_\nu^\dagger(n) \right. \\ \left. + U_\nu(n)U_\mu^\dagger(n + \hat{\nu})U_\nu^\dagger(n - \hat{\mu})U_\mu(n - \hat{\mu}) \right. \\ \left. + U_\mu^\dagger(n - \hat{\mu})U_\nu^\dagger(n - \hat{\mu} - \hat{\nu})U_\mu(n - \hat{\mu} - \hat{\nu})U_\nu(n - \hat{\nu}) \right. \\ \left. + U_\nu^\dagger(n - \hat{\nu})U_\mu(n - \hat{\nu})U_\nu(n + \hat{\mu} - \hat{\nu})U_\mu^\dagger(n) \right]. \quad (3.53)$$

Topological Charge

The gauge field in QCD exhibits particular topological properties that are believed to have important physical implications [19, 20]. The topological charge is an integer quantum number of the field in the continuum, similar to the concept of winding-number. In general fields with different topological charge are separated, but there exist non-zero instanton modes [21] that allow tunneling between them. On the lattice certain definitions can be used to reproduce the continuum properties, especially for the so called ‘‘topological susceptibility’’, that is the second moment of the distribution of the topological charge, which seems to be independent from the definitions of the base observable [22, 23].

The topological charge is the integral over all spacetime of the topological charge density

$$Q = \int d^4x q(x), \quad (3.54)$$

where

$$q(x) = \frac{1}{32\pi^2} \text{Tr}(G_{\mu\nu}(x)\tilde{G}^{\mu\nu}(x)). \quad (3.55)$$

Here with $\tilde{G}^{\mu\nu}$ the dual of the gauge field is intended and its definition is:

$$\tilde{G}^{\mu\nu}(x) = \frac{1}{2}\epsilon^{\mu\nu\rho\sigma}G^{\rho\sigma}(x), \quad (3.56)$$

with $\epsilon^{\mu\nu\rho\sigma}$ being the anti-symmetric Levi-Civita symbol. We can now see the link between this observable and the θ -term on the QCD Lagrangian we mentioned in section 2.1.2.

The topological charge can be estimated on the lattice, in the simplest way, by using the same clover definition of the field strength tensor we used before for the energy density:

$$Q[U] = \frac{1}{64\pi^2} \sum_{n \in \Lambda} \sum_{\mu < \nu} \epsilon^{\mu\nu\rho\sigma} \text{Tr}[G_{\mu\nu}^{(\text{clover})}(n) G_{\rho\sigma}^{(\text{clover})}(n)]. \quad (3.57)$$

As the lattice spacing is reduced, approaching the continuum, the topological sectors get more and more separated, preventing tunneling between them. This makes the Markov chain used to generate the ensemble less efficient in terms of growing autocorrelation times of the observables, as will be shown in section 6.2.

A derived quantity of great interest is the topological susceptibility χ , that is proportional to the expectation value of the square of the topological charge:

$$\chi = \frac{(\hbar c)^4}{a^4 |\Lambda|} \langle Q^2 \rangle. \quad (3.58)$$

This observable is particularly important for the properties of instantons on the lattice and its value can be related via the Witten-Veneziano formula [19] to the mass of the η' meson.

3.4 Modern Lattice QCD Calculations

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The state of the art calculations in Lattice QCD involve improved fermion and gluon actions. For the gluon sector the improvements are based on finding linear combinations of gauge invariant loops, such as 2×1 , 2×2 and 3×1 rectangles to systematically remove the higher order in the right-hand side of eq. (3.31). For fermions the improvements consist on actions that preserve certain symmetries more than others, like chiral fermion actions or staggered fermions. On the algorithm side, the currently most used method to generate gauge field configuration is the Hybrid Monte Carlo method, which combines stochastic sampling with hamiltonian dynamic updates in the gauge field space. Other interesting algorithmic problems regard the calculation of fermion determinants, which typically involve very large sparse matrix diagonalization.

A fundamental concept in Lattice calculations is to recover the continuum limit in a controlled manner at the end of the analysis. The error sources that come into play when discretizing spacetime on a fixed lattice are:

- Finite Volume Effects: caused by the non infinite domain of the simulation. Usually periodic boundary conditions are implemented to simulate infinite space, but one should always check if the total lattice volume is large enough, especially when dealing with large systems such as multiple baryons.

- State Isolation and Signal to Noise: when computing correlators between two points on the lattice it is possible to extract energy states of hadrons by looking at the exponential decay of the correlator in euclidean time. However this problem is largely affect by random noise making it hard to extract even the ground state most of the times; excited states are rarely considered.
- Chiral Limit: by which is intended the limit for which the masses of the quarks, and consequently of the computed hadrons, on the lattice approach the physical masses. One might question why this is even the case and that calculations should be performed at the physical masses only. However, the fermion propagator matrix becomes more and more sparse the lower the mass of the quark is, making the numerical algorithms that should diagonalize it slower to converge. The usual approach is to perform calculations on a set of pion masses (this is the usual reference for this problem) and afterwards the limit for $m_\pi^{(lat)} \rightarrow m_\pi^{(phys)}$ is taken to reproduce the physical observables.
- Continuum Limit: perhaps the most obvious thing to do to improve Lattice QCD calculations is to take finer and finer lattice spacings, to approach the continuum case. It should be noted however that this procedure has two downsides. First the lattice spacing in physical units is fixed by the quark masses, and as we previously mentioned, this affects the feasibility of the numerical simulation itself. Secondly, as the lattice approaches the continuum case the autocorrelation of observables in a given ensemble increases rapidly. This behavior, known as “critical slow-down” depends on the observable itself, some are more affected than others. The integrated autocorrelation time for the topological charge for example is believed to have either power-law with a large exponent or even an exponential relation to the lattice spacing.

As one can notice, Lattice QCD is still an open field of research in many aspects: the algorithmic/numerical, the extraction and improvement of the measurement of physical quantities and the theoretical model itself.

Chapter 4

Advanced Topics in Lattice QCD

In this sections we will present some problems and methods that are used in Lattice QCD and in particular for pure Yang-Mills case, the fermionless theory. A four-loop corrected expression for the β -function of QCD will be discussed as well, together with a tentative method of determining the scale parameter Λ from lattice calculations.

4.1 Scale Setting in the Quenched Approximation

On the lattice quantities are defined to be dimensionless and in order to link them with physical observables they need to be fixed by the correct power of the lattice spacing a to reproduce the physical units. In the quenched approximation, the pure Yang-Mills case, the fundamental question is then how to connect the lattice spacing to the coupling used in the gluonic action. The general strategy is to compute on the lattice a quantity that is known to be finite in the continuum limit.

In 1993 R. Sommer proposed to use the static quark potential as a reference in the case of SU(2) theory [24]. Later in another work [25] a length scale r_0 was suggested for SU(3) Yang-Mills theory to be set as the distance such that

$$r_0^2 F(r_0) = 1.65, \quad (4.1)$$

with $F(r_0)$ being the force between two static color charges. In [25] also a parametrization of the ratio a/r_0 is provided, in terms of the coefficients of the renormalization group equation. To leading order one finds:

$$\frac{a}{r_0} \propto e^{-\frac{\beta}{12b_0}}, \quad (4.2)$$

with inverse coupling $\beta = 6/g_0^2$ and b_0 the one loop coefficient of the perturbative expansion of the β function, see section 2.2.1 , $b_0 = 11/(4\pi)^2$. Using the ansatz:

$$\ln \left(\frac{a}{r_0} \right) = \sum_{k=0}^p a_k (\beta - 6)^k, \quad (4.3)$$

a parametrization up to $p = 3$ has been fitted to data by Guagnelli et al. [25]. The resulting interpolating function, plotted against the data in fig. 4.1, is:

$$\ln\left(\frac{a}{r_0}\right) = -1.6805 - 1.7139(\beta - 6) + 0.8155(\beta - 6)^2 - 0.6667(\beta - 6)^3. \quad (4.4)$$

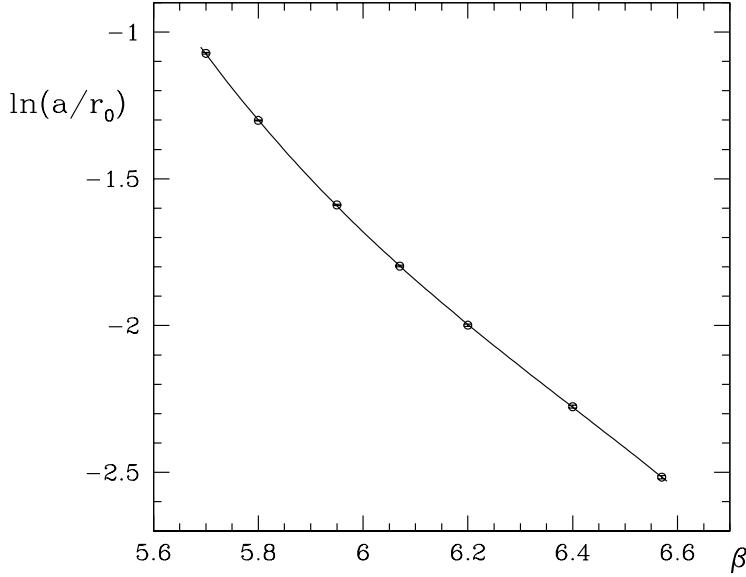


Figure 4.1: Data and interpolating function for $\ln(a/r_0)$ as found in [25].

This scale fixing procedure requires a proper estimate of the length scale from measurements of the static quark potential for various quark sources, which is an expensive calculation. For some simpler calculations, it is customary to set the length scale to a fixed value, namely $r_0 = 0.5$ fm. This value commonly called the Sommer scale parameter, is then used in eq. (4.4) to implicitly extract a .

More recently another procedure to fix the energy scale has been proposed, based on the Gradient Flow Method, which will be discussed in the following section. The key idea is again to choose an observable that is finite in the continuum limit and use it to set a reference scale on the lattice.

4.2 The Gradient Flow Method

The Gradient Flow Method, also known as the Wilson flow in the case of QCD, is a method for studying non-linear properties of the gauge fields. The key idea is to see how the theory evolves as it becomes less local, by driving it to the stationary points of the action, applying a “diffusion-like” differential equation to the field in a fictitious dimension called “flow time”. In more precise terms we apply a local gauge-covariant diffusion equation to the fields.

For the SU(3) gauge field case most of the theoretical foundation was set by M.Lüscher in [1, 26]

and for QCD in [27]. Starting from a field $A_\mu(x)$ the characterizing equations are:

$$\begin{aligned}\partial_{t_f} B_\mu &= D_\mu G_{\mu\nu} \\ G_{\mu\nu} &= \partial_\mu B_\nu - \partial_\nu B_\mu + [B_\mu, B_\nu] \\ D_\mu &= \partial_\mu + [B_\mu, \cdot],\end{aligned}\tag{4.5}$$

where the $B(t_f, x)$ field is the flowed version of the original field at flow time t_f . This is imposed by fixing:

$$B_\mu|_{t_f=0} = A_\mu.\tag{4.6}$$

The second equation of eq. (4.5) is the generalization of the field strength tensor of the field B ; the right-hand side of the first equation is then proportional to the gradient of the action. The effect of the gradient flow equations is to move the field along the direction of the minimum action, eventually reaching the stationary points. In a discretized theory taking the derivatives in the above equations is not a trivial procedure. However we can use instead the intuitive idea of the flow equations, that is to use the steepest descent method on the action of a field. We then introduce the flowed lattice gauge field $V_{t_f}(x, \mu)$ as the flow time evolved version of a field $U(x, \mu)$. Our flow equation then becomes:

$$\partial_{t_f} V_{t_f}(x, \mu) = -g_0^2[\partial_{x,\mu} S_G(V_{t_f})]V_{t_f}(x, \mu),\tag{4.7}$$

where S_G is the Wilson action, which was discussed in section 3.2.3, defined as:

$$S_G[V_{t_f}] = \frac{2}{g^2} \sum_{n \in \Lambda} \sum_{\mu < \nu} \text{ReTr}(\mathbb{1} - P_{\mu\nu}(n)).\tag{4.8}$$

generalized to flowed fields. The sum is performed over all sites of a lattice Λ and the 1×1 Wilson loop $P_{\mu\nu}(n)$ (the plaquette) is taken to from the action. The generalization is to the flowed field is straightforward. The definition is unchanged and the action is simply evaluated on the field V_{t_f} .

The flow equations for the fermion fields, although they have not been used in this work, are:

$$\begin{aligned}\partial_{t_f} \chi &= \Delta \chi, & \partial_{t_f} \bar{\chi} &= \bar{\chi} \overleftarrow{\Delta} \\ \Delta &= D_\mu D_\mu, & D_\mu &= \partial_\mu + b_\mu,\end{aligned}\tag{4.9}$$

note that Δ is the covariant Laplacian in this notation. The initial conditions are naturally

$$\chi|_{t_f=0} = \psi, \quad \bar{\chi}|_{t_f=0} = \bar{\psi}.\tag{4.10}$$

4.3 Perturbative Analysis of the Wilson Flow

An important thing to consider when applying the Wilson flow to a field is the renormalization of the observables: one has to check that expectation values of observables at non-zero flow time are renormalized quantities. Following the calculations performed in [1], we will consider the energy as our base observable.

First, we note that the flow equation is invariant under flow time independent gauge transformations, this prevents a detailed study of the renormalization. We consider a modified version of the flow equation by adding one term:

$$\partial_{t_f} B_\mu = D_\mu G_{\mu\nu} + \lambda D_\mu \partial_\nu B_\nu. \quad (4.11)$$

The original case is obtained again by setting $\lambda = 0$ and considering:

$$B_\mu = \Lambda B_\mu|_{\lambda=0} \Lambda^{-1} + \Lambda \partial_\mu \Lambda^{-1} \quad (4.12)$$

and with $\Lambda(t, x)$ now being a flow time dependent gauge transformation, set by:

$$\partial_{t_f} \Lambda_\mu = -\lambda \partial_\nu B_\nu \Lambda \quad \text{with } \Lambda|_{t_f=0} = 1. \quad (4.13)$$

The modified flow equation allows to asymptotically expand the field B_μ in powers of the bare coupling constant g_0 . In particular, to leading order and with the choice of $\lambda = 1$, it was shown in [26] that eq. (4.11) reduces to:

$$\partial_{t_f} B_\mu - \partial_\nu \partial_\nu B_\mu = 0, \quad (4.14)$$

which has a solution, in terms of the initial field $A_\mu(x)$ of the type:

$$B_\mu(t_f, x) = \int d^4y K_{t_f}(x-y) A_\mu(y). \quad (4.15)$$

with $K_{t_f}(z)$ being the heat kernel of the flow equation:

$$K_{t_f}(z) = \frac{e^{-z^2/4t_f}}{(4\pi t_f)^2} \quad (4.16)$$

This result shows that the flow equations, as a first approximation, are a smearing operation. The right-hand side can be seen as the integral average of the original field weighted with a gaussian distribution of variance $2t_f$ and mean-square-radius of $\sqrt{8t_f}$.

The energy E as defined in eq. (3.52) is a gauge invariant object. One can observe that the modified flow equation is invariant under gauge transformations, hence the energy, as any other gauge invariant observable, will preserve its invariance. Using the modified flow equation, the asymptotic expansion of the flowed field B_μ and the definition of the energy $\langle E \rangle = -\frac{1}{4} G_{\mu\nu} G^{\mu\nu}$ it is possible to write the expectation value of the energy as a function of the flow time and of the renormalized coupling [1]:

$$\langle E \rangle = \frac{3(N^2 - 1)g_0^2}{128\pi^2 t_f^2} [1 + \bar{c}_1 g_0^2 + \mathcal{O}(g_0^4)], \quad (4.17)$$

with:

$$\bar{c}_1 = \frac{1}{16\pi^2} \left[N \left(\frac{11}{3}L + \frac{52}{9} - 3\ln 3 \right) - N_f \left(\frac{2}{3}L + \frac{4}{9} - \frac{4}{3}\ln 2 \right) \right]. \quad (4.18)$$

The equation is for general N , the gauge group dimension, and number of quark flavors N_f . The coefficient \bar{c}_1 is in terms of a scale $L = \ln(8\mu^2 t_f) + \gamma_E$ where μ is the renormalization

energy scale and γ_E Euler's constant. From this last relation we can set a flow energy scale as $q = 1/\sqrt{8t_f}$ and give a definition of eq. (4.17) in terms of the running coupling $\alpha_s(q) = \frac{g_0^2}{4\pi}$:

$$\langle E \rangle = \frac{3(N^2 - 1)}{32\pi t_f^2} \alpha_s(q) [1 + k_1 \alpha_s(q) + \mathcal{O}(\alpha_s^2)] \quad (4.19)$$

where the coefficient k_1 is now:

$$k_1 = \frac{1}{4\pi} \left[N \left(\frac{11}{3} \gamma_E + \frac{52}{9} - 3 \ln 3 \right) - N_f \left(\frac{2}{3} \gamma_E + \frac{4}{9} - \frac{4}{3} \ln 2 \right) \right] \quad (4.20)$$

In the case of SU(3) symmetry group we have:

$$\langle E \rangle = \frac{3}{4\pi t_f^2} \alpha_s(q) [1 + k_1 \alpha_s(q) + \mathcal{O}(\alpha_s^2)], \quad k_1 = 1.0978 + 0.0075 \times N_f \quad (4.21)$$

The perturbative expansion is only valid for large values of q , that is when the flow time is small. We also can notice that $\sqrt{8t_f}$ is the inverse of an energy, or a length. This gives us a further intuitive picture of what the gradient flow equations do, that is to smear the field over a hypersphere of radius $\sqrt{8t_f}$. The gradient flow is particularly interesting to study quantities that are divergent on the lattice because of discretization effects, like the topological susceptibility, which become smooth at non-zero t_f . In fig. 4.2 an example of the smearing effect of the gradient flow on the topological charge is shown.

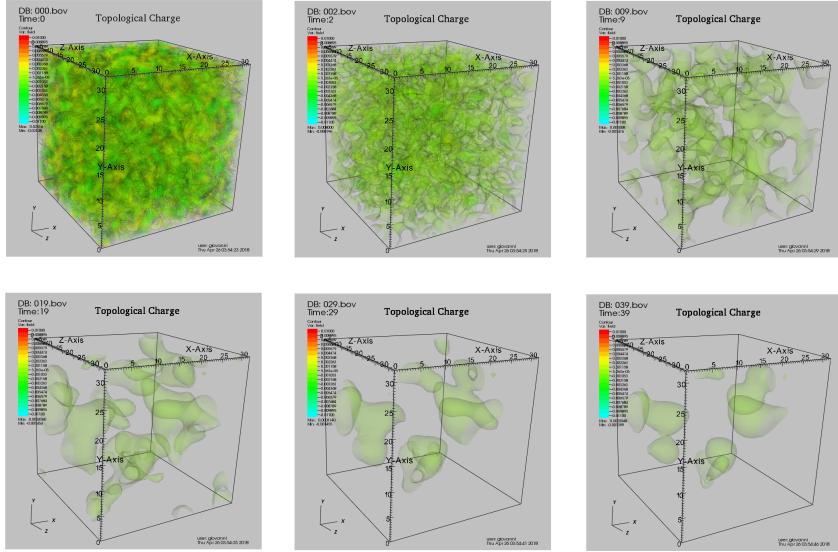


Figure 4.2: Topological charge computed ad one euclidean time of a lattice of size $32^3 \times 64$ with lattice spacing 0.06793 fm. The flow different plots are for flow times $\sqrt{8t_f} = 0, 0.14, 0.30$ fm on the first row and $\sqrt{8t_f} = 0.43, 0.52, 0.60$ fm on the second row. The visualization has been performed using a custom `python` interface (developed together with H.M.Vege), called “LatViz”, to the open source tool VisIt by LLNL [28].

The topological charge at $\sqrt{8t_f} = 0$ fm (the non flowed field configuration) is highly affected by large fluctuations over the lattice sites. This for example affects the topological susceptibility

causing it to be divergent. One can notice that the action of the flow equation is to remove such low distance fluctuations that at $\sqrt{8t_f} = 0.30$ fm are almost completely gone. For larger smearing radii the emergence of well separated topological sectors becomes visible.

4.3.1 Scale Fixing with the Gradient Flow

Because of the smearing properties of the flow equation the fields are driven towards the minima of the action, which are in general independent from the lattice spacing. For sufficiently large flow times, one can therefore decide to use the value of the dimensionless quantity $t_f^2 \langle E \rangle$ to fix the scale of the lattice. The choice of this quantity is motivated by the fact that $\langle E \rangle = -\frac{1}{4} G_{\mu\nu} G^{\mu\nu}$ is proportional to the Wilson gauge action eq. (3.33). The gradient flow minimizes the action, so independently of the lattice spacing, all field configurations will converge to the same value. It has been proposed by Lüscher that a value of $t_f^2 \langle E \rangle = 0.3$ is large enough to be used as a reference scale. In fig. 4.3 we can indeed observe that the value is constant as a function of the

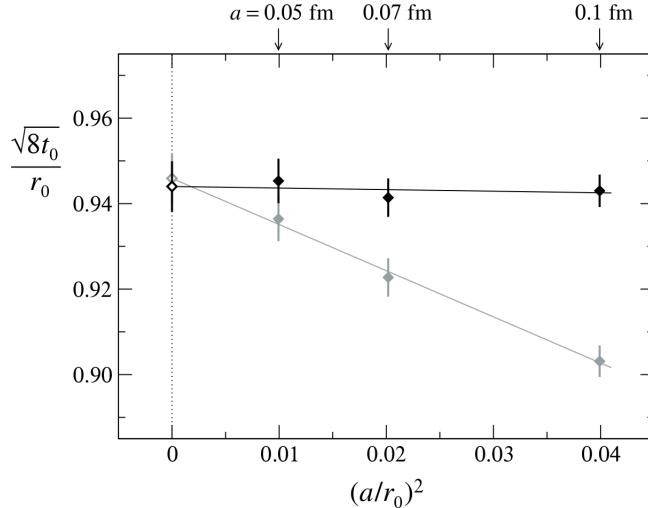


Figure 4.3: Plot of $\sqrt{8t_0}/r_0$ where t_0 is the value such that $t_0^2 \langle E \rangle = 0.3$ for different lattice spacings taken from [29]. The dimension less ratio $\sqrt{8t_0}/r_0$ shows the equivalence between the Sommer scale r_0 and the gradient flow reference scale t_0 . The continuum limit extrapolation is performed using $(a_0/r_0)^2$ as the dimensionless rescaling quantity. The two data series, better explained in the article, represent two different definitions of the gauge field strength tensor and consequently of the energy density.

lattice spacing (computed instead using the Sommer parameter). In section 8.1 this method will be applied and discussed more thoroughly.

4.4 Estimating the Λ Scale Parameter

Finally all the theoretical concepts and tools have been defined to state the main purpose of this work in a formal way. The goal is to use the perturbative expression of the energy as a

function of the flow time and of the running coupling, eq. (4.17), to estimate the scale Λ of the theory.

Since the algorithm for generating ensembles of gauge field configurations that has been developed in this work only considers pure gauge theory, we will restrict the analysis to Λ_{YM} . This is the scale parameter when one sets $N_f = 0$, no fermion flavors. In a future work the more interesting determination of Λ_{QCD} could be performed the same way.

On the lattice the quantity $t_f^2 \langle E \rangle$ can be computed easily, this leaves only the coupling on the right hand side of eq. (4.21) as the unknown variable. One can then choose an order for the expansion of the Renormalization Group Equation:

$$q^2 \frac{d\alpha(q)}{dq^2} = \beta(\alpha) = -(b_0\alpha^2 + b_1\alpha^3 + \dots) \quad (4.22)$$

When solving the equation for $\alpha(q)$ there is some freedom on the choice of the renormalization point, as we have seen in eq. (2.21), such that the equation in the end can be simplified by the introduction of a scale parameter.

Using the data computed on the lattice, we will first construct the continuum limit estimate for $t_f^2 \langle E \rangle$ in order to account for discretization effects. Then an estimate the scale will be found by finding an appropriate procedure to match the lattice results to the perturbation theory expansion.

4.4.1 4-loop Corrected Running Coupling

The equation eq. (4.21) needs to be have a functional form for $\alpha(q)$ as input, here we report an expression that is including diagrams up to 4-loops in the calculations of the RGE, parametrized by 4 coefficients. The equation was taken by the Particle Data Group review of QCD[30]:

$$\begin{aligned} \alpha(q) = & \frac{1}{b_0 t} \left[1 - \frac{b_1 \ln t}{b_0 t} + \frac{b_1^2 (\ln^2 t - \ln t - 1) + b_0 b_2}{b_0^4 t^2} \right. \\ & \left. - \frac{b_1^3 (\ln^3 t - \frac{5}{2} \ln^2 t - 2 \ln t + \frac{1}{2}) + 3b_0 b_1 b_2 \ln t - \frac{1}{2} b_0^2 b_3}{b_0^6 t^3} \right] \end{aligned} \quad (4.23)$$

here the convenient notation of $t \equiv \ln \frac{q^2}{\Lambda^2}$ has been introduced. The coefficients b_0, b_1, b_2 and b_3 can be computed analytically for a general theory on SU(3) with N_f quark flavors. While b_0 and b_1 have known exact values, the other two parameters are renormalization scheme dependent,

and we chose the ones computed in the \overline{MS} scheme found in [31]. The coefficients are then:

$$\begin{aligned} b_0 &= \frac{1}{(4\pi)} \left[11 - \frac{2}{3} N_f \right] \\ b_1 &= \frac{1}{(4\pi)^2} \left[102 - \frac{38}{3} N_f \right] \\ b_2 &= \frac{1}{(4\pi)^3} \left[\frac{2857}{2} - \frac{5033}{18} N_f + \frac{325}{54} N_f^2 \right] \\ b_3 &= \frac{1}{(4\pi)^4} \left[\left(\frac{149753}{6} + 3564\zeta_3 \right) - \left(\frac{1078361}{162} + \frac{6508}{27}\zeta_3 \right) N_f \right. \\ &\quad \left. + \left(\frac{50065}{162} + \frac{6472}{81}\zeta_3 \right) N_f^2 + \frac{1093}{729} N_f^3 \right] \end{aligned} \tag{4.24}$$

or in a numerical form, substituting also the Riemann zeta-function value $\zeta_3 = 1.202056903\dots$ one gets the more handy expression

$$\begin{aligned} b_0 &\approx \frac{1}{(4\pi)} (11 - 0.66667 N_f) \\ b_1 &\approx \frac{1}{(4\pi)^2} (102 - 12.66667 N_f) \\ b_2 &\approx \frac{1}{(4\pi)^3} (1428.50 - 279.611 N_f + 6.01852 N_f^2) \\ b_3 &\approx \frac{1}{(4\pi)^4} (29243.0 - 6946.30 N_f + 405.089 N_f^2 + 1.49931 N_f^3) \end{aligned} \tag{4.25}$$

with the aid of these coefficients the only unknown variable in eq. (4.17), using eq. (4.23) for $\alpha_S(q)$, is the scale parameter Λ .

Part II

Implementation

Chapter 5

Designing a Lattice $SU(3)$ Yang-Mills Theory Code

One of the major focuses of this work has been a completely new implementation of a program to generate and analyse $SU(3)$ lattice gauge field configurations. As it is common practice in Lattice QCD numerical implementations, the program is split in three parts:

1. *Generation of Gauge Fields*: in this case it is done using an implementation of the Metropolis algorithm using the gluonic Wilson action;
2. *Computation of Lattice Observables*: this includes applying the gradient flow to the gauge fields, as well as computing the energy density and the topological charge at every flow time; in other applications the computation of quark propagators and two- or three-point correlators are performed;
3. *Computation of Derived Quantities*: mainly post analysis, computation of secondary observables and uncertainties. Resampling and autocorrelation analysis are crucial to Lattice QCD given the typically small size of the ensembles.

In this chapter the main features of the program that was developed to deal with the first two steps, which are the most interesting, will be presented. The third step, the post analysis, which is a simpler calculation, did not require a very sophisticated, or challenging, program.

The problems are numerically very challenging and the calculations had to be performed on computing clusters. The programming language of choice was `C++` because of its high efficiency, high abstraction capabilities and for the easiness of the `MPI` integration (a standard HPC parallelization library). The analysis of data has been performed using `python` and in particular relying heavily on its standard data science packages such as `numpy` and `pandas`.

5.1 Generating Pure Gauge Fields

The task of generating lattice field configurations is extremely demanding in terms of computational requirements. The case of QCD is much more demanding than that of a pure gauge theory, but overall the latter calculation is still challenging. The main, and perhaps overwhelmingly simple, reason for this problem is the dimensionality. Dealing with a discretized space-time lattice, things tend to scale with powers of 2^4 , a trivial example is cutting in half the lattice spacing keeping a fixed total volume: this requires 16 more points in the global lattice.

To get a better feeling of the algorithm we first have to look at what the basic object of the program is: the lattice. On a computer the number of double precision floating point numbers to be stored for a field configuration is given by

$$\underbrace{N^3}_{\text{spatial dimension}} \times \underbrace{N_t}_{\text{time dimension}} \times \underbrace{4}_{\text{links per site}} \times \underbrace{9}_{\text{SU(3) matrix size}} \times \underbrace{2}_{\text{real and imaginary part}} \quad (5.1)$$

For example, if we choose $N = 48$, $N_t = 96$ (which is the largest system considered in this work) the resulting configuration is 6115295232 *bytes* large, that is 5.7 *GBytes*. The incredibly large size of the data to be handled limits greatly the possibility of simulating larger systems. The basic element at each lattice site is a set of 4 different $SU(3)$ matrices, one for each dimension, these are the links variables described in section 3.2. Since the links with negative orientation are related to the ones with positive orientation by the adjoint operation, only the positively oriented ones are stored in memory.

5.1.1 The Metropolis Algorithm

The main algorithm that has been used to generate an ensemble of gauge field configurations is the Metropolis Algorithm. It is a widely popular Markov Chain Monte Carlo Method to generate a sequence of random samples from a probability distribution [32][33].

In general, a Markov Chain is a sequence of randomly chosen variables X_1, X_2, \dots, X_t , in our case a lattice field configuration, with the Markov Property, that is: the probability of the step $t + 1$ depends only on the variable X_t :

$$P(X_{t+1} = x | X_1 = x_1, X_2 = x_2, \dots, X_t = x_t) = P(X_{t+1} = x | X_t = x_t). \quad (5.2)$$

In the case where the probabilities of moving from a state $X_i = i$ to a state $X_j = j$ is not known, the transition probability from the two states, $W(i \rightarrow j)$ can be split into two contributions: the probability $T(i \rightarrow j)$ for making the transition to state j being in state i and the the probability of accepting this transition $A(i \rightarrow j)$.

$$W(i \rightarrow j) = A(i \rightarrow j)T(i \rightarrow j). \quad (5.3)$$

If a Probability Distribution Function (PDF) is known for the process, we can label the probability of a given state at a fixed time $w_i(t)$. The transition probability to a state j at time $t + 1$ is the sum of the transition probabilities of moving to state j from state i plus the probability

of being in state i at time t and rejecting to move to any other state:

$$\begin{aligned} w_j(t+1) &= \sum_i [w_i(t)T(i \rightarrow j)A(i \rightarrow j) + w_j(t)T(j \rightarrow i)(1 - A(j \rightarrow i))] \\ &= w_j(t) + \sum_i [w_i(t)T(i \rightarrow j)A(i \rightarrow j) - w_j(t)T(j \rightarrow i)A(j \rightarrow i)]. \end{aligned} \quad (5.4)$$

For large times t , when equilibrium is reached, we require that $w_j(t+1) = w_j(t) = w_j$, and thus have:

$$\sum_i w_i T(i \rightarrow j) A(i \rightarrow j) = \sum_i w_j T(j \rightarrow i) A(j \rightarrow i). \quad (5.5)$$

Now, considering that the transition probability from a state j to all other states must be normalized $\sum_i W(j \rightarrow i) = \sum_i T(j \rightarrow i)A(j \rightarrow i) = 1$ we get:

$$\sum_i w_i T(i \rightarrow j) A(i \rightarrow j) = w_j. \quad (5.6)$$

At this point, the further constrain of Detailed Balance [34] is introduced, that is:

$$w_i W(i \rightarrow j) = w_j W(j \rightarrow i). \quad (5.7)$$

At equilibrium we then have:

$$\frac{w_i}{w_j} = \frac{W(i \rightarrow j)}{W(j \rightarrow i)} = \frac{T(i \rightarrow j)A(i \rightarrow j)}{T(j \rightarrow i)A(j \rightarrow i)}. \quad (5.8)$$

Making the approximation that the transition probability between states is the same, $T(i \rightarrow j) = T(j \rightarrow i)$, the brute-force approach, we are left with the acceptance ratio of a move to a new state to be the ratio of the PDFs.

In the case of pure gauge fields, the probability distribution is:

$$P(U) = \frac{e^{S_G[U]}}{Z} = \frac{e^{S_G[U]}}{\int \mathcal{D} U e^{S_G[U]}}, \quad (5.9)$$

which is the exponential of the gluon Wilson Action over the path integral of the same quantity over all space (the partition function). Luckily, we only need ratios of PDFs so the partition functions is not to be computed. We can rewrite any expectation value of an operator on the lattice as:

$$\langle O \rangle = \frac{\int \mathcal{D} U O[U] e^{S_G[U]}}{\int \mathcal{D} U e^{S_G[U]}} = \int \mathcal{D} U P(U) O[U], \quad (5.10)$$

which numerically becomes:

$$\langle O \rangle \approx \frac{1}{N} \sum_{i=1}^N O(U_i), \quad (5.11)$$

where, as described in section section 3.3, U_i are random configurations of a set chosen with the PDF of eq. (5.9).

Noting that, since we only care about ratios in the PDF, which turn out to be exponentials of the action difference between two configurations, a tentative algorithm can be written as:

Algorithm 1 Metropolis Algorithm

```

configuration ← initial configuration
for  $i < MonteCarloCycles$  do
    for  $j < N_{corr}$  do
        for  $U$  in Lattice do
            newConfiguration ← randomMove( $U$ ) + configuration
             $\Delta S$  ← action(newConfiguration) – action(configuration)
            if  $\Delta S > \text{random}(0, 1)$  then
                configuration ← newConfiguration
            end if
        end for
    end for
    saveConfiguration(configuration)
end for

```

The actual implementation of this algorithm on discretized space-time is however not as trivial as it seems. Firstly, a definition of what is a “random move” in configuration space is needed, secondly an efficient way of computing the action difference is to be found.

The ensemble, on which observables can be computed, is created by saving to disk, in the form of a simple binary file containing all the data, one configuration every N_C Monte Carlo updates. We need this in order to apply the gradient flow afterwards to the configurations and compute the observables we want. The choice of N_C turned out to be crucial for the autocorrelation of some observables, in particular the topological charge, see section 6.2.

5.1.2 Sampling the Configuration Space

To use Metropolis’ algorithm we need to define a random walk in the space of gauge field configurations. In particular, we will use the Wilson Action:

$$S_G[U] = \frac{\beta}{3} \sum_{n \in \Lambda} \sum_{\mu < \nu} \text{ReTr}(\mathbb{1} - P_{\mu\nu}(n)), \quad (5.12)$$

which is defined on plaquettes $P_{\mu\nu}(n)$ and contains the parameter $\beta = 6/g_0^2$ usually referred to as the inverse coupling. A random move in the configuration space can be seen as a random change of a single link of the lattice Λ . This change must be a unitary transformation in the $SU(3)$ group, but not a completely random one as it should have a strong real diagonal component, making it close to an identity matrix and thus producing a small change on the link.

In order to generate such transformation we use three random $SU(2)$ matrices that are as well “close to unity”. A recipe for generating these matrices is found in [16]. We consider three 2×2 unitary matrices R_2, S_2 and T_2 :

$$R_2 = \begin{pmatrix} r_{11} & r_{12} \\ r_{21} & r_{22} \end{pmatrix}, \quad S_2 = \begin{pmatrix} s_{11} & s_{12} \\ s_{21} & s_{22} \end{pmatrix}, \quad T_2 = \begin{pmatrix} t_{11} & t_{12} \\ t_{21} & t_{22} \end{pmatrix}, \quad (5.13)$$

The elements of the matrix are to be random, but with a controlled spread around the identity matrix. One starts by selecting four random numbers r_0 , \mathbf{r} (the latter is a three component vector) between $(-\frac{1}{2}, \frac{1}{2})$. Then a “spread parameter” ϵ is introduced, which controls how much to weight the off-diagonal terms. The random variables are scaled according to:

$$x_0 = \text{sign}(r_0)\sqrt{1 - \epsilon^2} \quad \mathbf{x}_\parallel = \epsilon \frac{\mathbf{r}}{|\mathbf{r}|}, \quad (5.14)$$

and these four coefficients, together with the generators of the $SU(2)$ group (the Pauli matrices) are used to build the element of the group:

$$U = x_0 \mathbb{1} + i \mathbf{x} \cdot \boldsymbol{\sigma} = \begin{pmatrix} u_{11} & u_{12} \\ u_{21} & u_{22} \end{pmatrix}, \quad (5.15)$$

One then embeds these $SU(2)$ matrices in three $SU(3)$ matrices by mapping them as:

$$R = \begin{pmatrix} r_{11} & r_{12} & 0 \\ r_{21} & r_{22} & 0 \\ 0 & 0 & 1 \end{pmatrix}, \quad S = \begin{pmatrix} s_{11} & 0 & s_{12} \\ 0 & 1 & 0 \\ s_{21} & 0 & s_{22} \end{pmatrix}, \quad T = \begin{pmatrix} 1 & 0 & 0 \\ 0 & t_{11} & t_{12} \\ 0 & t_{21} & t_{22} \end{pmatrix}. \quad (5.16)$$

These three matrices are clearly members of the $SU(3)$ group and so is their product $X = RST$, which is used as the “random move” in the Metropolis algorithm. We thus have defined a recipe for numerically generating random transformations, the key element for our algorithm.

5.1.3 Efficient Action Difference Calculation

An additional element that is needed in the Metropolis algorithm is the action difference ΔS , used in evaluating the ratio of the PDFs between configurations. When a random transformation $U_\mu(x) \rightarrow U'_\mu(x) = XU_\mu(x)$ is applied on a single link of the lattice, the total change in the action only depends on the plaquettes that contain the considered link variable. In four dimensions there are 12 such elements:

$$\Delta S = S[U'_\mu(x)] - S[U_\mu(x)] = -\frac{\beta}{N} \text{Re} \text{Tr} ([U'_\mu(x) - U_\mu(x)] A[U_\mu(x)]), \quad (5.17)$$

where $A[U_\mu(x)]$ is the sum of the “staples” of the link $U_\mu(x)$. They are the constant three sides of the plaquettes that contain $U_\mu(x)$:

$$A[U_\mu(x)] = \sum_{\nu \neq \mu} \left[U_\nu(x + \mu) U_{-\mu}(x + \mu + \nu) U_{-\nu}(x + \nu) + U_{-\nu}(x + \mu) U_{-\mu}(x + \mu - \nu) U_\nu(x - \nu) \right] \quad (5.18)$$

An intuitive pictorial representation is given in fig. 5.1.

$$S[U_\mu] = \sum_{\nu \neq \mu} \left(\begin{array}{c} \text{Diagram of two adjacent plaquettes} \\ \text{Red link } U_\mu \text{ highlighted} \end{array} \right) + \dots = \text{Red link } U_\mu \times \sum_{\nu \neq \mu} \left(\begin{array}{c} \text{Diagram of two adjacent plaquettes} \\ \text{Red link } U_\mu \text{ factored out} \end{array} \right)$$

Figure 5.1: Schematic representation of the symmetric definition of action $S[U_\mu]$ expressed as a function of the staples. The link in red is the one that is being considered for the action difference, U_μ . On the left-hand side there is the action defined as the sum of all the plaquettes containing U_μ , on the right-hand side the link U_μ is factored out. The sum on the right-hand side is $A[U_\mu(x)]$ as defined in eq. (5.18).

Algorithm 2 Metropolis Update

```

1: for  $x, \mu$  do
2:    $A \leftarrow \text{computeStaples}(x, \mu)$ 
3:   for  $i < N_H$  do
4:      $U_{\text{new}}(x, \mu) \leftarrow X \cdot U(x, \mu)$ 
5:      $\Delta S \leftarrow (U_{\text{new}}(x, \mu) - U(x, \mu)) \cdot A$ 
6:     if  $\text{realTrace}(\Delta S) > \text{random}(0, 1)$  then
7:        $U(x, \mu) \leftarrow U_{\text{new}}(x, \mu)$ 
8:     end if
9:   end for
10: end for

```

5.1.4 Updates Strategies

There is now some arbitrariness in what is defined as a Monte Carlo update, or cycle. In this work we call an update the procedure outlined in algorithm 2.

These updates are the ones we consider when we refer to Monte Carlo cycles (MC cycles) in the context of thermalization and autocorrelation time, for example. Note that each update includes a loop over all links on the lattice and that every link is "hit" N_H times before moving to the next one. This is done for computational efficiency, because computing the staples is the most expensive part of the algorithm. Once $A[U_\mu(x)]$ is computed for a link, the result is used to attempt multiple updates on the same variable. It can also be shown that if N_H is sufficiently large the algorithm becomes equivalent to the heatbath algorithm [16].

The order in which the links in the lattice are visited is also arbitrary. The simplest way, that is iterating on the links as they are stored in memory, was adopted. This however might have impacted the autocorrelation of the system, not allowing significant modification to the system as an update depends on the neighbors. An alternative choice is a checkerboard pattern, which has potential benefits to the autocorrelation time of the observables as well as on the performance, as it allows for better communication patterns in the case of a parallelized code.

5.1.5 Parallelization Scheme

Given the size of the lattice (the total number of sites in this work is from $|\Lambda| \approx 10^5$ to $V \approx 10^7$), it is almost necessary to split the computation of the updates on more than one processors. The most direct way is to divide the lattice into sub-blocks, or sub-lattices, and have each process handle its portion of the field alone.

In an operative way, given a lattice of size (N_x, N_y, N_z, N_t) each space-time dimension is split into even portions, of size (S_x, S_y, S_z, S_t) and mapped on N_{procs} processors, having that:

$$N_{procs} = \frac{S_x}{N_x} \times \frac{S_y}{N_y} \times \frac{S_z}{N_z} \times \frac{S_t}{N_t}. \quad (5.19)$$

A unique mapping can be performed from the “local” coordinates of a lattice site in one of the sub-blocks and the global coordinates once the position of the sub-lattice is known. In practice this can be done through the aid of the unique identifier of the processor that is given by the parallelization library. In particular, the Message Passing Interface (MPI) has a built-in function to map processors into a grid of arbitrary dimensions and size, called `MPI_CartCoord_Create`. The utility also allows for the easy identification of the rank (the jargon term for the unique identifier of a processor) of the nearest neighbors. Using this tools one can manage the distribution of the total lattice into N_{procs} sub-lattices.

However, because of the dependence of the action difference of a link on its neighboring site, when updating a link on the edge of a sub-lattice, information about another block is needed. Here is where MPI comes in use, substituting the usual periodic boundary conditions that are used in the calculations with periodic communication patterns between the processors. For this particular problem we decided to use a point to point communication scheme, the use of periodic boundary conditions allowed also for non-blocking communications, in particular the collective geometry based scheme shown in section 5.1.5 has been implemented.

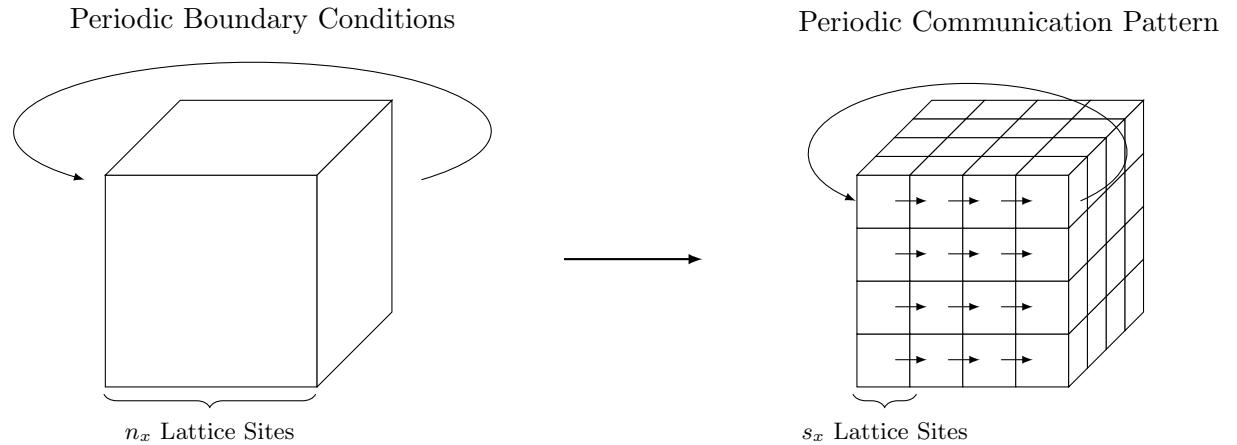


Figure 5.2: Schematic representation, in just 3 dimensions, of the splitting of the lattice into sub-lattices. The full system, which has periodic boundary conditions to emulate infinite volumes, is split into sub-blocks. These smaller regions of spacetime do not have periodic boundary conditions, but share faces with the neighboring blocks. To reproduce the periodicity, the communication pattern between the sub-lattices is made periodic, meaning that all sub-blocks aligned in one direction cyclically share faces with the neighbors.

When computing the staples, two links need to be fetched from the neighboring sites, since all processors are synchronized, when one processor is in need of these link variables one of its neighbors is requesting the corresponding links from it. The `SendRecv` function of MPI is exactly what is needed: it automatically handles, in the most efficient way, multiple data transfer between processors in situations when a single core needs to send and receive from different ranks.

Note that the communication is relevant only for the computation of the staples, this is another argument in favor of performing multiple hits on a link at every update. It is clear to see that the algorithm can be affected by large communication overhead problems. If the sub-blocks are too small, then most of the time would be spent on sending and receiving links from the neighbors. This causes the execution time to depend not linearly on the number of processors, instead the relation flattens at some value given by the communication overhead.

5.1.6 Summary of the Parameters for Generating Configurations

In total the algorithm we presented for generating pure gauge field configurations (using only the Wilson gluonic action) has four parameters as inputs. Of these, only one defines the physics of the system, the others have to be set and optimized in order to improve the acceptance ratio of the metropolis algorithm and the autocorrelation of the observables computed on the generated configurations. A summary is reported in table 5.1

Table 5.1: Parameters for the Monte Carlo generation of Yang-Mills gauge field configurations via the Metropolis Algorithm

Parameter	Description
β	Coupling parameter of the Wilson Action
ϵ	“Spread” of the random $SU(3)$ elements for the updates
N_C	Number of updates between one saved configuration (observable measurement)
N_H	Number of of “hits” per link at every update, with constant staple

5.2 Gradient Flow of Gauge Configurations

To study the flow time dependence of the observables a numeric implementation of eq. (4.7) is needed. The main challenge that this problem poses is how to numerically integrate the flow equation on the lattice while preserving the $SU(3)$ Lie Group structure of the field through the numerical procedure. The problem is solved by using the so called Structure-Preserving Runge-Kutta Methods, in particular the Runge-Kutta Munthe-Kaas (RK-MK) method that is designed to integrate first order ODE on a manifold [35, 36, 37].

In algorithm 3 the general scheme of the program to flow the gauge field configurations from time $t_f = 0$ to t_f^{MAX} in steps of size ϵ is outlined.

The main idea is to loop over all the configurations in an ensemble, provided as a set of input

Algorithm 3 Gradient Flow

```

1: for conf in ensemble do
2:   conf  $\leftarrow$  readInputFile()
3:   for  $t_f = 0, t_f < t_f^{MAX}$  do
4:     conf  $\leftarrow$  flowStep(conf)
5:     computeObservables(conf)
6:      $t_f \leftarrow t_f + \epsilon$ 
7:   end for
8: end for

```

files, and for each of them to numerically integrate the gradient flow equation eq. (4.5) in steps of ϵ (a dimensionless small step parameter). By using an explicit RK method, only one lattice is stored in memory at a time, and is updated at every iteration. At each step, the expectation values of all the desired observables are computed and their value stored to file for further analysis.

5.2.1 Numerical Integration of the Flow Equation

The problem of the gradient flow can be defined in more general terms as:

$$\dot{V}_t = Z(V_t)V_t. \quad (5.20)$$

here V_t is an element of a Lie group \mathcal{G} and $Z(V_t)$ is an algebra-valued function, in our case we have $Z(V_t) = -g_0^2 \partial_\mu S_G[V_t]$. The general ansatz for the RK-MK method is to write the integration step, with ϵ being the integration step, as:

$$V_{t+\epsilon} = \exp[\epsilon \Omega(V_t)] V_t, \quad (5.21)$$

where $\Omega(V_t)$ is a linear combination of $Z(V_t + \epsilon c_i)$, with c_i the RK coefficients, and of their commutators. EXPAND... Following the suggestion of [1] the third order integration scheme has been used:

$$\begin{aligned} W_0 &= V_t \\ W_1 &= \exp\left[\frac{1}{4}Z_0\right] W_0 \\ W_2 &= \exp\left[\frac{8}{9}Z_1 - \frac{17}{36}Z_0\right] W_1 \\ V_{t+\epsilon} &= \exp\left[\frac{3}{4}Z_2 - \frac{8}{9}Z_1 + \frac{17}{36}Z_0\right] W_2, \end{aligned} \quad (5.22)$$

where the shorthand notation $Z_i = i\epsilon Z(W_i)$ has been introduced. The choice of the coefficients for the Butcher's tableau [38] of the RK method is made in order to satisfy the requirements for a third order RK integrator and to cancel the commutator terms in $\Omega(V_t)$ when expanding it with the Baker-Campbell-Husdorff formula.

The choice of the steps size ϵ as been motivated by the results found in [39]. The article

suggests that for an integrator of third order the choice of $\epsilon = 0.01 - 0.02$ is sufficient to keep the numerical error under control.

What is then left to define are the derivative of the action at a given lattice site and the numerical exponentiation of a member of the $\mathfrak{su}(3)$ algebra, which returns an $SU(3)$ element.

5.2.2 The Action Derivative

The differential operator acting on an algebra-valued function $f(U)$ in a Lie group is defined, in terms of the generators t^a , as:

$$\partial_{\mu,x}^a f(U) = \frac{d}{ds} f(e^{sX} U) \Big|_{s=0}, \quad \text{with } X(\mu, x) = t^a \delta_{x,y} \delta_{\mu,\nu}. \quad (5.23)$$

In practice this operator leaves unchanged all the links except the one that is being derivated. When considering the action in eq. (3.33) and recalling the definition of staples eq. (5.18) we have:

$$\partial_{\mu,x}^a S_G(U_\mu(x)) = -\frac{2}{g_0^2} \sum_{x \in \Lambda} \sum_{\mu \neq \nu} \text{ReTr}(1 - P_{\mu\nu}) = -\frac{2}{g_0^2} \text{ReTr}(t^a U_\mu(x) A[U_\mu(x)]). \quad (5.24)$$

Defining $\Omega_\mu(x) = U_\mu(x) A[U_\mu(x)]$ and summing over all generators one has:

$$g_0^2 \partial_{\mu,x} S_G(U_\mu(x)) = 2t^a \text{ReTr}(t^a \Omega_\mu(x)) = t^a \text{Tr}(t^a (\Omega_\mu(x) - \Omega_\mu^\dagger(x))). \quad (5.25)$$

Invoking now the fact that any 3×3 matrix can be expanded as $M = \frac{1}{3}\text{Tr}M - 2t^a\text{Tr}(t^a M)$, we can rewrite the equation as:

$$g_0^2 \partial_{\mu,x} S_G(U_\mu(x)) = \frac{1}{2} (\Omega_\mu(x) - \Omega_\mu^\dagger(x)) - \frac{1}{6} \text{Tr} (\Omega_\mu(x) - \Omega_\mu^\dagger(x)). \quad (5.26)$$

This is our final expression for the action derivative. It should be noticed that computing this term is not anymore complex than computing the action difference for the gauge field configuration generation case, as in both cases the most complicated object to calculate on the lattice is the sum of the staples $A[U_\mu(x)]$. We can notice as well that the derivative is always a traceless hermitean matrix, thus an element of $\mathfrak{su}(3)$ as expected.

5.2.3 Exponential of an $\mathfrak{su}(3)$ Element

The last thing to define from eq. (5.22) is how to exponentiate the action derivative, that appears in the form $\exp[i\epsilon Z(W_i)]$. Following [40] we provide a numerical recipe for taking the exponential function of a traceless hermitean 3×3 matrix. The key idea is to use the Cayley-Hamilton theorem, that states that every matrix is a zero of its characteristic polynomial:

$$Q^3 - c_1 Q - c_0 \mathbb{1} = 0, \quad \text{with } c_0 = \det Q = \frac{1}{3}(Q^3), \quad c_1 = \frac{1}{2}(Q^2). \quad (5.27)$$

The exponential of iQ can then be written as:

$$e^{iQ} = f_0 \mathbb{1} + f_1 Q + f_2 Q^2, \quad (5.28)$$

where the f_i are functions of the eigenvalues of Q . These coefficients can be parameterized, because of the hermiticity of Q in terms of c_0 and c_1 . Labeling the eigenvalues of Q as q_1, q_2, q_3 and using the fact that the matrix is traceless a possible parametrization is:

$$q_1 = 2u \quad q_2 = -u + w \quad q_3 = -u - w, \quad (5.29)$$

with:

$$u = \sqrt{\frac{1}{3}c_1} \cos\left(\frac{1}{3}\theta\right) \quad w = \sqrt{c_1} \sin\left(\frac{1}{3}\theta\right) \quad \theta = \arccos\left[\frac{c_0}{2}\left(\frac{3}{c_1}\right)^{3/2}\right]. \quad (5.30)$$

As shown in [40] the coefficients f_i can be written as:

$$f_i = \frac{h_i}{9u^2 - w^2}, \quad (5.31)$$

with:

$$\begin{aligned} h_0 &= (u^2 - w^2)e^{2iu} + e^{-iu}\left[8u^2 \cos(w) + 2iu(3u^2 + w^2)\frac{\sin(w)}{w}\right], \\ h_1 &= 2ue^{2iu} - e^{-iu}\left[2u \cos(w) - i(3u^2 - w^2)\right], \\ h_2 &= e^{2iu} - e^{-iu}\left[\cos(w) + 3iu\frac{\sin(w)}{w}\right], \end{aligned} \quad (5.32)$$

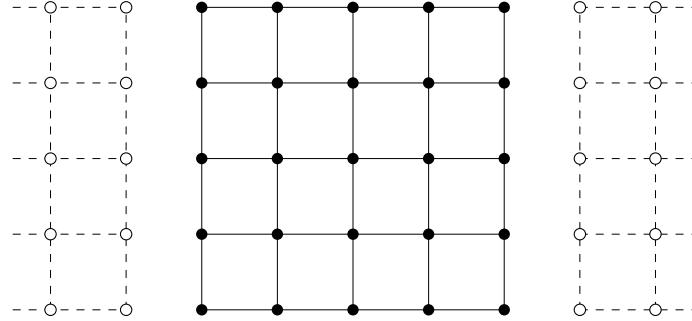
This algorithm is the core of the gradient flow implementation, because fundamentally the integration of the flow equation is an alternating sequence of staples calculation and matrix exponentials.

5.2.4 Parallelization Scheme

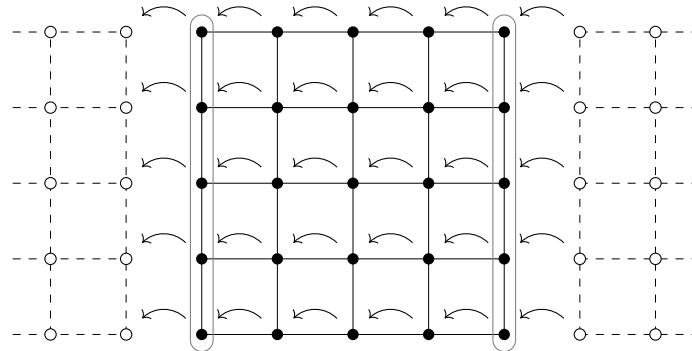
Since this problem is very expensive in computational terms, as was the gauge field generation, it is worth designing a parallelization scheme for it. The idea is still to split the lattice into sub-blocks as in section 5.1.5 and have each processing unit handle one of them.

By looking at eq. (5.22) we see that all operations can be defined on a lattice-wise scale. This allows us to define a “lattice shift” operation, that consists in creating an additional lattice in memory that is the translation of the original along an axis. Mathematical operations can then be performed on corresponding elements of the original lattice and of the shifted one. In order to create the shifted lattice some communication between the processors is needed, when the elements on the edge need to be shifted. The huge advantage is that in order to perform this shift operation only one communication instance can be used (following the previous scheme in section 5.1.5 for the periodic boundary conditions) in which the message is whole shared cube between two neighboring processors. This is much more efficient compared to the single link variable communication that is used in the gauge field generation. A simple 2D representation of the shift operation is drawn in fig. 5.3

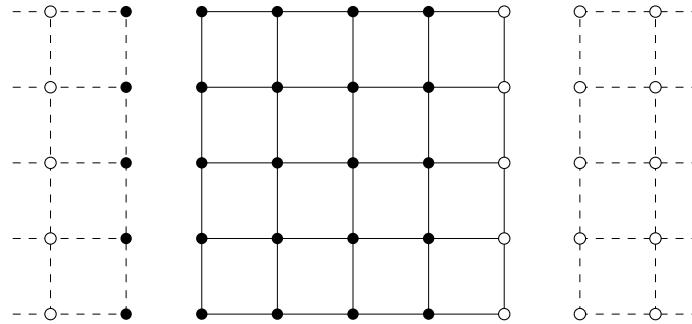
The data to be shifted is on the order of hundreds of kilobytes and sending such an amount of data using MPI is a time consuming operation. An possible optimization that has been implemented is to rewrite the algorithm in order to use non-blocking communications. These allow



(a) Normally, one processor only has a local lattice and no data from the neighboring sub-lattices. An empty copy of the sub-lattice is made, to store the shifted lattice.



(b) The shifted lattice is filled with the neighboring links, some of which are received by the “forward” neighbor (so some are sent as well to the “backward” neighbor)



(c) The shifted lattice is now a local object mapped with the same indices as the original one. A site-by-site multiplication is now possible between the original and the shifted lattices.

Figure 5.3: Parallelization scheme based on lattice shifts. When, for example in the staples computation, a link needs to be multiplied by a neighboring one, a new lattice is created. It is filled with the shifted link, which is fetched either from the local sub-lattice if it is in the central region, or from the neighbors if it is on the edge. This schematic drawing shows how a shift along the “right” direction is performed. A new lattice containing the data of the target link is created by filling it with the target links. The circled regions in fig. 5.3b, the edges, are shared using MPI by first creating a “buffer” of the shape of the edge (it is a cube in the real simulation) and it is sent in a single message to the neighbor which unpacks it in the right location. The new lattice is the multiplied with the original one on a site by site basis, which now is only a local operation.

the program , while the Message Passing Interface is handling the communication, to continue with the execution instead of waiting for the message to be sent. One must however be careful not to use any of the data that is being sent or received before checking that communication has been finished.

In the case of lattice shifts the use of non-blocking communications is simple and highly beneficial. First let us recall that, because of periodic boundary conditions, sending operations are always paired to a receive one. This means that a processor sends data to its neighbor on one side and receives the corresponding data from the neighbor on the other side along the same direction (this becomes clearer by looking at fig. 5.3). To start a lattice shift one first instantiates a non-blocking send instruction, then shifts the inner points of the sub-lattice (which do not require links from the neighbors) while the message is being sent. To complete the shift operation the last thing to do is to wait for the message from the neighbor from the receiving side to arrive.

This lattice shift operation has been implemented for the action derivative lattice and the gauge field itself. The massive reduction in communication overhead, compared to the single link exchange in section 5.1.5, improved the scaling of the algorithm greatly, making this part of the problem much more efficient.

5.2.5 Summary of the Parameters for the Gradient Flow

As it has been done for the gauge field configuration generation algorithm, in table 5.2 we report a summary of all free parameters of the program that performs the numerical integration of the gradient flow equation on the lattice.

Table 5.2: Parameters for the numerical integration of the gradient flow equations on Yang-Mills gauge field configurations.

Parameter	Description
β	Coupling parameter of the Wilson Action, fixed by the input configuration
ϵ	Runge-Kutta integration step
t_f^{MAX}	Maximum value of the flow time reached for every configuration.

5.3 Structure and Tools

The programs that have been developed to generate gauge fields and to apply the gradient flow to them have been written in C++. The full code can be found on the web under the link <https://github.com/GioPede>, where source code for both is hosted. The technical documentation is found at <https://giopede.github.io/LatticeYangMills/html/index.html>, but it should be noted that this is still in active development, though sufficient for this thesis, so the structure of the source code might vary from what is presented in this section.

The choice of programming language was mainly due to it being the most high-performing language with a high enough abstraction level. The high performance of C++ is given by the tight link that it has to the hardware, allowing precise management of the memory and instructions

optimizations.

One of the main features that distinguishes C++ from the C programming language (a valid and equally performing alternative), is that it is an Object-Oriented language. This allows to create abstractions that give an organized structure to the source code, making the developing process easier and the code more readable and intuitive.

5.3.1 Object-Oriented Programming

One of the most successful programming paradigms is Object-Oriented Programming. An object, or “class”, in this context is a collection of data, or “attributes”, and procedures, “methods”, that act on them.

The greatest advantage of using OOP is the introduction of an abstraction level from simple numerical variables, like integers, floating point variables or arrays, to more complex structures. A simple example of this is the class, that was implemented, to represent $SU(3)$ matrices. Instead of dealing every time with the allocation of 18 variables (3×3 real and imaginary) one can instantiate a variable of type `SU3`. This example is also useful to introduce the concept of operator overloading, which fundamentally means to implement mathematical operation between classes.

Finally, in OOP the concepts of inheritance and polymorphism are very important. The calculation of the observables is a good example for these. An observable, in the program, has a very well defined behavior and structure, for instance it stores his value and has a method that can be called to compute it given a gauge field lattice as input. However, the operations to be executed in order to compute the value depend on the observable type. An external class that wants to compute multiple observables should be able to call the compute method without knowing the exact type of the observable. The solution is to create an abstract base class, the observable, and define derived types for each different type that “inherit” the general behavior from the base class. For an external class, the derived objects can all be used in the same way, though the operations they perform are different, even if the exact type is unknown; this is called polymorphism.

5.3.2 Source Code Structure

It is possible to identify a layered structure in the class hierarchy of the program, outlined in fig. 5.4. At the time of writing the actual design of the source-code differs slightly from what is presented here (some classes are merged together, for example the observables), but development is being done to reach this cleaner and fully modular structure.

There is a set of “Core Classes”, the second layer in figure, on top of which all the rest is built:

- the $SU(3)$ matrix implementation, that specifies basic operation for the most fundamental object of Lattice QCD, the link variables. It also includes the algorithm for generating random matrices and the matrix exponential found in sections 5.1.2 and 5.2.3

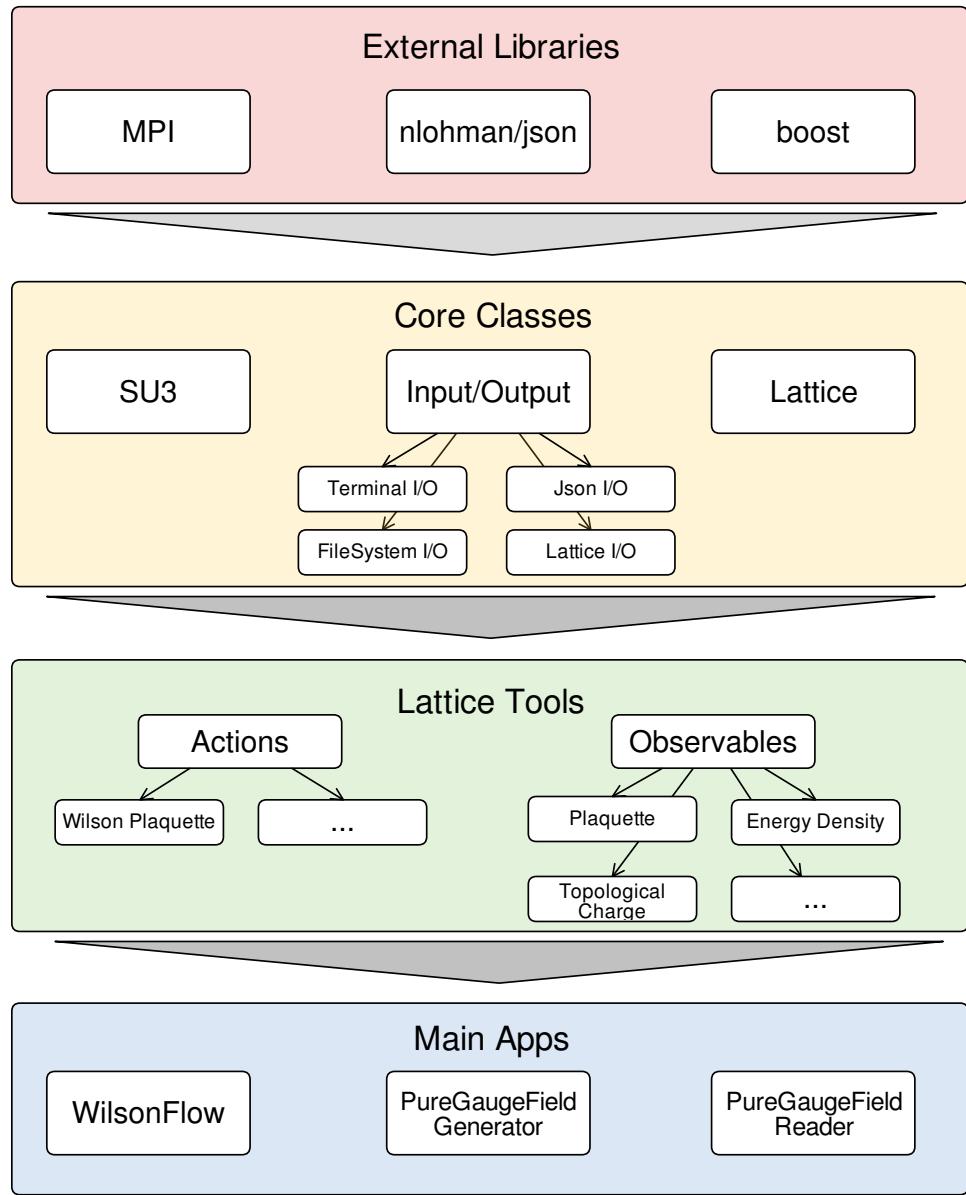


Figure 5.4: Class structure of the code, showing the approximative hierarchy between the different components.

- the Lattice class, which specifies operations on four dimensional grids. In particular it handles most of the parallelization tools to allow distributed memory computation, like on computing clusters. The parallelization is introduced by the use of the Message Passing Interface, MPI, to implement for example the lattice shift function, section 5.2.4.
- Input/Output tools, to handle input parameter files, gauge configuration parallel reading and writing, observable output files handling and standard command-line output. For the input files the choice has to write a `json` parsing class, based on the project found at [41], to provide an easier user interface via the usage of a structured input file.

With the tools described just above the main classes of the program were written. These “Lattice Tools”, the third layer in fig. 5.4, represent general physical concepts in Lattice QCD. They are:

- the Action abstract base class, of which the Wilson Plaquette action derived type has been implemented, is an object that if given a gauge field lattice can compute: the value of the action at every single lattice site; the staples of a link (for the action derivative and the Metropolis algorithm); the action difference of a link if provided with a random transformation for it.
- the Observable abstract base class. It contains simply the value of the observable and a method to compute it. The derived types that have been implemented are the Plaquette, the Energy Density and the Topological Charge.

For both of these base classes new derived types could be implemented, extending the capabilities and the use cases of the program.

Finally the “Main Applications” that have been developed are: the Pure Gauge Field Generator, which is an implementation of the metropolis algorithm as presented in this chapter; the Wilson Flow, a program to apply the gradient flow equation to a set of gauge field configuration and compute observables at every flow-time; a Pure Gauge Field Reader, mainly intended for testing, it is a useful application to view the single links of a lattice configuration.

Chapter 6

Tests and Runs Description

When developing a program for numerical simulations from scratch, it is crucial to design test cases and benchmarks to verify the goodness of the implementation. The easiest case is that of a fully deterministic program (for which the output depends only on some initial parameters) because one can compare the results with other implementations to validate the code. In the case of stochastic processes, for example that of a Markov chain, it is only possible to compare the average properties and results between different implementations.

In this chapter we will give an overview of the test cases, the properties of the ensembles used in this thesis and how the parameters that were needed to generate them have been set.

6.1 Generated Ensembles

In order to study the energy scale it has been necessary to choose a set of decreasing lattice spacings to then be able to take the continuum limit. This procedure is common to most Lattice QCD calculations as it is the most straightforward way to take recover the continuum theory. The lattice spacing is linked to the coupling g_0 , or the inverse coupling β , by the relation found in eq. (4.4).

We chose 4 values of β that span lattice spacings from approximately 0.1 fm to 0.05 fm in approximately equal steps.

For the calculation to be consistent however, the total volume of the lattice should be kept constant, so the choice of the lattice spacings also determined the number of lattice sites per dimension, having $L = aN \approx const.$. The time dimension has been take to be twice as big as the three spatial dimensions. The number of lattice points per dimension, for the sake of parallelization, were chosen in order to have many integer divisors, to allow for different and small sub-blocks. Table table 6.1 summarizes the physical properties of the ensembles that were generated.

For each value of β a statistical ensemble of randomly chosen gauge field configurations was needed. Ideally, one would take as many configurations as possible for each value. However it is clear from the discussion in chapter 5 that the number of lattice sites affects computation

Table 6.1: Physical properties of the ensembles used for this work. The parameter β is the inverse coupling, and its value fixed the lattice spacing a through the usage of eq. (4.4). The value of N represents the number of lattice sites per spatial dimension and N_T , here fixed to $2N$, the number of sites in the time dimension. The size $L = aN$ of the system is also reported in fermi.

β	a	$N^3 \times N_T$	LN [fm]
6.00	0.09314	$24^3 \times 48$	2.23536
6.10	0.07905	$28^3 \times 56$	2.21367
6.20	0.06793	$32^3 \times 64$	2.17405
6.45	0.04781	$48^3 \times 96$	2.29488

times and the capability of storing the configurations dramatically. Moreover, as will be shown in section 6.2.1 the autocorrelation time of the observables has a non-trivial, power-law or exponential behavior with the lattice spacing, making the generation of the larger β ensembles even more challenging.

table 6.2 reports the final values for the parameters of the Metropolis algorithm for the different ensembles that were generated. These values have been chosen after many tests, mainly checks

Table 6.2: Parameters used for the generation of the ensembles on which the results of this work are based on. In the table β is the inverse coupling; “MC Cycles” stands for the total number of Monte Carlo updates performed, a detailed description of an update is found in section 5.1.4. The following parameters are the algorithm specific ones described in table 5.1.

β	MC Cycles	N_{conf}	N_{corr}	N_{hit}	$\epsilon_{SU(3)}$
6.00	600000	1000	600	30	0.25
6.10	300000	500	600	30	0.25
6.20	400000	500	800	30	0.25
6.45	400000	250	1600	30	0.25

of the autocorrelation time of the observables presented in section 3.3.2. All the parameters are free in principle and no reference study on their impact on the resulting ensemble properties was found for the case of the Metropolis algorithm. In the next section the trial and error approach that led to this decision will be briefly discussed.

Evidently, there is less freedom on the choice of the parameters in table 5.2 for the gradient flow compared to the gauge field generation. This is because the numerical integration of the flow equation is fully deterministic. There are only two quantities to fix: the value of t_f^{MAX} , the final flow time point; and the size of the integration step ϵ . In the calculations the flow time is expressed in lattice units, $t_f \rightarrow t_f/a^2$, so the quantities we set were in this dimensionless form. To rescale the final flow time to physical units it must be multiplied again by the square of the lattice spacing, but it is more interesting to report the final smearing radius of the flow equation, remembering that it is defined as: $\sqrt{8t_f^{MAX}}$. In table 6.3 we report the choice of the parameters for the flow of the ensembles in table 6.1.

The choice of ϵ was motivated by the work by Cè et al. [39], and the same values for the

Table 6.3: Parameters used for the numerical integration of the gradient flow equations eq. (4.7) for the ensembles in table 6.1. The inverse coupling is β and its value implicitly defines the lattice spacing a . ϵ is the size of every Runge-Kutta integration step. The value of t_f^{MAX}/a^2 is the last point of integration in dimensionless units and the corresponding maximum smearing radius is shown as $\sqrt{8t_f^{MAX}}$ in fermi.

β	ϵ	t_f^{MAX}/a^2	$\sqrt{8t_f^{MAX}}$ [fm]
6.00	0.01	10.00	0.8330
6.10	0.01	10.00	0.7070
6.20	0.01	10.00	0.6076
6.45	0.02	20.00	0.6047

combination of ϵ and t_f^{MAX}/a^2 have been chosen for all ensembles, regardless of the physical maximum smearing radius. The exception is the $\beta = 6.45$ case, for which the total number of integration steps of the Runge-Kutta Munthe-Kaas integrator has been kept equal to the other lattice sizes, but the integration step length is doubled. This was needed to obtain results for large enough smearing radii out of the ensemble (which is by far the largest one) in reasonable computing time.

6.2 Test Runs

Running some test calculations with a completely new program is obviously necessary. The simplest benchmark was to check the expectation values of the observables. To perform these checks we computed the energy density and the topological charge on two configurations generated with the CHROMA lattice QCD code from the USQCD collaboration[42]. The results were found to match to machine precision.

The gradient flow implementation could also be tested numerically by comparing our results to the ones generated by an extension of CHROMA called FlowOps (courtesy of T.Luu and A.Shindler)[43]. This program as well applies the gradient flow to gauge field configurations with the same integrator scheme that we used eq. (5.22), thus by setting the same integration parameters we were able to compare the values of the observables at every flow time. Also in this case our implementation gave results equal to machine precision to the ones by FlowOps at all flow times for every configuration we tested.

Checking the validity of expectation values of observables is a solid indication that the overall core libraries of the new framework we developed has been implemented correctly.

Testing the algorithm to generate gauge field configurations and assessing the quality of the generated ensembles is much harder than testing the gradient flow, because it involves stochastic computations, hence no numerical check can be easily defined. One has instead to look at average properties of the set of configurations. Unfortunately the generation of gauge field configurations as we saw in table 5.1 has three parameters that need to be set that do affect the statistical properties of the output ensemble, mainly the autocorrelation time of the observables. The simplest parameter to set was $\epsilon_{SU(3)}$, which controls the spread of the $SU(3)$ random elements around the identity matrix. This directly affects how much each link is changed from

one MC cycle to the other, a larger value of $\epsilon_{SU(3)}$ implies larger changes. The drawback is the decrease of the acceptance ratio of the Metropolis test and consequently of the efficiency of the program. The value that was chosen, 0.25 was a good trade-off between the two things. The other parameters two parameters to set were: N_{corr} , the number of configurations between observable measurements (in practice between two saved configurations); (N_{hit} , the number of random unitary transformations applied to every link per MC cycle). Some tests runs with different combinations of the two were performed on the three largest lattice spacings, as the system for the smallest one is much larger and requires a considerably longer time to compute.

6.2.1 Tests for the Autocorrelation of Observables

The most important test has been the assessment of the autocorrelation time for different observables at various lattice spacings varying the parameters of the ensemble generation algorithm. As expected, the most problematic quantity is the topological charge, especially at large flow times. A good measure of how much a data series is autocorrelated is the integrated autocorrelation time τ_{int} , see section A.2, which is expected to be 1/2 if the data is uncorrelated. In general, a larger τ_{int} implies an underestimation of uncertainties, so the variance of a quantity is corrected as $\tilde{\sigma}^2 = 2\tau_{int}\sigma^2$.

The first test, in fig. 6.1 shows the integrated autocorrelation time for the topological charge at fixed $N_{corr} = 200$ and $N_{hit} = 10$ for different lattice spacings:

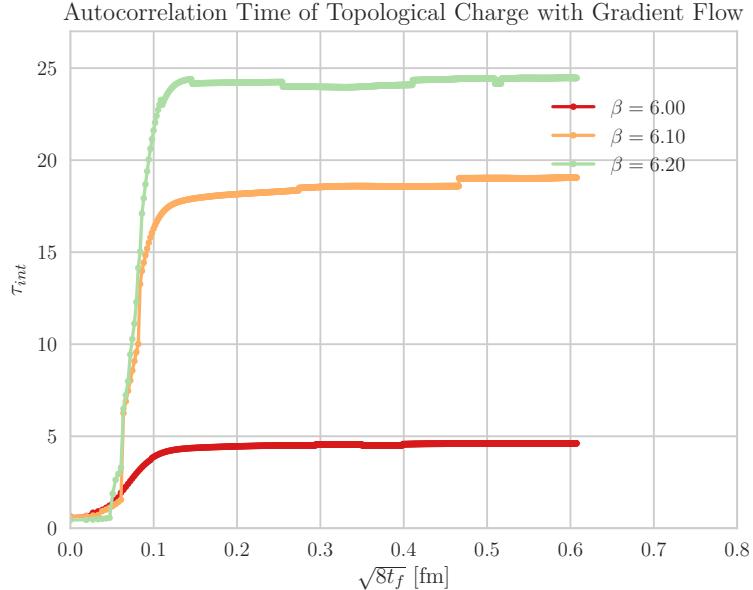


Figure 6.1: Integrated autocorrelation time τ_{int} for fixed values of $N_{corr} = 200$ and $N_{hit} = 10$. The different data series represent different inverse coupling values, the system sizes are the ones of table 6.1.

From the plot it is clear that the values that were guessed for the parameters N_{corr} and N_{hit} ,

which could be acceptable for the lowest beta value, are not at all fine for the other lattice spacings. One can observe that for increasing values of β the τ_{int} grows rapidly if the parameters of the generation algorithm are kept fixed. This is the reason why in table 6.1 the different ensembles have different parameters.

On a general note we observe how the flow quickly removes the low range noise from the observable through the smearing, making it clear that the topological charge is highly correlated in Monte Carlo time even though it might seem uncorrelated if one only looks at $t_f = 0$.

A second test was to look at a single lattice spacing and try to vary the parameters N_{corr} and N_{hit} . The results for the $\beta = 6.10$ case are shown in .

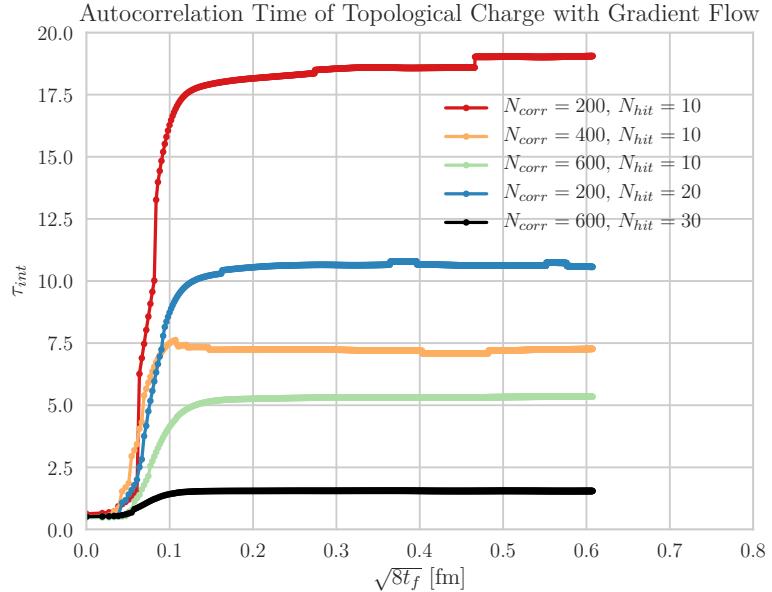


Figure 6.2: Integrated autocorrelation time for $\beta = 6.10$ for a lattice of size $28^3 \times 58$. The different data series represent different combinations of the parameters for the gauge field generation N_{corr} and N_{hit} . The black data series shows the set of parameters that were chosen for the generation of the larger ensemble (the one used for the analysis of the energy scale Λ_{YM}).

The results of fig. 6.2 suggests a possible cure for the autocorrelation of the observables. It is almost tautological that increasing the value of N_{corr} decreases the autocorrelation time, as the parameter represents how many MC updates are performed between two measurements of the observable. The dependence of τ_{int} from N_{hit} is also clear since the more attempts of transforming a single link are performed, the more the configuration will be different from the previous one.

From the plot we can observe that increasing both parameters improves the autocorrelation, so the choice was to set both to a larger value from our initial guesses. This proved to be the right choice.

The black data series in fig. 6.2 is the data that we used for our analysis. One can notice that it is still autocorrelated, but to a degree that can be handled with a reasonable correction to the variance of the data.

It is important to stress again that the autocorrelation is a huge problem mainly for the topo-

logical charge. The energy density operator is much less affected and for all the ensembles we generated the integrated autocorrelation time is never a value much greater than $1/2$ (see fig. 7.4).

6.2.2 Thermalization

When computing expectation values using the Monte Carlo integration, it is important to ensure that the sampling of the observables is performed when the Metropolis algorithm reached equilibrium, or in jargon when it is thermalized.

The Markov Chain to generate gauge field configurations can be initialized in two distinct ways. These represent the initial configuration of the gauge field to which Monte Carlo updates are later applied. The two possibilities are:

- Cold Start: all links of the gauge field configuration are set to the $SU(3)$ identity element;
- Hot Start: all links are set to random $SU(3)$ matrices.

It is easy to see that for the cold start the initial value for the plaquette operator would be 1, when normalized on the lattice volume and so the action starts in its minimum. For the hot start the initial value for the plaquette expectation value is a random number, centered around 0. The Metropolis algorithm is supposed to drive the Markov chain towards the peak of the PDF, eq. (5.9), that corresponds to an intermediate value of the two.

In fig. 6.3 we show that only a small number of MC updates is needed to reach equilibrium, especially when compared to the value of N_{corr} from the discussion in section 6.2.1.

To ensure thermalization, all results of this work are generated discarding the first 10^4 MC cycles, which is much larger than the value we found, but still a small fraction, so an affordable exaggeration, of the total number of MC cycles of the Markov chain.

6.2.3 Strong and Weak Scaling

One important thing to look at when developing a program intended to be highly parallelized are the scaling properties. We can now check these for the two sections of the code, the generation of configurations and the gradient flow. We will distinguish the analysis in the two usual quantities used in High Performance Computing: strong and weak scaling.

Strong Scaling

Strong scaling is the performance of a program as the total problem size is kept fixed and the number of parallel units, the number of processors in this case, increases. The parallel efficiency

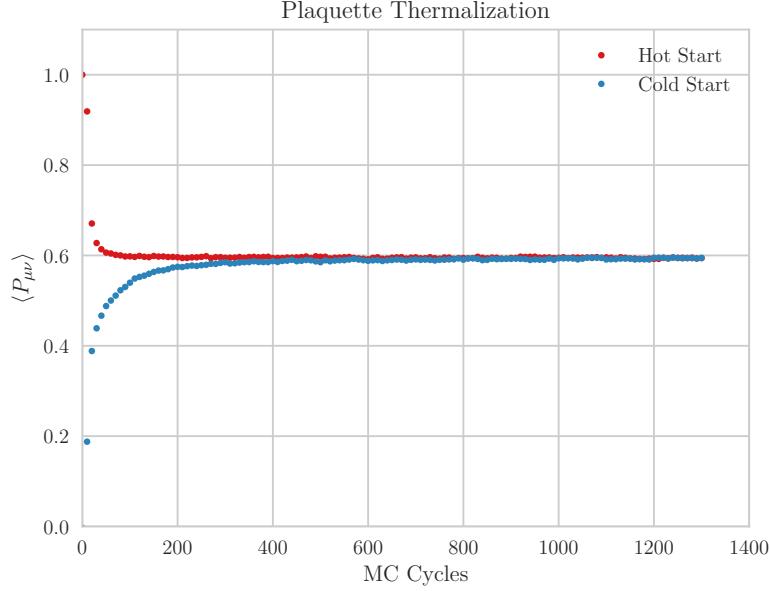


Figure 6.3: Expectation value of the Plaquette as a function of the number Monte Carlo cycles performed, for hot and cold starts. Equilibrium is reached in both cases after ≈ 600 updates. The data is taken from a simulation with $\beta = 6.10$.

is computed as:

$$\eta_s = \frac{t_0}{t_N N} \cdot 100\% \quad (6.1)$$

where t_0 is the execution time for one unit of computation, t_N the execution time for the case of N units computing the same system size. In the case for perfect parallelization the efficiency is 100% as an increase of some factor in the number of processors would decrease by the same factor the execution time. In real cases this is seldomly true, the execution time is in fact usually larger than expected.

The test was set up in this way:

- Gauge Field Generation: one lattice of size $24^3 \times 48$ has been randomly generated and 10^4 MC cycles were performed. The system was split in sub-lattices of decreasing size, so to increase the number of processors.
- Gradient Flow: ten lattice of size $24^3 \times 48$ have been flowed for a total of 1000 integration steps each. Again, the sub-lattice size was decreased and the number of processor increased accordingly.

Figure 6.4 shows that the performance scales well with the number of processors, but the efficiency is dropping for very large processor numbers. This is expected, the reason is that the sub-lattice the surface to volume ratio of the sub-lattice makes the fraction of operations that require inter-processor communication become comparable to the "inner" points calculations.

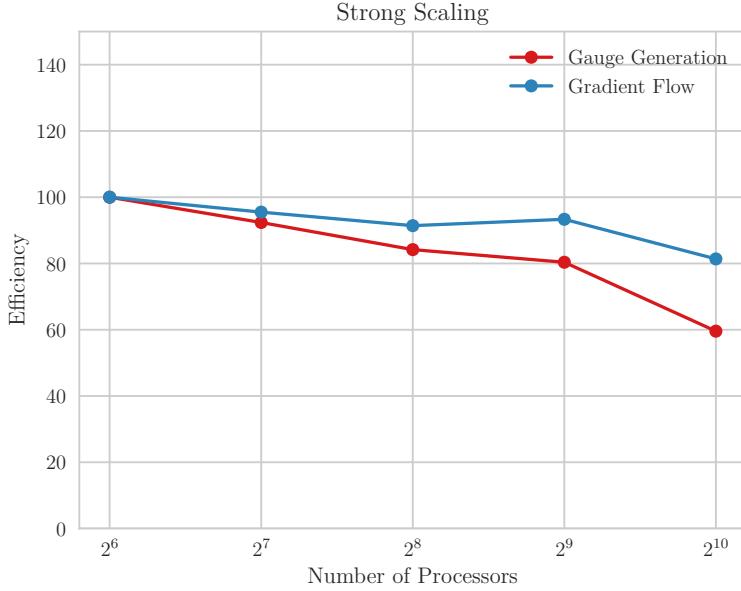


Figure 6.4: Strong scaling of the two parts of the program. The efficiency, eq. (6.1), is normalized to the first point plotted, that is 64 processors.

It is to be noted that smaller number of processors could not be used, as the processors in that case would have been all on the same node (the cluster used for testing has 28 cores per node), so any timing would have excluded the inter-node communication time, which is significantly higher than the inter-node one.

Overall the gradient flow program scales better than the ensemble generation one. This is most likely explained by the fact that the communication overhead introduced by MPI is larger in the case of the generation program (which shares one link variable at a time) than in the gradient flow code (which uses lattice shifts and shares one whole shared edge between processors at once, see section 5.2.4 for details).

Weak Scaling

By weak scaling of a parallelized problem it is intended the measure of the performance as the local workload is kept fixed and the total volume of the problem is increased by adding processors. The efficiency is computed as:

$$\eta_W = \frac{t_0}{t_N} \cdot 100\% \quad (6.2)$$

where t_0 is the execution time of the smallest case and t_N the one of a system N times larger computed on N times the parallel units. The test set up was:

- Gauge Field Generation: the sub-lattice has been fixed to 4^4 , and volumes from $16^3 \times 32$, a total of 32 processes, to $32^3 \times 64$, 512 processors, have been considered. Intermediate

volumes are built by multiplying one spatial dimension at a time by a factor of 2. For each volume 10000 MC cycles have been computed.

- Gradient Flow: the gauge fields generated in the previous point have been flowed for a total of 1000 integration steps by the same number of processors that generated them, so the sub-lattice size was set again to be 4^4 .

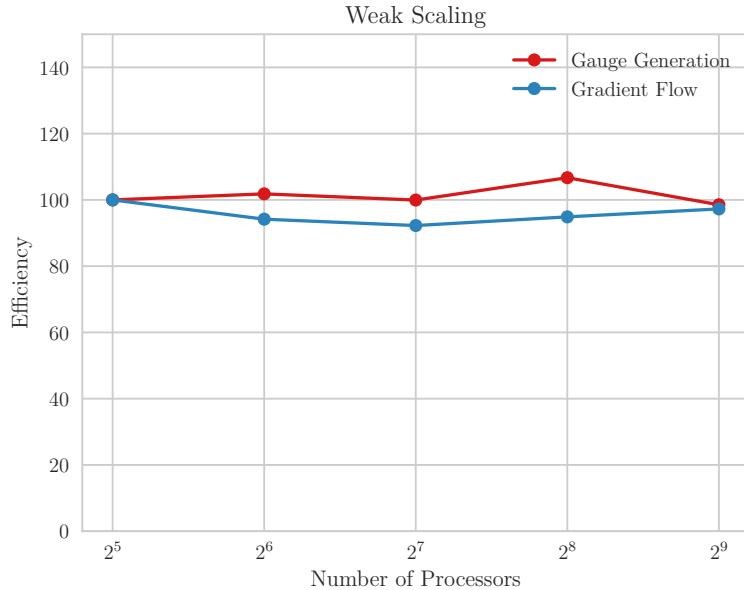


Figure 6.5: Weak scaling of the two parts of the program. The efficiency is the ratio of eq. (6.2) with respect to the first data point.

The results in fig. 6.5 show that both the programs are perfectly scalable. The communication overhead, introduced by the need of sharing link variables across the edges, is constant with respect to the number of processors. We should note that already the first point plotted contains a large amount of communications between processors. What the plot is truly suggesting is that the addition of other processors keeps the overall overhead constant. This implies that the parallelization procedure has been implemented correctly, though it could still be potentially improved by further optimizations.

The fact that some points have an efficiency larger than 100% is most likely due to the fact that the first data point, for the gauge field generation, is unreliable. This could be due to some issue related to the machine where the test was run that resulted in a slightly longer than average execution time.

These results for strong and weak scaling are very comforting, as they show that the hardest part of the development process, the parallelization of the code, has been well designed and implemented.

6.3 Main Runs and Timing

All runs of the programs, to generate the four ensembles in table 6.1 and to apply and integrate the gradient flow equations to all their configurations, were carried out on the High Performance Computing Center at Michigan State University (MSU), with the support of the Institute for Cyber-Enabled Research (iCER). Development was performed on local machines and on the small cluster SMAUG located at the Department of Physics of the University of Oslo (UiO) and some larger benchmarks were run on the Abel Computer Cluster also at UiO.

The ensembles, on which the results in part III were generated using the parameters in table table 6.2. A summary of the computational resources used for the generation of all four ensembles is presented in table 6.4.

Table 6.4: Execution times for generating ensembles from table 6.1 with parameters found in table 6.2. All the runs were performed on the iCER cluster at MSU. N_{procs} is the number of processors used for each ensemble; the Wall Time is the “wall clock” time (the time spent in the parallel execution); Time per MC cycle is the time spend in each step of the Markov Chain; Time per Configuration is the time spent between saved configurations, to be used for observables, so it includes the adjustment made for the longer autocorrelation times of larger β values.

β	N_{procs}	Wall Time [h]	Time per MC Cycle [s]	Time per Configuration [s]
6.00	1024	27	0.16104	96.6
6.10	512	35	0.42102	252.6
6.20	1024	53	0.4780825	382.4
6.45	1024	204	1.836	2937.6

It is clear from the table that the largest set, out of which only 250 configurations were saved, had execution times larger by one order of magnitude compared to the second largest; this made increasing the ensemble size unfeasible with the time and resources for this work.

To conclude, in table 6.5 we present the summary of the computational resources used by the Gradient Flow program to integrate the flow equation on all the configurations of the ensembles.

Table 6.5: Execution times for the numerical integration of the flow equations on the ensembles from table 6.3 with parameters found in table 6.3. All the runs were performed on the iCER cluster at MSU. N_{procs} is the number of processors used for each ensemble; the Wall Time is the “wall clock” time (the time spent in the parallel execution); Time per Configuration is the time spent flowing one single gauge configuration.

β	N_{procs}	Wall Time [h]	Time per Configuration [s]
6.00	1024	15	109.3
6.10	512	30	218.6
6.20	1024	33	243.6
6.45	1024	96	1326.9

Part III

Data Analysis and Results

Chapter 7

Raw Observables

In this chapter the results for simple observables obtained from the ensembles in table 6.1 are presented. Throughout the chapter, all error estimates associated to expectation values of observables have been computed using the bootstrap method, see section A.1.1, which is a popular resampling method used when the sample size of a statistical population is small, as is our case. To handle the autocorrelation of the observables, introduced by the Markov chain, is handled using the procedure found in section A.2 as a correction to estimate of the variance proportional to the integrated autocorrelation time.

7.1 Plaquette and Energy Density

The plaquette, eq. (3.27), and the energy density, eq. (3.52), are tightly related as they both are estimates of the action of the gauge field. Since the system is expected, after thermalization, to be at equilibrium, the sampling of the plaquette and of the energy density should be rather constant in Monte Carlo time. Figure 7.1 shows that this is indeed the case.

When applying the gradient flow equation:

$$\partial_{t_f} V_{t_f}(x, \mu) = -g_0^2 [\partial_{x,\mu} S_G(V_{t_f})] V_{t_f}(x, \mu), \quad (7.1)$$

the configuration is evolved towards the minimum of the action. That implies that the values for the plaquette and the energy density at sufficiently large flow times are both lattice spacing and flow time independent. The results plotted in fig. 7.2 for the plaquette expectation value as a function of flow time, show that the gradient flow is indeed driving the gauge fields towards the stationary points. From the expression of the Wilson gluonic action, eq. (3.33), it is clear that the minimum of the action is reached when all the plaquettes are equal to the identity matrix. The expectation value of the average plaquette is therefore expected to be exactly one.

The energy density has an even simpler proportional relation to the action and fig. 7.3 proves that the action is minimized.

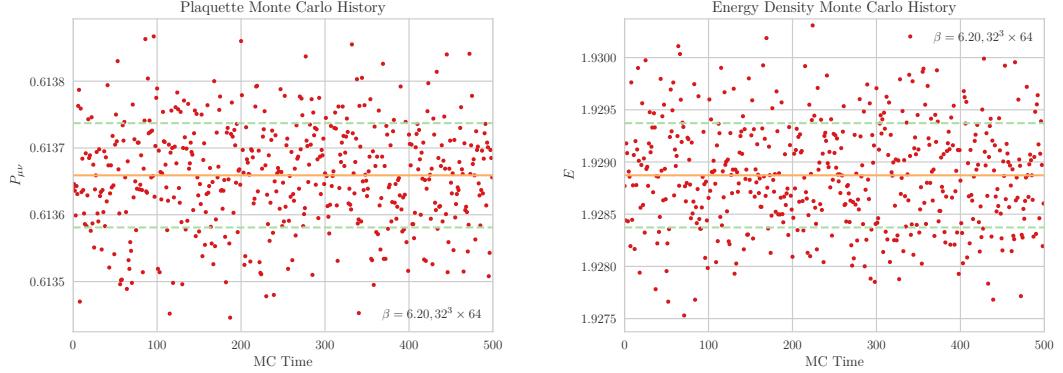


Figure 7.1: Values for the Plaquette (left) and Energy Density (right) as a function of Monte Carlo time. The blue line is the average and the green dashed lines are the one σ interval around it. The points represent the values of the two observables computed on every configuration of the ensemble. The average is thus the expectation value.

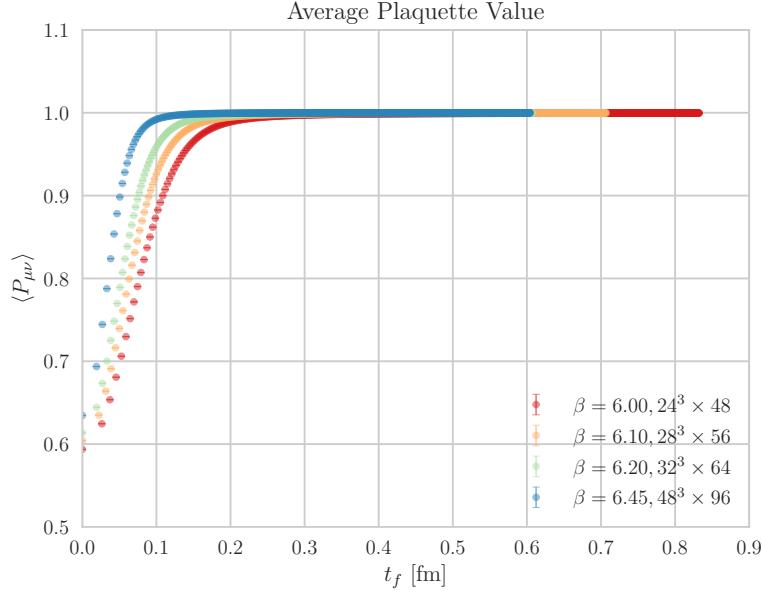


Figure 7.2: Plaquette expectation value $\langle P_{\mu\nu} \rangle$ as a function of flow time t_f , here expressed in the form of the smearing radius $\sqrt{8t_f}$. The data for all the four ensembles of table 6.1 is shown. Errorbars are too small to be visible on the plot.

As discussed in section 6.2.1, the integrated autocorrelation time of the observables is a crucial quantity to assess the quality of the results. We therefore computed τ_{int} for all the different ensembles at every flow time. From fig. 7.4 we conclude that for all ensembles we generated the value of τ_{int} for the energy density never exceeds the value of one. This result is suggesting that the data we sampled for this observable is uncorrelated. Nevertheless, all results of this work (including fig. 7.2 and fig. 7.3) have variances corrected by $\tilde{\sigma}^2 = 2\tau_{int}\sigma^2$. **DA CALCOLARE**

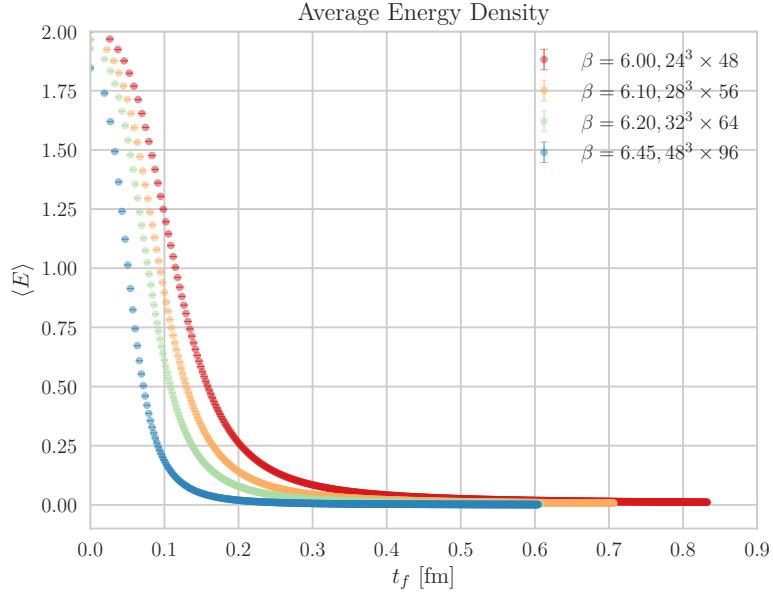


Figure 7.3: Energy density expectation value $\langle E \rangle$ as a function of gradient flow smearing radius $\sqrt{8t_f}$. The different data series represent the four different ensembles that have been generated, see table 6.1. The errorbars not visible on the plot as they are small.

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7.2 Topological Charge

A more interesting quantity to measure on the ensembles of gauge field configurations that we have generated is the topological charge. In the continuum theory it assumes only integer values, with a distribution around zero that resembles a gaussian, though there are studies that prove that indeed it is not a normal distribution [44].

The topological charge on the lattice is affected by short range fluctuations given by discretization effects. The smoothing properties of the gradient flow on this observable remove the fluctuations and reveal the “true” value of the charge. An illustrative example is to plot the topological charge of one gauge field configuration as a function of the smearing radius. In fig. 7.5 we show the evolution of the topological charge for three different gauge field configurations of lattice spacing $a = 0.049$ fm ($\beta = 6.45$).

We can notice that for all the three considered configurations the topological charge has a plateau at a smearing of roughly 0.15 fm. This value corresponds to approximately three times the lattice spacing.

For some configurations however, especially for larger lattice spacings, the plateau is not clear and defined. The explanation could be the too simple definition that has been used for the gauge field strength tensor. Perhaps an improved definition of the tensor, given for example by

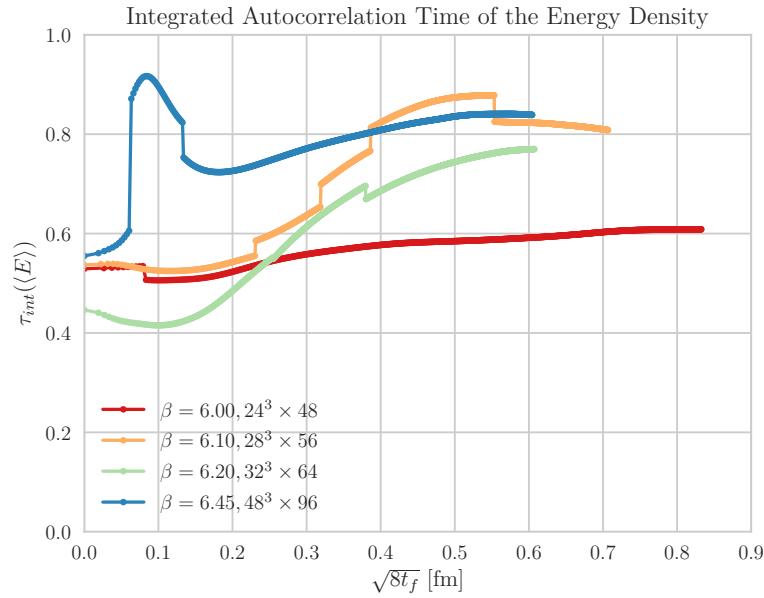


Figure 7.4: Integrated autocorrelation time of the Energy Density as a function of the smearing radius $\sqrt{8t_f}$. The discontinuity of the data is given by the approximation given by the truncation procedure described in section A.2

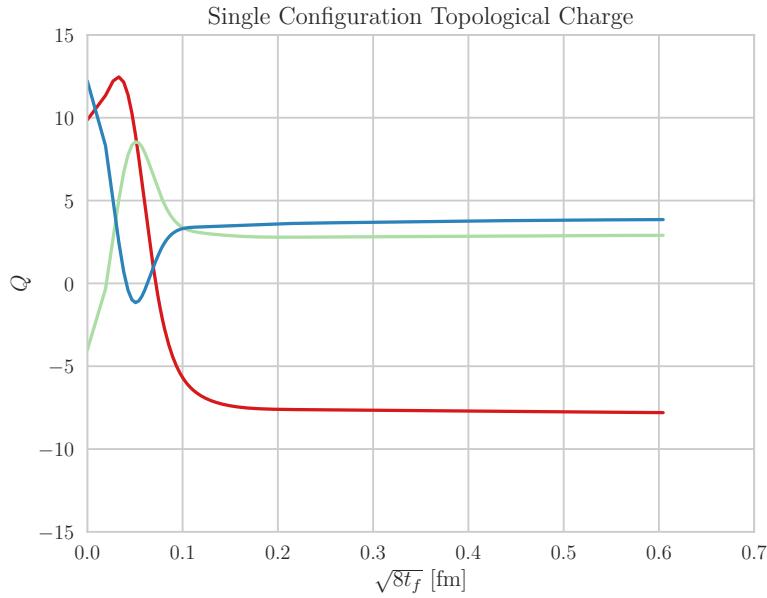


Figure 7.5: Topological charge for three single gauge field configurations, taken at random from the $\beta = 6.45$ ensemble, plotted against the smearing radius $\sqrt{8t_f}$ of the gradient flow.

a linear combination of plaquettes and larger Wilson Loops could be beneficial[22].

When considering the expectation value of the topological charge of an ensemble, we expect its value to be zero [44]. From fig. 7.6 one can check that our data for the expectation value of $\langle Q \rangle$ is indeed within error-bars always compatible with zero.

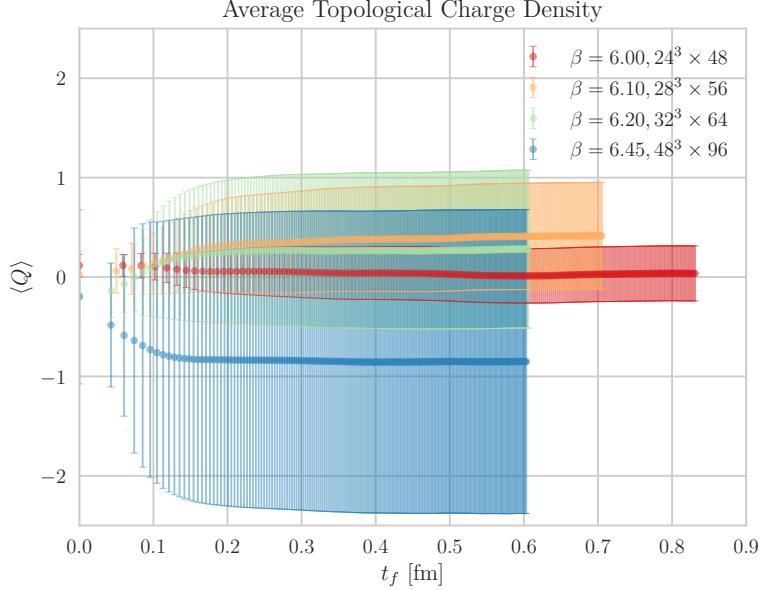


Figure 7.6: Topological charge expectation value $\langle Q \rangle$ as a function of the smearing radius $\sqrt{8t_f}$. Errorbars are computed using bootstrap and corrected with the integrated autocorrelation time correction factor.

We should note that the increasing error-bars, for $\beta = 6.2$ and $\beta = 6.45$ in particular, are given by the decreasing ensemble size and the increasing autocorrelation time τ_{int} . The latter is plotted in fig. 7.7. By comparing it with the equivalent plot for the energy density (fig. 7.4) the difference is clear. The topological charge not only has a much larger autocorrelation time on the lattice, but it grows rapidly with the inverse lattice spacing. It should be noted that the value of N_{corr} , the Monte Carlo updates between two data points, is much larger for the $\beta = 6.45$ ensemble than all the others, but nevertheless the τ_{int} is by far larger.

One last interesting consideration on the topological charge is the comparison of its distribution at zero flow time and the one in the large flow time limit. In fig. 7.8 we show two histograms of such distributions for the $\beta = 6.20$ ensemble at $\sqrt{8t_f} = 0$ fm and $\sqrt{8t_f} = 0.6$ fm respectively. The bins are taken to be centered on integer and half-integer values in order to capture differences of the distribution of the topological charge. In the continuum only integer values should be allowed, while on the lattice, due to discretization, the topological charge is in general not an integer. However, from fig. 7.8 it is possible to see that the gradient flow estimate of the charge has a distribution with much sharper peaks around integer values, rather than half-integer ones, if compared to the zero flow time distribution.

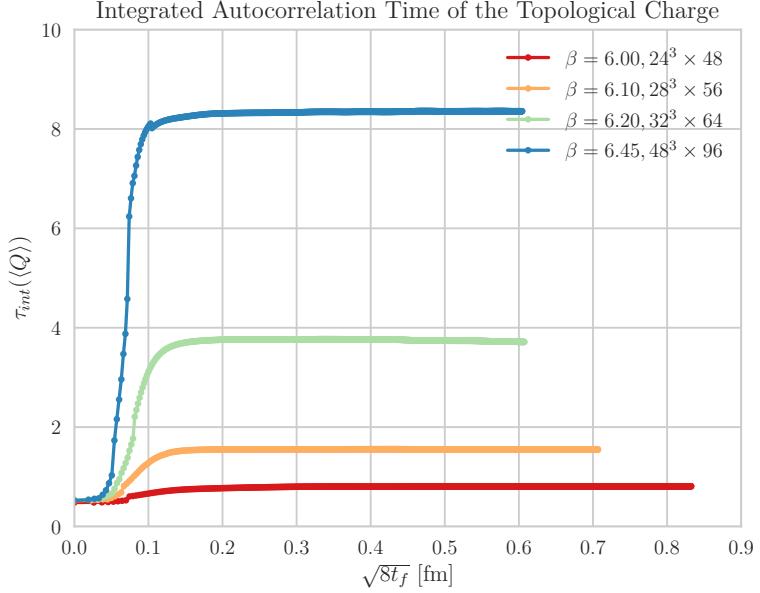


Figure 7.7: Integrated autocorrelation time τ_{int} of the topological charge as a function of the gradient flow smearing radius $\sqrt{8t_f}$. The ensembles are the ones reported in table 6.1.

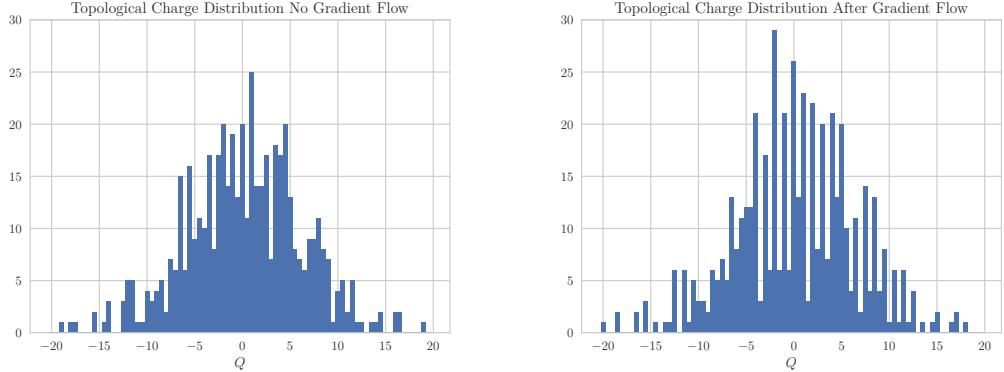


Figure 7.8: Histograms of the topological charge values at $\sqrt{8t_f} = 0$ fm (left) and $\sqrt{8t_f} = 0.6$ fm (right) for the $\beta = 6.20$ ensemble. The bins are centered on integers and half-integer values.

7.3 Topological Susceptibility

Another quantity that is interesting to measure from the gauge field configurations that we generated is the topological susceptibility. Being proportional to the expectation value of the squared topological charge, it is related to the width of the distributions in fig. 7.8. The evolution with the gradient flow of the susceptibility is not trivial. At low flow times the high-frequency fluctuations in the topological charge make the expectation value of the susceptibility diverge. The smoothing property of the gradient flow remove the fluctuations revealing the real value of

the quantity. In fig. 7.9 the flow time evolution of the topological susceptibility is shown.

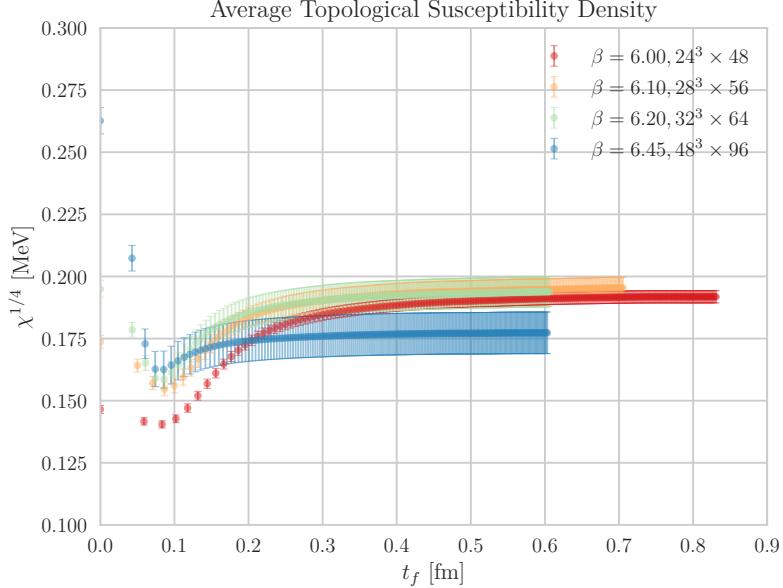


Figure 7.9: Topological susceptibility $\chi^{1/4}$ as a function of the smearing radius $\sqrt{8t_f}$, for all the ensembles generated for this work.

The value of $\chi^{\frac{1}{4}}$, as defined in eq. (3.58) has the physical units of energy, hence it is reported in MeV. We notice that for all ensembles in the large flow time limit the topological charge has a well defined plateau. It is then possible to take the continuum limit in a region between 0.4 fm and 0.6 fm.

The behavior at zero flow time is not what we were expecting, comparing for example to ref. [43] we don't observe a clear divergency at $\sqrt{8t_f} = 0$ fm. The behavior seems to be more and more divergent (a true divergency is impossible, it is always a finite quantity on the lattice) for larger values of β . A possible explanation of this behavior is the usage of the clover definition of the gauge field strength tensor. Being the topological charge, at zero flow time, dominated by high-frequency fluctuations, the use of only 1×1 Wilson loops (the four plaquettes of the clover, fig. 3.3) in the definition of $G_{\mu\nu}$ could be a problem. In the reference the field strength tensor was computed including loops up to size three. This perhaps allows to remove low distance noise from the lattice, which contribute to an oversampling of the low topological charge values. This low flow time dependence of the susceptibility on the definition of the field strength tensor should be investigated more.

From fig. 7.9 one can infer that the continuum limit, to recover the physical value of the topological susceptibility, can be taken by considering the value of the observable over an interval of the smearing radius between $\sqrt{8t_f} = 0.4$ and 0.6 fm. The extrapolation to the continuum has to be performed by using the dimensionless quantity $(a/r_0)^2$, with a being the lattice spacing and $r_0 = 0.5$ fm the Sommer parameter [25]. The ratio is taken squared in order to match the lattice units of the quantity that is being extrapolated, the susceptibility in this case (which is

proportional to an energy). In fig. 7.10 the continuum limit of the topological susceptibility is taken using the values found for $\sqrt{8t_f} = 0.5$ fm.

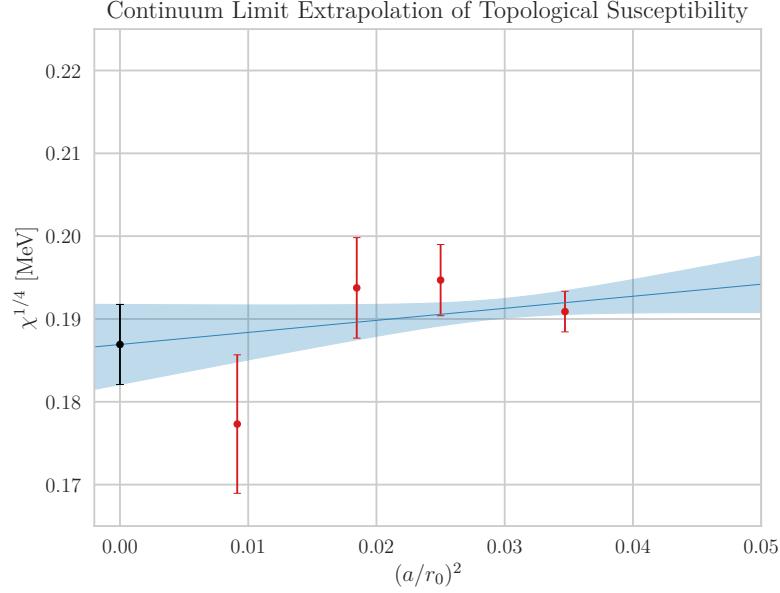


Figure 7.10: Continuum limit extrapolation of $\chi^{\frac{1}{4}}$ from the values at $\sqrt{8t_f} = 0.5$ fm. The fit is performed using the dimensionless quantity $(a/r_0)^2$ as a reference. The black point is the extrapolated point.

The value we obtain for the extrapolated topological susceptibility is:

$$\chi^{\frac{1}{4}} = 186.9(4.9) \text{ MeV}. \quad (7.2)$$

This is compatible with the value in [43] of $\chi_{ref}^{\frac{1}{4}} = 195.9(4.9)$ MeV. **PROBLEMS; B=6.0 POINT; MATHIAS**

As an additional result we can use this extrapolation to compute the mass of the η' meson taken from the Witten-Veneziano formula [19]:

$$\chi = \frac{F_\pi^2 m_{\eta'}^2}{2N_{flavors}}. \quad (7.3)$$

Using as inputs $F_\pi = 92$ MeV, the pion decay constant, and having $N_{flavors} = 3$, we find the mass of the η' meson to be:

$$m_{\eta'}^{(lattice)} = \sqrt{\frac{\chi^{\frac{1}{4}} 2N_{flavors}}{F_\pi^2}} = 921(16) \text{ MeV}, \quad (7.4)$$

which we compare with an experimental value $m_{\eta'}^{(exp)} = 957.78(6)$ MeV, so we are off by more than 2σ , but probably the result is largely influenced by the data point obtained from $\beta = 6.45$, which has less statistics (only 250 configurations) and larger uncertainty.

Chapter 8

Running Coupling and Scale Fixing

In section 4.3 the perturbative behavior of the expectation value of $t_f^2 \langle E \rangle$ was introduced. In this chapter the same quantity is studied from the data generated from our ensembles of lattice gauge field configurations, for which the code-base was developed. The matching between the perturbative and lattice results will also be discussed.

8.1 Discretization Effects on $t_f^2 \langle E \rangle$

Plotting the data for the average value of the energy as a function of the dimensionless quantity t_f/r_0^2 ($r_0 = 0.5$ is the Sommer parameter, described in section 4.3.1 and has unit of length) many interesting properties can be observed. Firstly, the data suggests that for $t_f > 0.05r_0$ is lattice spacing independent, and so coupling independent since they are related. This allows, as it was suggested by [1] that this quantity can be used to set the reference scale of the lattice, similarly to r_0 . By selecting a value for $t_f^2 \langle E \rangle$ one can create a reference scale. It was proposed to use the value of 0.3, which from the plot we can see that is in the region of best matching between the different data series.

In an operative way one finds the value of the flow time t_0 for which:

$$t_0 \langle E \rangle = 0.3 \quad (8.1)$$

This value, in this work, is found by selecting the 20 points, for each data series corresponding to a different β value, and fitting a straight line through them. By inverse regression the value of t_0 and its uncertainty are found.

We can then check the validity of the assumptions, plot the analog of fig. 4.3 in which the ratio of $\sqrt{8t_0}/r_0$, that is the rate of the tentative scale based on the gradient flow and the Sommer parameter. If this ratio is independent from the lattice spacing then $\sqrt{8t_0}$ can be used as a reference scale r_0 . With this procedure we fix a value of $\sqrt{8t_0}/r_0$ to be:

$$\sqrt{8t_0}/r_0 = 0.9499(8) \quad (8.2)$$

which is very close to the value found in [44] of $(\sqrt{8t_0}/r_0)_{ref} = 0.941(7)$.

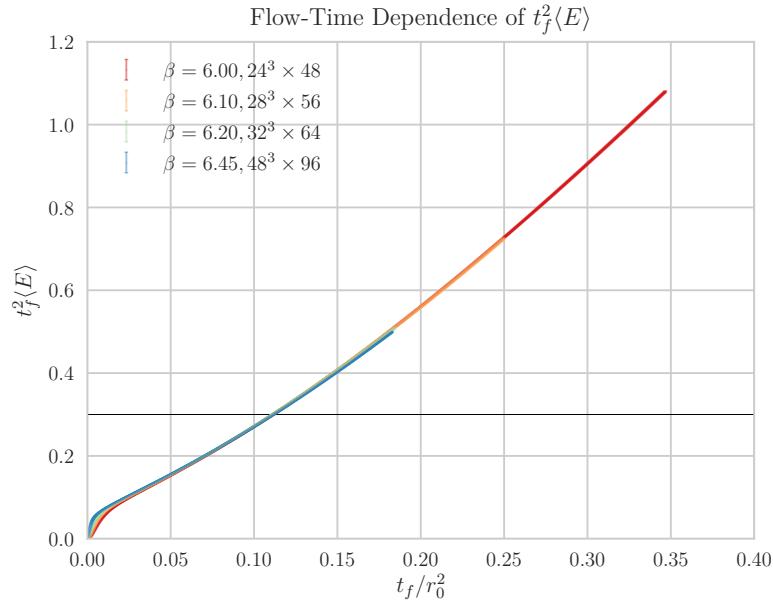


Figure 8.1: $t_f^2 \langle E \rangle$ computed on the four ensembles that were generated (table 5.1) as a function of the flow time in units of r_0^2 . The solid line is for $t_f^2 \langle E \rangle = 0.3$

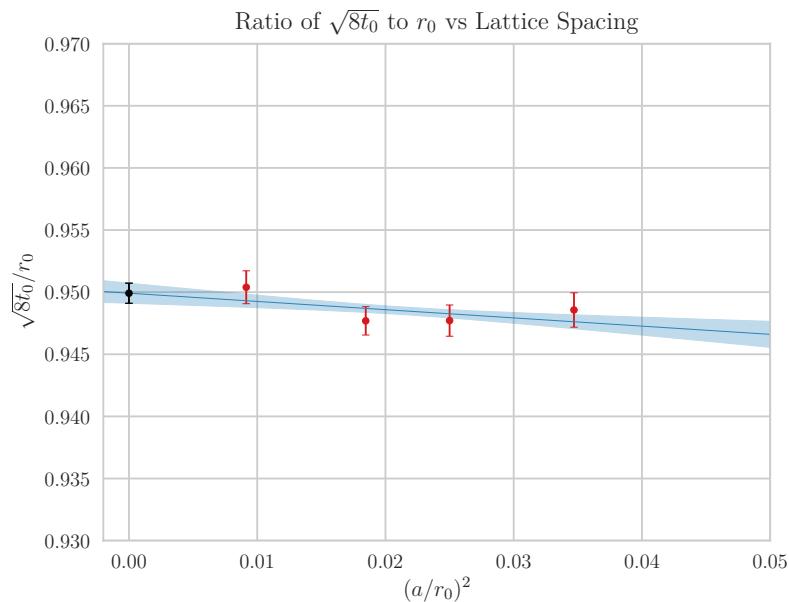


Figure 8.2: Continuum limit extrapolation of the ratio $\sqrt{8t_0}/r_0$. The solid line is the linear fit of the data points, each representing a different β value. The errorband is the 1σ confidence interval and the black point is the extrapolated continuum point.

8.2 Matching Perturbative and Lattice Results

When focusing on the low flow time section of the $t_f^2 \langle E \rangle$ plot, one can study the matching between the perturbative expansion of the observable, performed in [1], and the lattice results that we generated.

The greatest challenge however is that the flow time interval in which this matching happens is unknown. Certainly for large t_f perturbation theory fails to describe the system: any perturbative expansion in $\alpha_s(q)$ becomes meaningless, as $q = 1/\sqrt{8t_f}$ becomes small the coupling grows and approaches one. For small flow times the lattice results become unreliable, in the region where the smearing radius is much smaller than the lattice spacing; this is clear by looking at a zoomed version of fig. 8.1. Both of these two bounds, no matter how intuitive they appear, are also not clearly defined and so no unique procedure can be found to constrain the matching region.

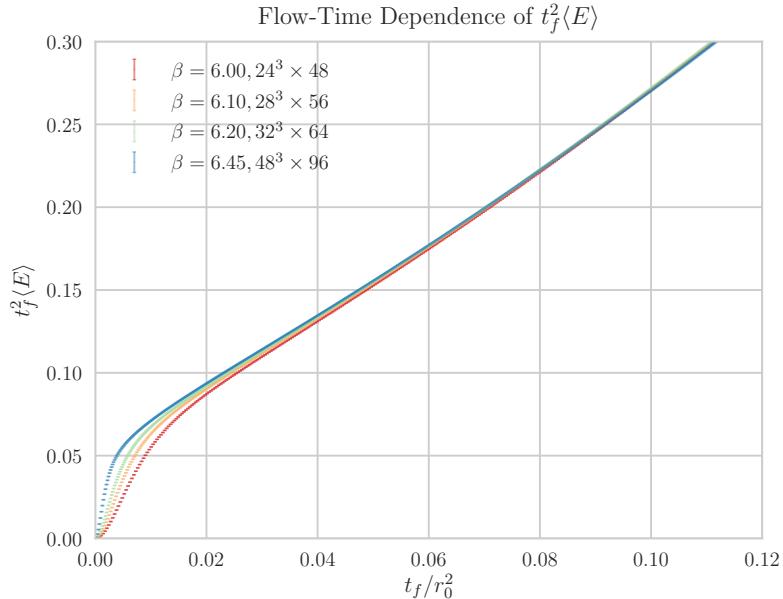


Figure 8.3: Detail of fig. 8.1 for small flow times.

8.2.1 Strategy for the Estimation of the Scale Energy Λ

The problem we are interested to solve is to extract the value of the scale energy of Yang-Mills theory from the lattice data. The only link between the two is eq. (4.21). If the equation is rewritten as:

$$t_f^2 \langle E \rangle = \frac{3}{4\pi} \alpha_s(q) [1 + k_1 \alpha_s(q) + \mathcal{O}(\alpha_s^2)], \quad k_1 = 1.0978 \quad (8.3)$$

one can compute the left hand side on the lattice and fit it to the analytic expression on the right hand side from perturbation theory.

As mentioned earlier the problem is to determine matching region, which now affects the fit range. This must be done carefully in order to prevent possible biases and to assess the error on Λ properly. The first thing to do is to reconstruct the continuum limit curve for $t_f^2\langle E \rangle$. This has been done by selecting linearly spaced values of t_f/r_0^2 from 0.005 to 0.085 with spacing of $5 \cdot 10^{-4}$. For each of the β values, the ten closest points around every value of the linear space were used to fit a straight line and interpolate the value of $t_f^2\langle E \rangle$ exactly on the point. An example for one of those points is shown in fig. 8.4

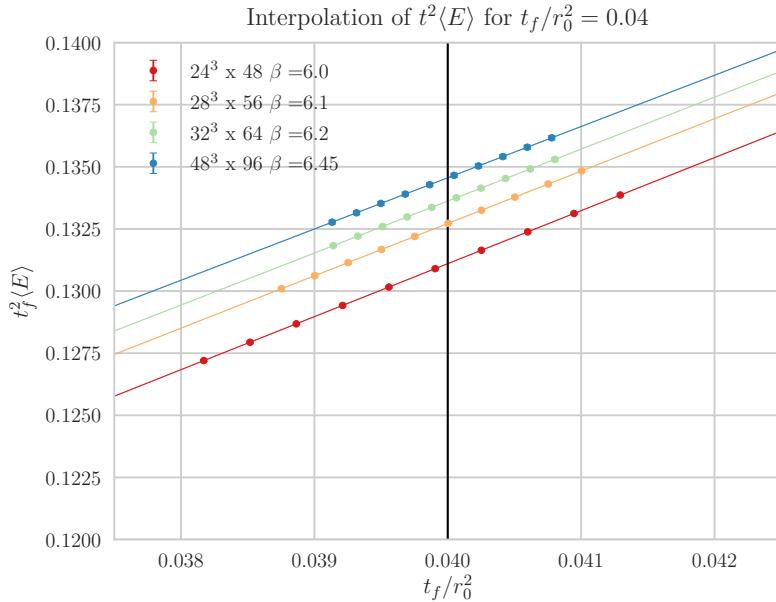


Figure 8.4: Interpolation of the lattice data for $t_f^2\langle E \rangle$ around the point $t_f/r_0^2 = 0.04$. The solid color lines represent the linear fits on the data for one β value each. The data points are the ten closest points to the desired value, represented by the black solid line.

With the interpolated data the continuum limit is taken for each value of the linear space by extrapolating the interpolated points to zero, fig. 8.5

Once this procedure is performed for all points one can plot the data series for the continuum limit. In fig. 8.6 it is plotted together with the data series from the lattice. Notice that the data is plotted against $q = \hbar c / \sqrt{8t_f}$, which is an energy. As a rough approximation, neglecting the $k_1 \alpha_s^2$ term, the continuum limit plot should be proportional to the running coupling.

The estimation of the momentum scale Λ can now be performed on the continuum data. By plugging eq. (4.23), using the values of eq. (4.25) with $N_f = 0$, into eq. (8.3) one can fit the numerical data. However, the range in which to perform the fit is unknown and there is some freedom in the choice of the order of the $\alpha_s(q)$. Under such circumstances, in order to get an unbiased result for Λ a method inspired by the analysis of nucleon masses found in [45] was applied. In that paper, multiple definitions of the nucleon masses were used, but no one could be a priori be preferred over the others, and fits to multiple data with different ranges were performed; a situation somewhat resembling ours.

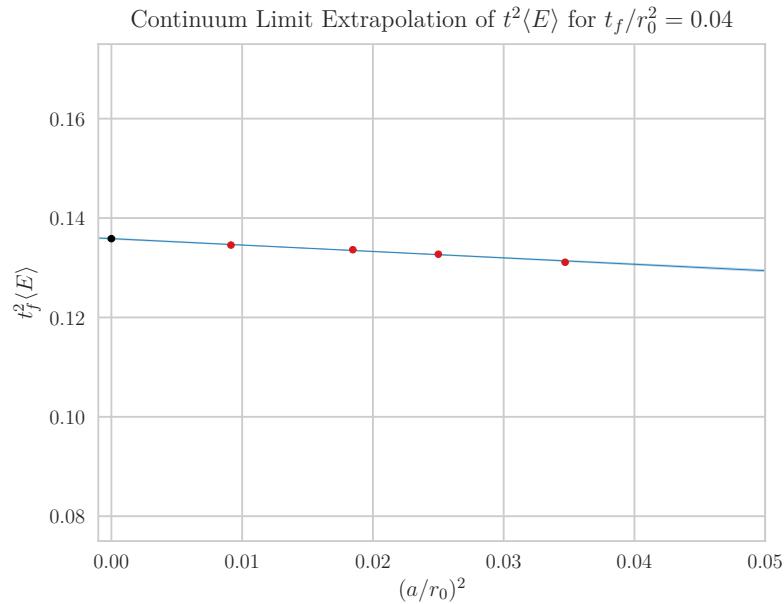


Figure 8.5: Extrapolation of the continuum limit for $t_f^2 \langle E \rangle$ around the point $t_f/r_0^2 = 0.04$. The data points are the result of the interpolation procedure in fig. 8.4.

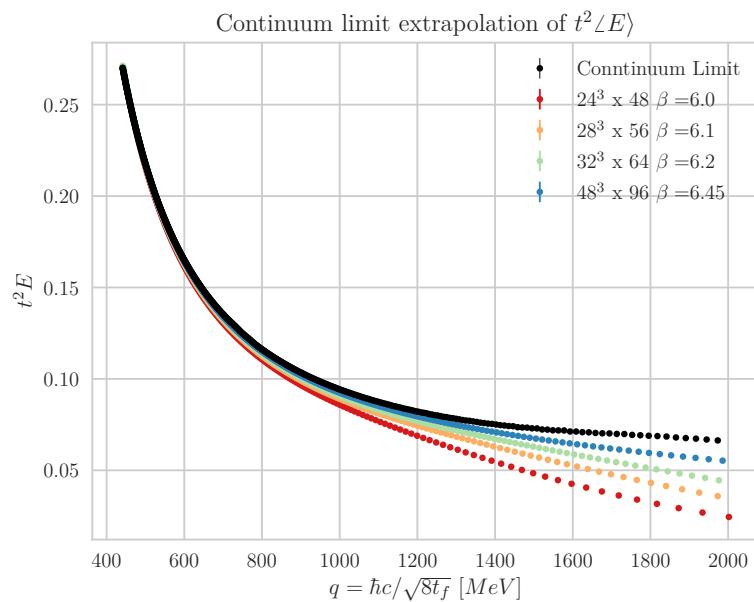


Figure 8.6: Data for $t_f^2 \langle E \rangle$ from lattice observables for the ensembles in table 6.1 and the extrapolated continuum plotted against the energy scale $q = \hbar c / \sqrt{8t_f}$.

The key idea is to effectively perform fits between all possible fit ranges, possibly using also multiple data series if it makes sense to, and construct a histogram of the distribution of the fit parameter. The points in the distribution are then weighted with a quantity that represents the goodness of the fit and the weighted median of the points is taken as the estimator for the fit parameter.

8.3 Results for the Momentum Scale Λ

As outlined in the previous section, the aim is to estimate the momentum scale Λ by performing multiple fits of eq. (8.3). In total four different values for the momentum scale have been calculated, one for every order, from 1 to 4, of loop corrections in $\alpha_s(q)$. The difference between the various functions for the coupling are shown in fig. 8.7.

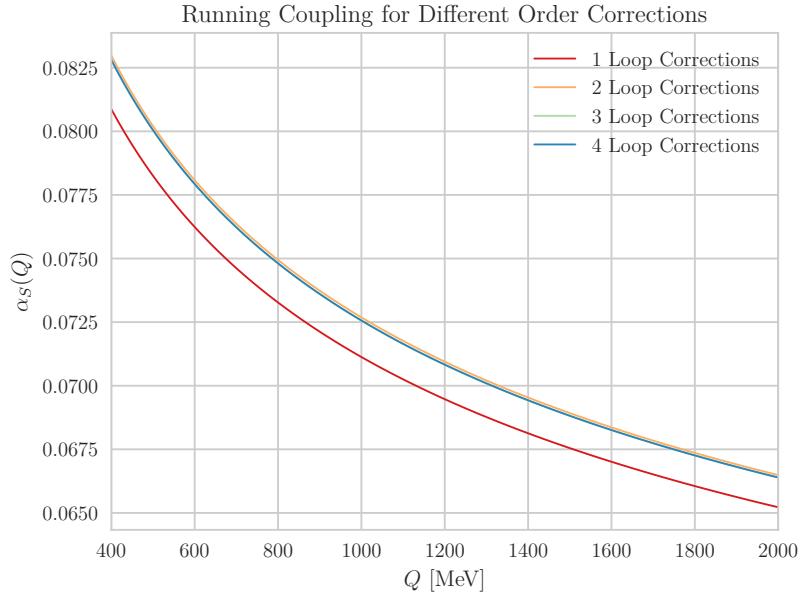


Figure 8.7: Running coupling $\alpha_s(Q)$ of pure gauge Yang-Mills theory for different loop order corrections.

The fit ranges were chosen in terms of the flow time: from $t_f/r_0^2 = 0.005$ to 0.085, of all possible lengths with minimum size 0.005. In total this gave around 3000 different fit ranges for every order of α_s . In fig. 8.8 the distributions of the results of the fits are shown, both unweighted and weighted. The choice of the weight, which has to depend on the goodness of the fit, was $1/\tilde{\chi}^2$, where $\tilde{\chi}^2$ is the reduced chi.square function:

$$\tilde{\chi}^2 = \frac{1}{\nu} \sum_i \frac{(y_i - f(x_i))^2}{\sigma_i^2} \quad (8.4)$$

where $f(x)$ is the function to be fitted (eq. (8.3) in this case), y_i are the data to be modeled (the continuum limit of $t_f^2(E)$) which lie at coordinates x_i and have variance σ_i^2 .

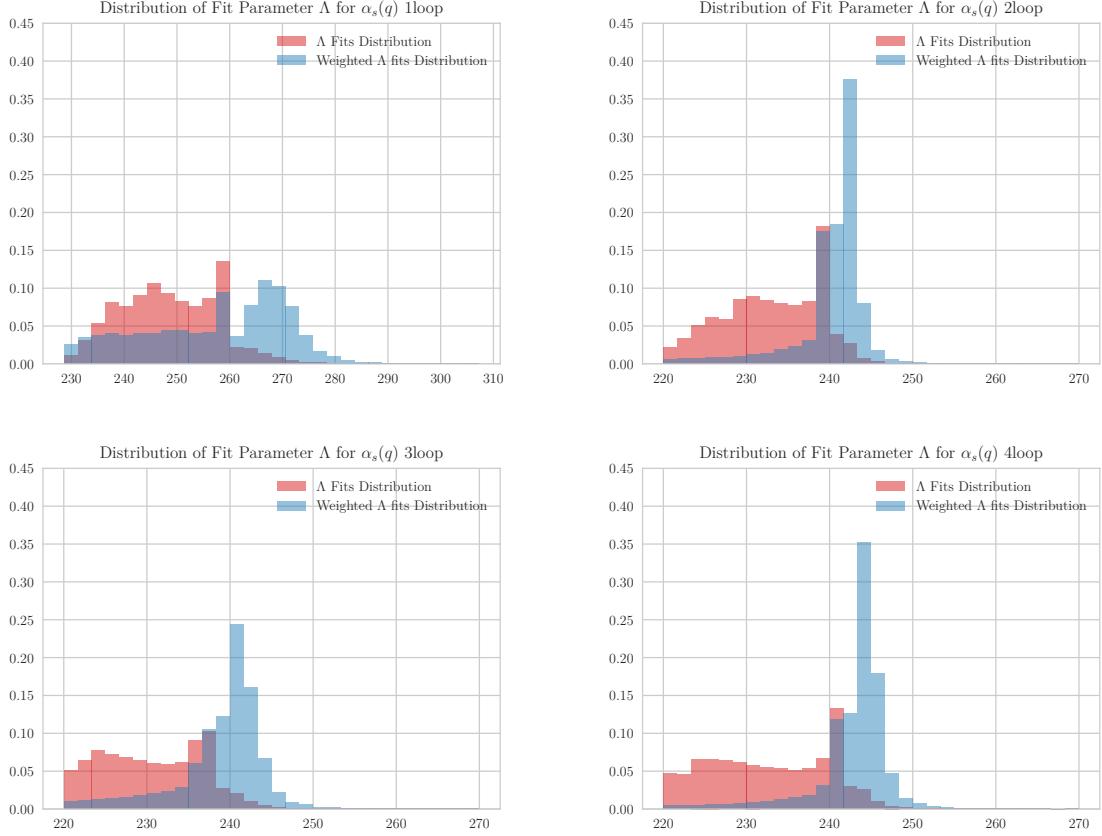


Figure 8.8: Distributions of the fit parameters, Λ , obtained by all possible fit ranges between $t_f/r_0^2 = 0.005$ to 0.085 of minimum size 0.005 . The four different plots are for the four different orders of corrections for $\alpha_s(q)$. Data for the distribution weighted with $1/\chi^2$, normalized to the number of fits, is also reported.

The fact that the weighted distributions have a peak means that there is a certain region in the energy spectrum where the function in eq. (8.3) matches the data better. This is expected since we know that for low energies perturbation theory does not hold anymore, so the function does not represent the data correctly. At high energies on the other hand, by looking at fig. 8.3 for example, the discretization effects introduced by the lattice method become relevant, so the continuum limit extrapolation that we made is unreliable in that region.

From the plot of the 1-loop corrected function, one can see that there is no sharp peak in the weighted distribution, compared to the others. The difference in the behavior is not surprising if one looks at fig. 8.7 and considering that the higher order functions for $\alpha_s(q)$ are all close together in the energy range considered.

The estimation of the Λ momentum scale is taken by considering the weighted median of the data, which is defined as the 50% weighted percentile, the element x_k of an ordered set of values

x_1, x_2, \dots, x_N that satisfies:

$$\sum_{i=0}^{k-1} w_i \leq \frac{1}{2} \quad \text{and} \quad \sum_{i=k+1}^N w_i \geq \frac{1}{2} \quad (8.5)$$

The systematic uncertainty has been taken as the 68% confidence interval of the weighted median distribution, as it was done in [45]. To estimate the statistical error, the whole procedure has been applied to 500 bootstrap samples of the lattice data, constructing the continuum limit limit for each sample and performing fits on all fit ranges for sample. The results are shown in section 8.3.

Fit Function Order	Λ [MeV]	σ_{sys} [MeV]	σ_{stat} [MeV]
1-loop	258.9	23.5	6.3
2-loop	241.6	9.4	0.7
3-loop	240.1	14.0	1.2
4-loop	243.9	10.2	0.7

Table 8.1: Final results for the momentum scale parameter from the distribution of the fits in fig. 8.8.

The statistical error is consistently smaller than the systematic one proves that the method for the extrapolation of the continuum limit data is solid, but also implies that almost all the uncertainty is given by the method used to estimate Λ .

This value can now be compared with ones found in literature for the same quantity using different methods. In particular we compare it with the one found in [2] obtained using a recursive finite-size technique, which is reported as $\Lambda_{ref} = 238(18)$ MeV. The values found in section 8.3 agree very well with this value.

One last result to show is the plot of $t_f^2 E(q)$ using our values of Λ as input. In fig. 8.9 it is plotted, order by order, together with the continuum limit data extrapolated from the lattice. We can see that good agreement is found between $q \approx 700$ and 1400 MeV, or in flow time between $t_f/r_0^2 \approx 0.01$ to $t_f/r_0^2 \approx 0.045$

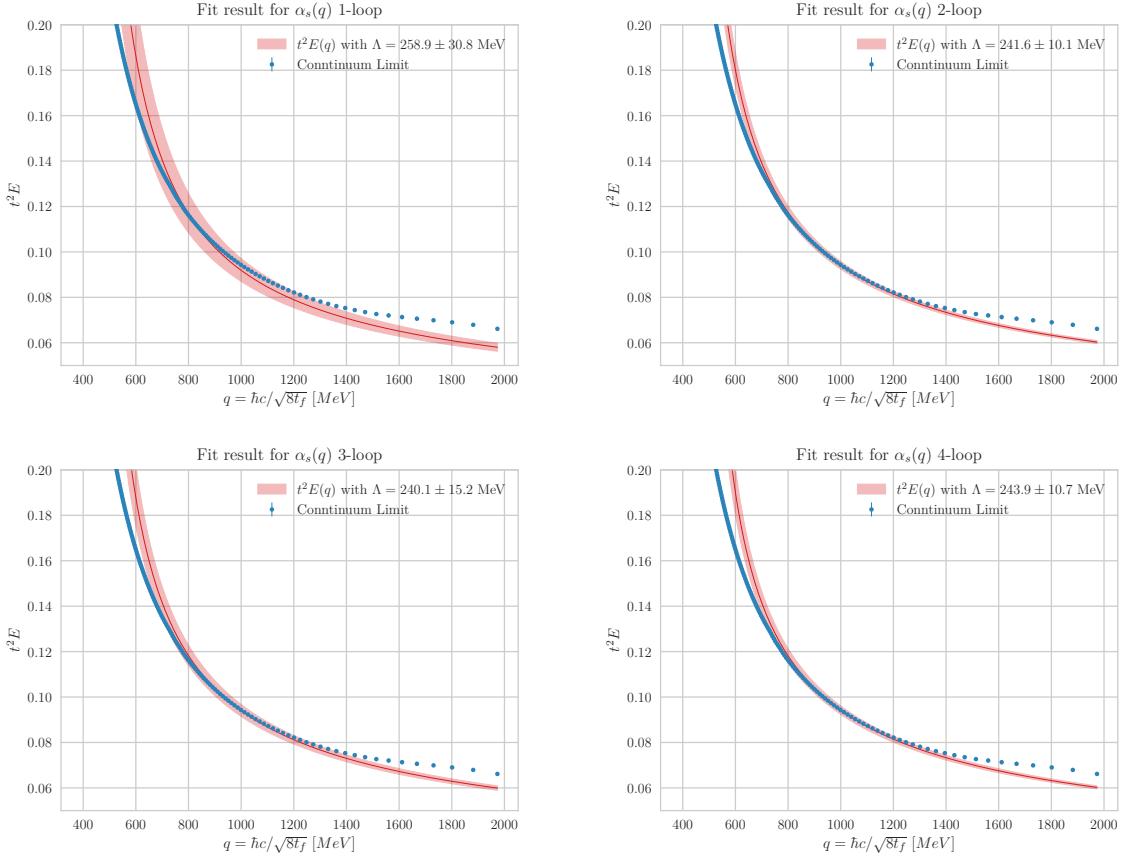


Figure 8.9: Plot of $t_f^2 E(q)$, eq. (8.3), using the values of the momentum scale found in section 8.3. The error band is the sum of the statistical and systematic errors from the same table. The data series in blue is the continuum limit extrapolated from the lattice.

Part IV

Conclusions

Chapter 9

Summary and Outlook

This work had two main goals: develop a completely new framework to perform lattice calculations in Yang-Mills theory, in particular using the gradient flow method; and to perform a calculation of the momentum scale Λ by matching the perturbative analysis of the energy density as a function of flow time and the continuum limit extrapolation of lattice results.

The main features of Quantum Chromodynamics and of how to discretize it on a lattice have been discussed. In chapter 5 the details for the implementation of the Metropolis algorithm and the numerical integration of the flow equation have been presented, together with the structure of the source-code that has been developed.

For the computational part of the work, the benchmarks and the performing scaling properties the program for both the generation of gauge field configurations and the gradient flow are a strong indication of the goodness of the implementation. [MORE ON SCALING...](#)

The calibration of the parameters of the Metropolis algorithm for generating gauge field configuration was performed by looking at the integrated autocorrelation time τ_{int} of the topological charge at different flow times. It was found that, as expected, the autocorrelation for smaller lattice spacings grows non-linearly. Consequently the natural solution has been to increase the distance in Monte Carlo time of the measurements of the observables, bu this meant a much larger computational cost so the ensemble size was also reduced. The final results for τ_{int} for the worse case ($\beta = 6.45$ and $a = 0.0478$ fm) are contained. The energy density showed no particular issues in this regard.

When looking at the topological charge, computed on all of the four ensembles that were generated, it was found to be always consistent with 0, but given the very large volumes and the relatively small ensembles that were considered the uncertainties are rather large. The analysis of the topological susceptibility, on the other hand, proved to be a good estimate, the value we found is $\chi^{\frac{1}{4}} = 186.9(4.9)$ MeV, in good agreement with what reported in [1, 43] and also, through the Witten-Veneziano formula with the experimental result from the mass of the η' meson.

Lastly, the main result of this work is the estimate of the momentum scale of Yang-Mills theory,

starting from the lattice results for $t_f^2 \langle E \rangle$, taking the continuum limit function and iteratively fitting the perturbative expansion function. The numerical procedure that has been applied produced results, depending on the order of loop correction for the renormalized coupling, that are consistent with other methods found in literature. The fact that the total uncertainty of our results is lower for the high order definitions of $\alpha_s(q)$ could mean that this method is valid and that is more precise, and one could add much simpler, than others. However it could also be the result of the underestimation of the systematic error, so further work is needed.

9.1 Future Work

Most of the possible future work is related to the code-base for lattice calculations. Planned short-term improvements include: higher order calculations of gluonic observable (that is larger Wilson loops), improved actions such as the Conjugate Wilson Action or the Symanzik or Ywasaki improved actions, different sampling algorithms, like the Hybrid Monte Carlo algorithm. Another possible, but probably more time-consuming, new development is the introduction of dynamical fermions, allowing for full QCD calculations.

On a lower level, improvements in the handling of parallelization are planned. In particular some initial steps have been done in the creation of a low level library for SIMD calculations, using processor specific instruction sets. Hybrid memory parallelization, especially using the new features offered by the MPI 3.0 standard, is being considered as a possible feature.

Regarding the estimation of the momentum scale from lattice calculations, we are planning to test in the very near future the method that has been used in this work to data of the energy density of field configurations with $2 + 1$ dynamical fermions computed with the gradient flow method. The goal is to verify if the approach holds and if a value for Λ_{QCD} can be obtained. The most obvious future addition would be to perform the calculation on other ensembles of gauge field configurations, to check the stability with respect to lattice spacing, finite volume effects and, more in general, to improve the continuum limit extrapolation of the data. A more detailed analysis of the systematic error estimation should also be performed, to finally confirm the results presented in this thesis.

Part V

Appendices

Appendix A

Statistical Tools

The determination of the uncertainty of the results of a numerical experiment is Extremely important. In cases of small datasets, as in this work or in Lattice QCD in general, resampling methods are a popular solution for improving the estimate of the uncertainty. Another very important statistical property that is to be considered is the autocorrelation, introduced by the use of Markov chains.

A.1 Resampling: Bootstrap and Jackknife

Given a set of N measurements of an observable O , a first estimate of the expectation value and of the uncertainty are the mean and the standard deviation of the data:

$$\bar{O} = \frac{1}{N} \sum_{i=1}^N O_i, \quad \sigma_O = \sqrt{\frac{1}{N} \sum_{i=1}^N (O_i - \bar{O})^2} \quad (\text{A.1})$$

where O_i are the measured values of the observable. Resampling methods use the data O_i to construct multiple data series, and the average statistical properties of these series are used to estimate the properties of the original set. Both the bootstrap and the jackknife methods assume uncorrelated data. The observable O can also be a derived quantity, such as a fit on some raw data.

A.1.1 Bootstrap

The Bootstrap method consists in building N_b bootstrap samples by choosing N points at random from the the original values (repetition are allowed). The average of a bootstrap sample is then:

$$\hat{O}_b = \frac{1}{N} \sum_{i=1}^N O_{r(N)} \quad (\text{A.2})$$

where $r(N)$ stands for a random number between 1 and N . One then has N_b such averages and constructs:

$$\tilde{O} = \frac{1}{N_b} \sum_{b=1}^{N_b} \hat{O}_b, \quad \tilde{\sigma}_{\tilde{O}} = \sqrt{\frac{1}{N_b} \sum_{b=1}^{N_b} (\hat{O}_b - \tilde{O})^2} \quad (\text{A.3})$$

The expectation value of the observable is then $\langle O \rangle = \tilde{O}$ and its associated uncertainty is $\tilde{\sigma}_{\tilde{O}}$.

A.1.2 Jackknife

The Jackknife sample are not constructed at random, as is the case for bootstrap, but in a systematic manner. One creates N jackknife samples by removing for each of them the j th element of the data:

$$\hat{O}_j = \frac{1}{N-1} \sum_{i \neq j}^N O_i \quad (\text{A.4})$$

One then computes:

$$\tilde{O} = \frac{1}{N} \sum_{j=1}^N \hat{O}_j, \quad \tilde{\sigma}_{\tilde{O}} = \sqrt{\frac{N-1}{N} \sum_{j=1}^N (\hat{O}_j - \tilde{O})^2} \quad (\text{A.5})$$

The final estimator for the expectation value is $\langle O \rangle = \tilde{O}$ and the uncertainty is $\tilde{\sigma}_{\tilde{O}}$.

Both of these resampling methods have been implemented and used; they were found to be consistent with each other. The jackknife is deterministic, the bootstrap, for sufficiently large N_b (usually ≈ 500 are sufficient) had error estimates compatible with jackknife and of the same order of magnitude. The error estimate when using resampling compared to the standard deviation is smaller by orders of magnitude.

A.2 Autocorrelation

The data generated via a Markov chain Monte Carlo method is always affected by autocorrelation because the data points intrinsically depend on each other. One can tune the algorithm to reduce the autocorrelation time until it becomes negligible, but this is not the case for Lattice calculations because of the computational cost involved.

We followed the procedures in [46] to get an algorithm to estimate the integrated autocorrelation time and the error correction factor due to autocorrelation.

The autocorrelation function of an ordered set variables $\{x_1, x_2, \dots, x_N\}$ with expectation value \bar{x} is defined as:

$$\Gamma(t) = \Gamma(-t) = \langle (x_i - \bar{x})(x_{i+t} - \bar{x}) \rangle \approx \frac{1}{N-t} \sum_{i=1}^{N-t} (x_i - \bar{x})(x_{i+t} - \bar{x}) \quad (\text{A.6})$$

where t is the “lag” between two points. The integrated autocorrelation time is given by:

$$\tau_{int} = \frac{1}{2} \sum_{t=1}^{\infty} \frac{\Gamma(t)}{\Gamma(0)} = \frac{1}{2} \sum_{t=1}^{\infty} \rho(t) \quad (\text{A.7})$$

For sufficiently large N , the true variance of the data can be expressed as:

$$\tilde{\sigma}^2 = 2\tau_{int}\sigma^2 \quad (\text{A.8})$$

This result is very neat, because it implies that in one can calculate the integrated autocorrelation time correctly, the inclusion of the effects on the uncertainty estimate is trivial. In order to truncate the infinite summation in eq. (A.7) one can look at the deviation squared of $\rho(t)$, which as suggested by Sokal (**CITATION NEEDED**) can be written as:

$$\langle \delta\rho(t)^2 \rangle \approx \frac{1}{N} \sum_{k=1}^{\infty} [\rho(k+t) + \rho(k-t) - 2\rho(k)\rho(t)]^2 \quad (\text{A.9})$$

all these terms, for a sufficiently large value of k should all vanish, hence one can choose a cutoff Λ and truncate the sum up to $t + \Lambda$. The integrated autocorrelation time, if the deviations of $\rho(t)$ become small, plateaus, so one can truncate the sum in eq. (A.7) as well:

$$\tau_{int} = \frac{1}{2} \sum_{t=1}^W \rho(t) \quad (\text{A.10})$$

where W is the first lag t for which $\rho(t) < \sqrt{\langle \delta\rho(t)^2 \rangle}$, when the contribution to the integration of τ_{int} from that lag become smaller than the deviation of that same lag.

The integrated autocorrelation time has been computed for all quantities presented in this work and statistical uncertainties associated with results have all been corrected via eq. (A.8)

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