

Tuning NRQCD b-quarks for Bottomonium Spectroscopy Using Lattice QCD

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Abstract

Quantum Chromodynamics (QCD) is an integral component of the Standard Model (SM). At low energies QCD cannot be solved perturbatively, however it is possible to find solutions by coarse-graining and computing on a lattice. Bottomonium mesons can be studied using Non-Relativistic QCD (NRQCD) as the b-quarks are much heavier in mass compared to the lighter quarks and need super fine lattices to simulate them relativistically which are not easily available. This project will involve modifying an existing NRQCD codebase to support an anisotropic lattice, with finer graining in the time component. The b-quark mass m_b will be tuned, along with the degree of lattice anisotropy, in order to produce spectra for both vector and scalar B -mesons. These will be compared using established lattice QCD statistical techniques (`lsqfit` and `corrfit`) with existing spectra. (TODO: Correct this *no longer studying anisotropy, nor modifying code.*)

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1 Introduction

1.1 Standard Model

The Standard Model is currently the most accurate model of particle physics. It describes fermions: quarks (q), leptons (l), and their antiparticles, denoted with an overline (\bar{q}). It also describes the gauge bosons which mediate the strong, weak, and electromagnetic (EM) fields.

There are 6 flavours of quarks, in 3 different generations. Quarks are coupled to the strong, weak, and EM interactions. The leptons, also in 3 generations, are coupled to the EM field if they are charged, and to the weak force. This project is to study the spectra of bottomonium mesons. These are hadrons composed of a bottom quark and any antiquark (or an antiparticle of this). The weak and EM interactions are completely neglected, leaving only the strong interaction which is governed by the theory of Quantum Chromodynamics (QCD).

1.2 Quantum Chromodynamics

QCD is a $SU(3)$ gauge field theory governing the interactions of quarks and gluons. As quarks are fermions, they are described by a Dirac lagrangian coupled to an external gauge field. For a quark field annihilation operator ψ_f , Dirac field creation operator $\bar{\psi} = \psi_f^\dagger \gamma^0$ with quark flavour f , we have[1, 2]:

$$\mathcal{L} = \sum_f \psi_f^\dagger (i \not{D} - m_f) \psi_f - \frac{1}{4} F_{\mu\nu}^a F_a^{\mu\nu} \quad (1)$$

In this Lagrangian we have a gauge covariant derivative $\not{D} = \not{\partial} + ig_s \not{A} \lambda_a / 2$. Here and in the rest of this section, we leave the indices referring to colour implicit. However to be clear, the quarks carry a conserved *colour charge* analogous to the electric charge in QED. There are three separate *colours* which are often called *red*, *green*, and *blue*¹. The colour of a quark is represented by a colour spinor, which is the space acted on by the λ_a . It is these indices, referring to the matrix multiplication, that are omitted.

The other important quantity is the gluon field strength tensor:

$$F_{\mu\nu}^a = \partial_\mu A_\nu^a - \partial_\nu A_\mu^a - g_s f_{abc} A_\mu^b A_\nu^c \quad (2)$$

Here the latin indices label the 8 generators² ($t_a = \lambda_a / 2$) of the QCD $\mathfrak{su}(3)$ Lie algebra with (totally antisymmetric) structure constants $[t_a, t_c] = i f_{abc} t_c$. These generators are considered to be linearly independent colour charges of gluons. A_μ^a are the respective 4-potentials, which are coupled to the quarks.

Unlike in QED, QCD allows for self-interaction of the gauge bosons. As a result, QCD is known as an asymptotically-free theory: as interactions are taken to higher energies, the forces involved become weaker. This leads to colour confinement, where only bound states with zero overall colour charge can be isolated. Traditional perturbation theory

¹Although, the choice of basis is arbitrary and can be transformed by any matrix $U \in SU(3)$.

² λ_a are the Gell-Mann matrices, which form a basis of $SU(3)$.

expands a system at low energy, but in the case of QCD this would give divergent forces which would invalidate such a method. Instead, a common approach is solving systems on a lattice.

1.3 Lattice QCD

Lattice QCD is a method to solve QCD systems by placing the fields on a spacetime lattice. To see how this is done we first consider the path integral formulation of QFT. To compute the expectation of an operator, we take a functional integral of the Lagrangian over all possible field configurations[3, 2]

$$\langle \mathcal{O} \rangle = \frac{1}{\mathcal{Z}} \int \mathcal{D}\psi \mathcal{O} \exp\left(i \int d^4x \mathcal{L}(\psi, \partial\psi)\right), \quad (3)$$

$$\mathcal{Z} = \int \mathcal{D}\psi \exp\left(i \int d^4x \mathcal{L}(\psi, \partial\psi)\right), \quad (4)$$

where \mathcal{Z} is the partition function. This is named in analogy with the statistical partition function, which can be obtained by performing a Wick rotation $t \mapsto it$.

Clearly it is not usually feasible to integrate over an infinite-dimensional space of functions, so we use a technique called regularisation. To do this we replace the $4D$ continuum integral by a spacetime lattice. Often an isotropic lattice is chosen (with lattice spacing a), however the main aim of this project is to use an anisotropic lattice the spatial and temporal lattice spacings differ.

We define the quark fields[2] $\psi(x)$ only on lattice points x_μ , and the gluon fields $U_\delta(x) = U(x, x + \hat{\delta})$ defined on the links between the lattice points. The links are specified as written, by a point x^μ and a unit vector $\hat{\delta}^\mu$ specifying the direction along the lattice. The conjugate field $U_\delta^\dagger(x) = U(x + \hat{\delta}, x)$ gives the link from $x + \hat{\delta}$ back to x . Once we have our particle fields, it is possible to replace derivatives with differences in fields at adjacent lattice sites. One can then construct an appropriate gauge invariant action, called the Wilson action.

We hope that by using this lattice model of QCD, we can produce sensible results in the limit of a short lattice spacing. Unfortunately, when the computations are carried out there are a lot of discretisation errors[4]. By a careful choice of discrete derivative, we can remove $\mathcal{O}(a)$ error terms. The $\mathcal{O}(a^2)$ error terms can also be removed by adding and subtracting appropriate counterterms[5]. This leaves us with a discretised theory with errors in $\mathcal{O}(a^3)$, which we hope is small enough that our extrapolation to $a \rightarrow 0$ is accurate.

1.4 NRQCD

In this project we are looking specifically at bottomonium mesons. In a relativistic approach, there are certain constraints placed upon both our lattice spacing and our lattice size. These constraints require a lattice with lattice spacing equal or smaller than 0.45 fm to compute b-quarks relativistically[2]. In practice sometimes these fine lattices are not easily available, or even if they are the computation becomes very expensive. Instead we take a non-relativistic approach. This NRQCD method is justified by the

relatively large mass of the b-quark: we can expand our action in factors of $1/m_b$ and neglect the appropriately higher order terms.

This is the first step in the Foldy-Wouthuysen-Tani (FWT) transformation. Carrying on with the remaining steps gives us our lattice NRQCD, which to first order in v^2 (the first term in the non-relativistic expansion) is the Schrödinger Lagrangian:

$$\mathcal{L}_0 = \psi^\dagger(x) \left(iD_t + \frac{\mathbf{D}^2}{2m_b} \right) \psi(x), \quad (5)$$

where D_t is the time component of the gauge invariant derivative, and \mathbf{D} is the spatial 3-vector component. The expansion to v^4 is required for an interesting theory, and involves terms that interact[6] with the chromoelectric $E_j = F_{0j}$ and chromomagnetic $B_j = -\frac{1}{2}\epsilon_{jkl}F^{kl}$ fields, which is exact to $\mathcal{O}(1/m_b^3)$. In this project, we will in fact use an action which is expanded to $\mathcal{O}(v^6)$. TODO: This section needs updating, we are not only looking at bottomonium

1.5 Relating lattice and physical

TODO: This is for chiral-continuum limit etc. Any mass specified in lattice units can be converted into physical units by $m = (\hbar c)m^{\text{lattice}}/a$.

2 Statistical methods and Data

2.1 Jackknife averaging

Given a large sample of data with completely unknown errors, we often want to compute unbiased estimators for statistical parameters. One such way is to use Jackknife averaging[7]. Consider a sample of n i.i.d. values $x_i \sim F$. Construct n mean estimates $\bar{x}_i = \sum_{j \neq i} \frac{x_j}{n-1}$. The average of these sample means $\bar{x} = \sum_i \frac{\bar{x}_i}{n}$ is precisely the mean of the whole sample, and the error on this estimate is given by the variance $\text{Var}[\bar{x}] = \frac{n-1}{n} \sum_i (\bar{x}_i - \bar{x})^2$ of the distribution of \bar{x}_i . That is, we have an unbiased estimator for the mean and standard deviation of F given by $\bar{x} \pm \sqrt{\text{Var}[\bar{x}]}$. In fact by replacing the means \bar{x}_i above by (almost) any estimator $\bar{\theta}_i$, it is possible to construct a new estimator, using a slightly more laborious method, which has an asymptotically smaller bias[8].

2.2 Bayesian analysis

Using the `corrfit`[9] Python package, it is possible to perform a least-squares regression on our 2-point correlator data. For a given set of priors and a data set, the least-squares regression gives best-fit values for parameters of choice. In the context of this project, we are interested in fitting the amplitudes a_n , a_{on} and the log of the energy differences, $\log(dE_n)$, $\log(dE_{on})$. In particular, the log of energy differences allows us to implicitly force the energy levels of the fit to strictly increase.

For a fit of this type, we expect a good quality fit to have $\chi^2/\text{dof} \approx 1$. Values much larger than this are a poor fit, and values closer to 0 could indicate overfitting. There

Label	Spacing a (fm)	Lattice size (time)	m_l/m_s
Very Coarse	0.1509	48	0.036
Coarse	0.124 04	64	0.2
Fine	0.090 23	96	0.2

Table 1: Parameters of the data used in the main part of this project. Corresponding to sets 1, 5, 6 from table I[10] respectively. The lattice size is the number of discrete time slices that are recorded.

are other measures of goodness-of-fit such as the Gaussian Bayes Factor and the Q value, however their use is unnecessary here and we will continue only using the χ^2 measure.

2.3 Data

Data for this whole project has been provided by Dr. B. Chakraborty. This data has been produced by isotropic lattice NRQCD on a variety of lattice spacings, described in Table 1. Each correlator sample corresponds to a randomised gauge configuration of gluon fields.

By using a finite lattice we have discretisation errors (discussed in Section 1.3), which we can only eliminate by extrapolating from multiple datasets, to the continuum limit. Additionally, for computational reasons it is not possible to use physical light quark masses for all lattice sizes. As a result, we must also extrapolate to the chiral limit, where $m_l = 0$, which is close to the physical value of $m_l \approx 0.02m_s$.

3 Preliminary study

Using Jackknife averaging, an existing dataset[11] was analysed. This is a set of 2-point correlators for a D -meson ($c\bar{q}$). In this dataset there are $n = 489$ gauge configurations, with $T = 96$ time slices. The average over Jackknife averages is shown in Figure 1. From[12] we know that these correlators are expected to follow the following relationship,

$$G(t) = \sum_n a_n^2 (e^{-E_n t} + e^{-E_n (T-t)}) + (-1)^t \sum_n a_{on}^2 (e^{-E_{on} t} + e^{-E_{on} (T-t)}), \quad (6)$$

where E_n gives the energy (or mass) of the n^{th} excited state, and a_n gives its amplitude. In this expression, the reflected $T-t$ terms are due to the periodic boundary conditions, and the E_{on} terms are oscillatory contributions due to the fermion doubling in the staggered quark analysis.

If we assume the oscillatory terms are small then one can look near the middle of the data (shown in Figure 1). Here the only relevant term is the ground state (a_0, E_0). As a result, in this region we should be able to compute an estimate for the ground state effective mass,

$$m_{\text{eff}}(t) = \log \left(\frac{G(t)}{G(t+1)} \right). \quad (7)$$

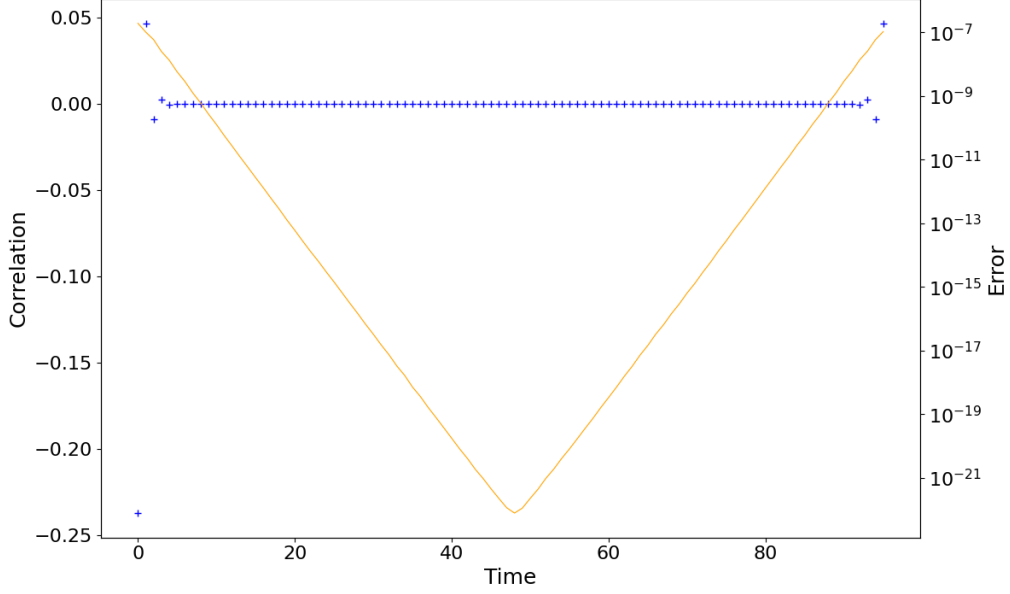


Figure 1: Average 2 point correlator for $n = 489$ gauge configurations on a fine lattice. Computed by using Jackknife averaging. Away from the endpoints correlator values are very small, which is consistent with our assumption that our exponentials decay sufficiently quickly.

When actually computing this value, we must make sure that we do not go past halfway through the data, as the second half (with rising exponentials) will lead to a negative effective mass with the same magnitude. As shown in Figure 2, for the provided (non-Gold D meson on a fine lattice), the effective mass is approximately $m_{\text{eff}} \approx 0.87 \approx 1900 \text{ MeV}$. To compute this value we took the average of the time-dependent values for time steps $t = 24 - 36$, where the data has clearly plateaued but still has small errors.

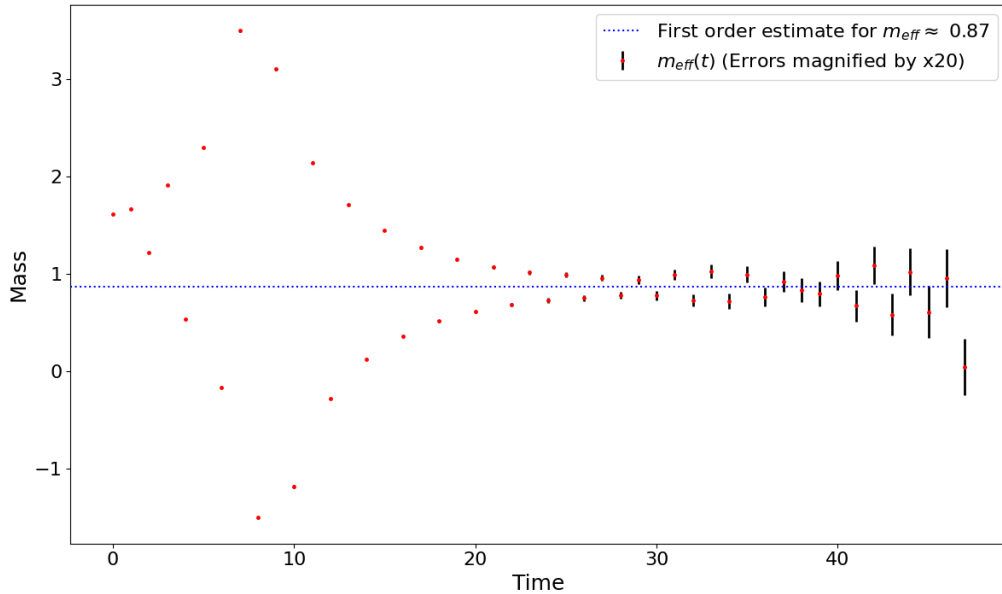


Figure 2: Estimate for m_{eff} using Equation 7.

4 Data Analysis

Throughout this section, all correlator data has been Jackknife averaged to give an estimated Gaussian variable at each time slice. The analysis code is a very heavily modified version of code again provided by Chakraborty, which includes optimised prior values. Throughout this analysis we are fitting up to 6 terms in the correlator (Equation 6). It is not reasonable to extract physical results from most of these terms, however we can certainly look at the ground state of each meson. Taking the most reliable value of meson mass for each of the data sets, it is possible to extrapolate to the chiral and continuum limits to compute final estimates for the neutral K and D meson masses.

During the analysis, there are two main fitting parameters that are varied. t_{\min} tells the fitting software how many time slices to ignore from the start and the end of each set of correlator data. This is useful as there are extra effects happening at these early time scales, and we would like to ignore these for simplicity. The other parameter, n , corresponds to the number of exponentials for which we are providing prior data. Studying energy levels larger than n is completely unreasonable as these are only used to achieve a better fit, and the values are unphysical.

For all datasets that are analysed, we limit the fitting for K mesons to only have $t_{\min} = 3$. For D mesons we vary this as $t_{\min} \in \{3, 4, 5, 6, 7, 9\}$. For both mesons, we also vary $n \in \{1, \dots, 6\}$. In all of the following plots, $n = 4$ was determined to be the most reliable point, so all values given on the plots are evaluated here. For plots showing fitting results, error bars are indicated in red. When a data point (error bar) appears missing it is either too large (small) to be shown on the scale. For convenience, the first 4 data points of the fitting quality plots have also been reproduced on the plots, as the majority of these points are far too large to be shown on the same scale as the rest of the points.

4.1 Fit convergence

The fine dataset corresponds to a lattice spacing $a \approx 0.09$ fm and a light quark mass $m_l \approx 0.2m_s$. For the K meson, it is clear that for any value of $n \geq 4$, the fit has converged (Figure 3³). Inspection of the results for D meson show that the fits tend to converge first for $n = 3$ or 4. Now we turn to the fit quality, where as we can see in Figure 4, the fitting converged well for all t_{\min} and for all $n \geq 4$.

The coarse dataset corresponds to a lattice spacing $a \approx 0.12$ fm and a light quark mass $m_l \approx 0.2m_s$. In this dataset, we can see that the K meson values (Figure 5) actually converge for $n = 3$. More generally, the fit quality (Figure 6) suggests that $n = 3$ or 4 are the earliest parameters that converge, although this is more reliable for higher t_{\min} .

The very coarse dataset corresponds to a lattice spacing $a \approx 0.15$ fm and a light quark mass $m_l \approx 0.036m_s$. Here we see that the K meson fit (Figure 7) appears to converge even for $n = 2$. The fit quality plots (Figure 8) show a different result however. Even at higher t_{\min} , the χ^2 parameter does not reach a reasonable value until at least $n = 3$.

³Plots of this form have been produced for all combinations of dataset and t_{\min} , however where they do not add to the discussion, they have been omitted.

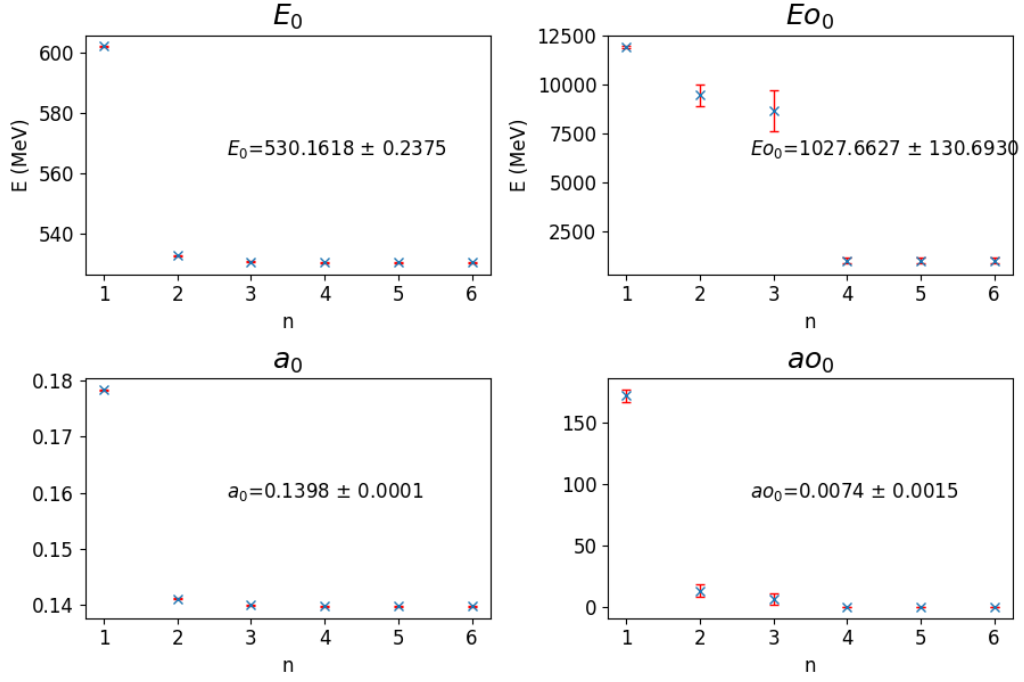


Figure 3: K meson fitting results on a fine lattice.

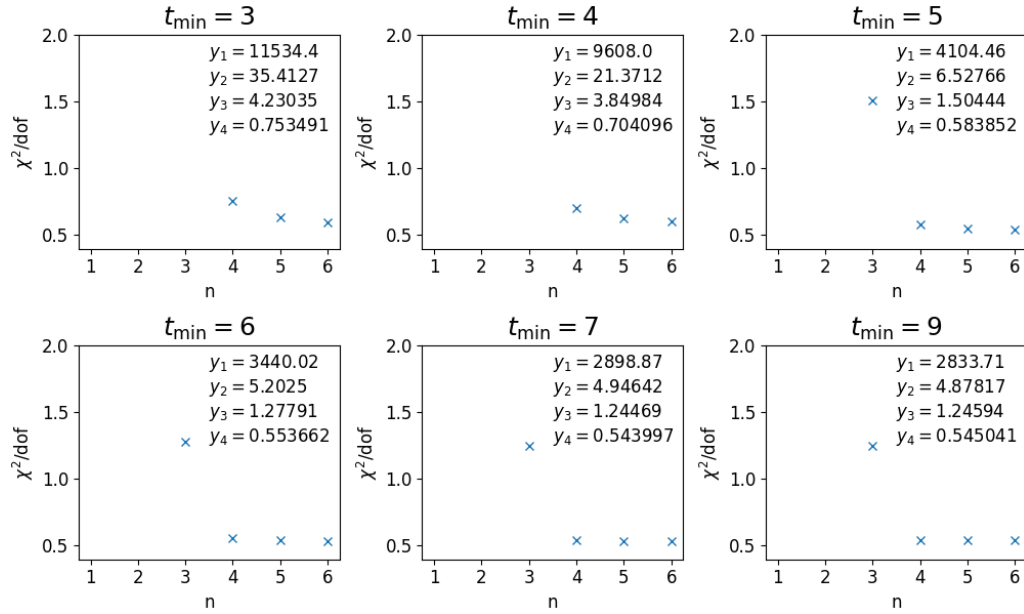


Figure 4: Fitting quality for fine lattice.

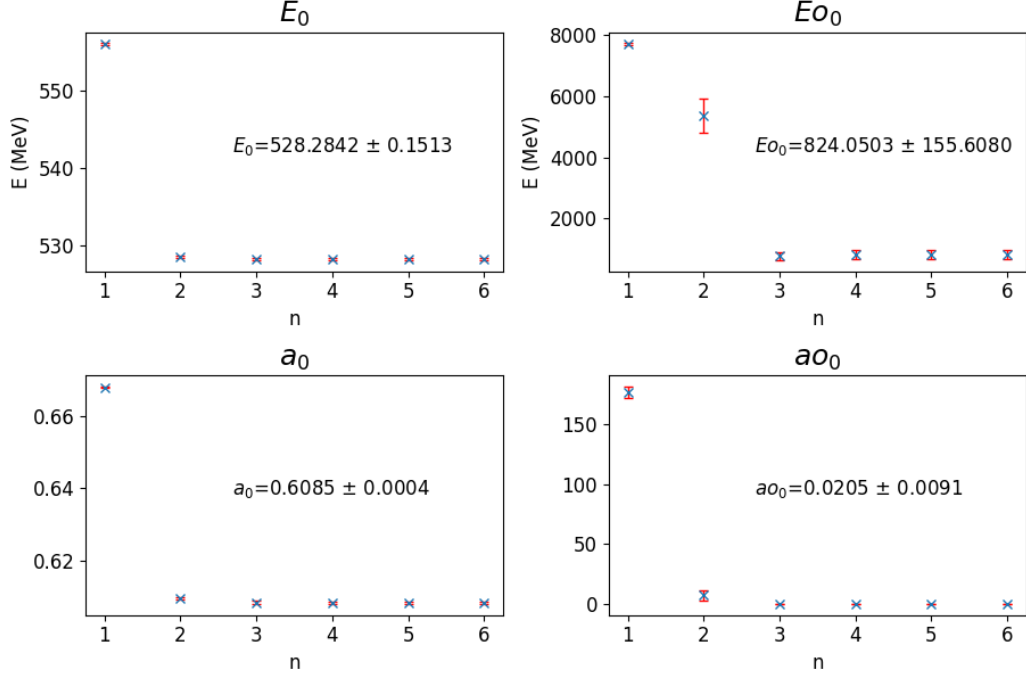


Figure 5: K meson fitting results on a coarse lattice.

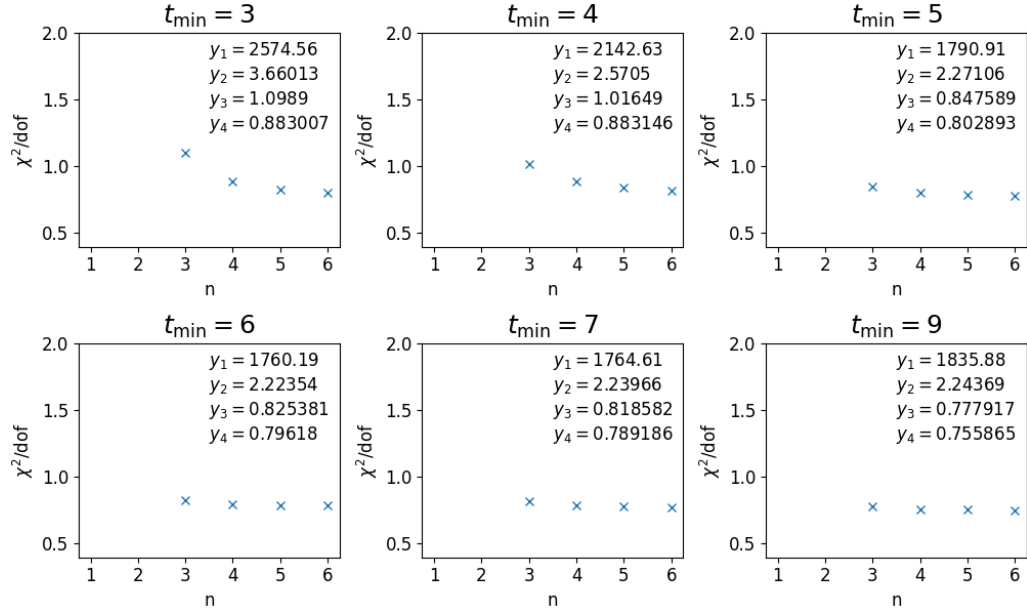


Figure 6: Fitting quality for coarse lattice.

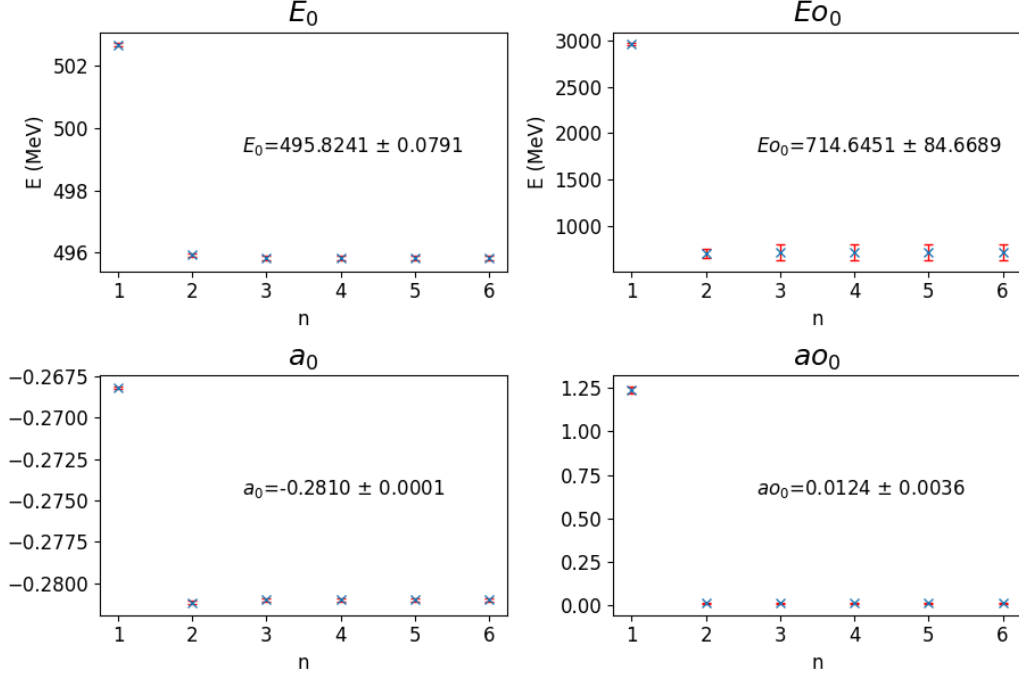


Figure 7: K meson fitting results on a very coarse lattice.

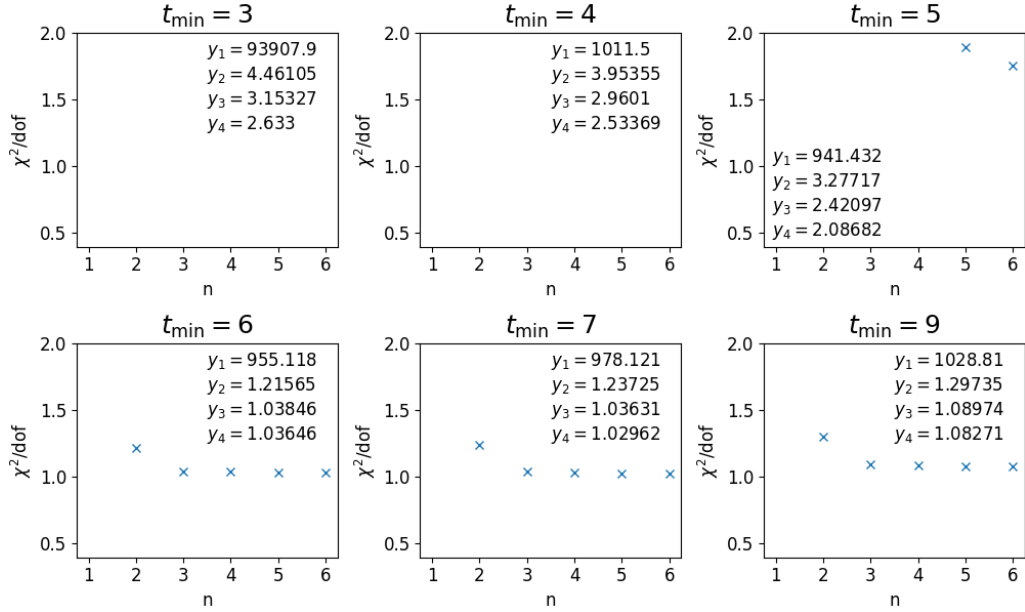


Figure 8: Fitting quality for very coarse lattice.

Label	a (fm)	m_l/m_s	E_0 (MeV)		a_0	
			K	D	K	D
Very Coarse	0.15	0.036	495.92(08)	1887.3(51)	-0.281 01(11)	-0.2286(65)
Coarse	0.12	0.2	528.28(15)	1893.0(15)	+0.608 53(37)	+0.1880(12)
Fine	0.09	0.2	530.16(24)	1889.3(11)	+0.139 78(12)	+0.1242(07)

Table 2: Results from fitting all datasets. $t_{\min} = 6$ for D mesons, and $n = 4$ for all results. Uncertainties given in parentheses are statistical.

4.2 Results

Taking into account the values of parameters for which all the datasets had reliable fits, the smallest (and least likely to lead to overfitting) viable choices for the parameters are $n = 4$ and $t_{\min} = 6$. The collection of E_0 and a_0 values that are found with these parameters is given in Table 2. The energy values here seem reasonable as they correspond roughly to the masses provided in [13]: $m_K \approx 497$ MeV and $m_D \approx 1865$ MeV. Of course, these values are far outside the provided statistical error bounds, but this is expected as we have both unphysical quark masses, and discretisation errors. Our next step will extrapolate our values to the chiral-continuum limit, which should give accurate results.

5 Chiral-Continuum Extrapolation

TODO: Extrapolation

6 Conclusion

TODO: Conclusion

References

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