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Effective spin models from the Yang-Mills action

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I present a method for finding the variation of the effective spin (or Polyakov line) action along any path in the configuration space of effective spins, given only the action of the underlying lattice gauge theory. The method can be applied, in principle, to extract the effective spin model from the underlying gauge theory.

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1. Introduction

The study of the QCD phase diagram at large baryon density calls for the introduction of a quark chemical potential into the QCD Lagrangian. This results in an intrinsically complex action, with the consequence that importance sampling, essential to all lattice Monte Carlo simulations, can no longer be applied in a straightforward way. It is possible that one may get around the problem by simulating instead the corresponding Polyakov line action.

The Polyakov line action S_P , also known as the "effective spin" action, is obtained from the QCD action S_{QCD} by integrating out all degrees of freedom in the theory, under the constraint that the Polyakov line holonomies are fixed. This gives us, in principle, a D = 3 dimensional theory whose phase structure ought to parallel that of full QCD. The integration can be carried out analytically at strong lattice couplings and heavy quark masses via the hopping parameter expansion, to give an action of the following form:

$$S_P = \beta_P \sum_{\mathbf{x}} \sum_{i=1}^{3} [\mathrm{Tr} U_{\mathbf{x}}^{\dagger} \mathrm{Tr} U_{\mathbf{x}+\hat{i}} + \mathrm{Tr} U_{\mathbf{x}} \mathrm{Tr} U_{\mathbf{x}+\hat{i}}^{\dagger}] + \kappa \sum_{\mathbf{x}} [e^{\mu} \mathrm{Tr} U_{\mathbf{x}} + e^{-\mu} \mathrm{Tr} U_{\mathbf{x}}^{\dagger}], \qquad (1.1)$$

The sign problem in this model seems relatively mild, for a wide range of parameters β_P , κ , μ , and can be handled in a number of different ways. These include the "flux representation" [1], reweighting [2], stochastic quantization [3], and mean field methods [4]. If we knew S_P corresponding to S_{QCD} at realistic gauge couplings and light quark masses, and if the sign problem remains mild, then in principle we could determine the phase diagram of S_{QCD} by computing the phase diagram of S_P . The first problem, however, is to find S_P in the parameter range of interest. In this talk I will discuss a novel approach to this problem; more details can be found in ref. [5].¹

2. The relative weights approach

Consider QCD on an $L^3 \times N_t$ lattice volume, with $\mu = 0$ for now. It is convenient to fix to temporal gauge, with timelike links set to the unit matrix everywhere except on one timeslice, say at t = 0. Then, by definition

$$\exp\left[S_P[U_{\mathbf{x}}]\right] = \int DU_0(\mathbf{x}, 0) DU_k D\overline{\psi} D\psi \left\{\prod_{\mathbf{x}} \delta[U_{\mathbf{x}} - U_0(\mathbf{x}, 0)]\right\} e^{S_{QCD}} .$$
(2.1)

Let us choose a set of "effective spin" configurations

$$\left\{ \{ U_{\mathbf{x}}^{(i)}, \text{all } \mathbf{x} \in V_3 \}, \ i = 1, 2, ..., M \right\}.$$
(2.2)

where V_3 is a three-dimensional lattice of volume L^3 , and imagine restricting the timelike link variables at t = 0 to this set. Define the partition function

$$\mathscr{Z} = \int DU_0(\mathbf{x}, 0) DU_k D\overline{\psi} D\psi \sum_{i=1}^M \left\{ \prod_{\mathbf{x}} \delta[U_{\mathbf{x}}^{(i)} - U_0(\mathbf{x}, 0)] \right\} e^{S_{QCD}} , \qquad (2.3)$$

¹For alternative approaches based on (i) strong-coupling expansions, cf. [2]; and on (ii) the inverse Monte Carlo method, cf. [6] and [7].

and consider the ratio

$$\frac{\exp\left[S_{P}[U^{(j)}]\right]}{\exp\left[S_{P}[U^{(k)}]\right]} = \frac{\int DU_{0}(\mathbf{x},0)DU_{k}D\overline{\psi}D\psi\left\{\prod_{\mathbf{x}}\delta[U_{\mathbf{x}}^{(j)}-U_{0}(\mathbf{x},0)]\right\}e^{S_{QCD}}}{\int DU_{0}(\mathbf{x},0)DU_{k}D\overline{\psi}D\psi\left\{\prod_{\mathbf{x}}\delta[U_{\mathbf{x}}^{(k)}-U_{0}(\mathbf{x},0)]\right\}e^{S_{QCD}}}$$

$$= \frac{\frac{1}{\mathscr{Z}}\int DU_{0}(\mathbf{x},0)DU_{k}D\overline{\psi}D\psi\left\{\prod_{\mathbf{x}}\delta[U_{\mathbf{x}}^{(j)}-U_{0}(\mathbf{x},0)]\right\}e^{S_{QCD}}}{\frac{1}{\mathscr{Z}}\int DU_{0}(\mathbf{x},0)DU_{k}D\overline{\psi}D\psi\left\{\prod_{\mathbf{x}}\delta[U_{\mathbf{x}}^{(k)}-U_{0}(\mathbf{x},0)]\right\}e^{S_{QCD}}}, \quad (2.4)$$

Both the numerator and denominator have the interpretation of a probability, i.e.

$$\operatorname{Prob}[U^{(j)}] = \frac{1}{\mathscr{Z}} \int DU_0(\mathbf{x}, 0) DU_k D\overline{\psi} D\psi \left\{ \prod_{\mathbf{x}} \delta[U_{\mathbf{x}}^{(j)} - U_0(\mathbf{x}, 0)] \right\} e^{\mathcal{S}_{QCD}}$$
(2.5)

is the probability that timelike links at t = 0 are found to belong to the *j*-th member of the set (2.2).

Numerical simulation of the system described by \mathscr{Z} is carried out by updating all degrees of freedom in the usual way, except for the timelike links at t = 0. These timelike links are updated simultaneously, by choosing at random one of the *M* members of the set (2.2), and accepting or rejecting it via the Metropolis algorithm. We keep track of the number of times N_n that the *n*-th member of the set is accepted, out of the total $N_{tot} = \sum_n N_n$. Then $\operatorname{Prob}[U^{(n)}] = N_n/N_{tot}$, in the limit $N_{tot} \to \infty$, and therefore

$$\frac{\exp\left[S_P[U^{(j)}]\right]}{\exp\left[S_P[U^{(k)}]\right]} = \lim_{N_{tot} \to \infty} \frac{N_j}{N_k}$$
(2.6)

From this information we can either test any proposal for S_P or, possibly, determine S_P from the data.

Let λ parametrize a path $U_{\mathbf{x}}(\lambda)$ through the space of Polyakov line configurations. We can use (2.6) to compute path derivatives $dS_P/d\lambda$ at $\lambda = \lambda_0$. Let configuration $U_{\mathbf{x}}^{(j)}$ correspond to the point at $\lambda_0 + \Delta \lambda$ and $U_{\mathbf{x}}^{(k)}$ to the point at $\lambda_0 - \Delta \lambda$. Then it follows, from (2.6), that

$$\left(\frac{dS_P[U_{\mathbf{x}}(\lambda)]}{d\lambda}\right)_{\lambda=\lambda_0} \approx \frac{1}{2\Delta\lambda} \left\{ \log \frac{N_j}{N_{tot}} - \log \frac{N_k}{N_{tot}} \right)$$
(2.7)

Numerical accuracy can be improved by choosing a set of configurations $\{U_{\mathbf{x}}^{(n)} = U_{\mathbf{x}}(\lambda_n)\}$ along the path with

$$\lambda_n = \lambda_0 + \left(n - \frac{M+1}{2}\right) \Delta \lambda \quad , \quad n = 1, 2, \dots, M \quad , \tag{2.8}$$

Then for sufficiently small $\Delta \lambda$ we have

$$\left(\frac{dS_P[U_{\mathbf{x}}(\lambda)]}{d\lambda}\right)_{\lambda=\lambda_0} \approx \text{ slope of } \log \frac{N_n}{N_{tot}} \text{ vs. } \lambda_n .$$
(2.9)

Our long-term program is to first determine S_P from S_{QCD} at $\mu = 0$. Then we obtain S_P at finite μ from the substitution $U_{\mathbf{x}} \rightarrow e^{N_t \mu} U_{\mathbf{x}}$, $U_{\mathbf{x}}^{\dagger} \rightarrow e^{-N_t \mu} U_{\mathbf{x}}^{\dagger}$. From there, the idea is to solve the effective spin theory by any of the means (flux representation, reweighting, stochastic quantization, mean field) that have been suggested so far. The immediate goal, however, is to see if the relative weights method can extract S_P from the underlying lattice gauge theory, and this talk will concentrate on the simplest example of pure SU(2) gauge theory.

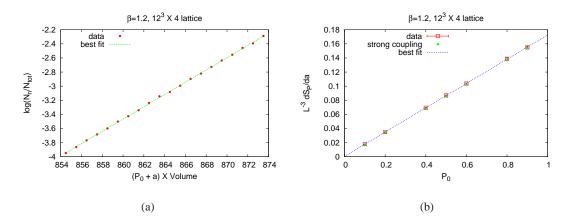


Figure 1: (a) The slope of the straight-line fit to the data shown gives an estimate for the derivative $L^{-3}dS_P/da$ of S_P with respect to the amplitude of spatially constant effective spin configurations. (b) A plot of the values for $L^{-3}dS_P/da$ vs. P_0 . Each data point is extracted from a plot similar to (a). Also shown are the corresponding strong-coupling values, and a best linear fit to the data points.

3. A test at strong coupling

The first question is whether the relative weights method can be used to determine S_P in a case where the answer is known. Let us take $\beta = 1.2$, $N_t = 4$ in lattice SU(2) pure gauge theory. S_P can then be readily computed by strong-coupling character expansion methods:

$$S_{P} = \left[1 + 4N_{t} \left(\frac{I_{2}(\beta)}{I_{1}(\beta)}\right)^{4}\right] \left(\frac{I_{2}(\beta)}{I_{1}(\beta)}\right)^{N_{t}} \sum_{\mathbf{x}} \sum_{i=1}^{3} \mathrm{Tr} U_{\mathbf{x}} \mathrm{Tr} U_{\mathbf{x}+\hat{i}} = \beta_{P} \sum_{\mathbf{x}} \sum_{i=1}^{3} P_{\mathbf{x}} P_{\mathbf{x}+\hat{i}}, \qquad (3.1)$$

where $P_{\mathbf{x}} \equiv \frac{1}{2} \text{Tr} U_{\mathbf{x}}$ and

$$\beta_P = 4 \left[1 + 4N_t \left(\frac{I_2(\beta)}{I_1(\beta)} \right)^4 \right] \left(\frac{I_2(\beta)}{I_1(\beta)} \right)^{N_t} .$$
(3.2)

The action can be divided into a potential and kinetic part

$$S_{P} = K_{P} + V_{P}$$

$$K_{P} = \frac{1}{2}\beta_{P}\sum_{\mathbf{x}}\sum_{i=1}^{3}(P_{\mathbf{x}}P_{\mathbf{x}+\hat{i}} - 2P_{\mathbf{x}}^{2} + P_{\mathbf{x}}P_{\mathbf{x}-\hat{i}})$$

$$V_{P} = 3\beta_{P}\sum_{\mathbf{x}}P_{\mathbf{x}}^{2},$$
(3.3)

First we determine V_P by relative weights. We choose the set of spatially constant configurations

$$U_{\mathbf{x}}^{(n)} = (P_0 + a_n) \mathbb{1} + i\sqrt{1 - (P_0 + a_n)^2} \sigma_3$$

$$a_n = \left(n - \frac{1}{2}(M+1)\right) \Delta a \ , \ n = 1, 2, ..., M \quad ,$$
(3.4)

So in this case *a* is the λ parameter, and it is easy to see that $dS_P/da = dV_P/dP_0$. The data for $\log(N_n/N_{tot})$ vs $L^3(P_0 + a_n)$ at $P_0 = 0.5$ is shown in Fig. 1(a), and the slope of this line gives us

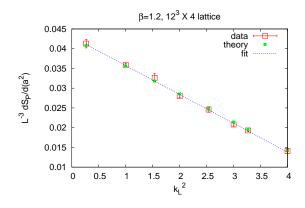


Figure 2: Derivative of the action w.r.t. path parameter a^2 vs. squared lattice momentum. Data is taken at strong gauge coupling $\beta = 1.2$ for plane-wave deformations. Squares indicate the relative-weights values, while green dots are the values obtained from the strong-coupling expansion.

 $L^{-3}dV_P/dP_0$. All data shown in this article was obtained at L = 12. Repeating the calculation for various values of P_0 (Fig. 1(b)) we find that dV_P/dP_0 is linear in P_0 . Integrating, and dropping an irrelevant constant of integration, the result is

$$V_P = \begin{cases} 0.1721(8) \sum_{\mathbf{x}} \frac{1}{2} P_{\mathbf{x}}^2 \text{ relative weights method} \\ 0.1710 \sum_{\mathbf{x}} \frac{1}{2} P_{\mathbf{x}}^2 \text{ strong-coupling expansion} \end{cases},$$
(3.5)

For the kinetic term, which by definition is zero for constant configurations, we choose a set of small plane wave deformations of a constant background

$$U_{\mathbf{x}}^{(n)} = P_{\mathbf{x}}^{(n)} \mathbb{1} + i\sqrt{1 - (P_{\mathbf{x}}^{(n)})^2} \sigma_3$$
$$P_{\mathbf{x}}^{(n)} = P_0 + a_n \cos(\mathbf{k} \cdot \mathbf{x})$$
(3.6)

where $k_i = 2\pi m_i/L$ are the components of the lattice wavevector, and it is convenient to choose $\lambda_n = a_n^2$ in this case. We compute $L^{-3}dS_P/d(a^2)$ from the slope of $\log(N_n/N_{tot})$ vs L^3a^2 , at fixed **k** and P_0 , with the result, at fixed $P_0 = 0.5$ but different choices of **k**, shown in Fig. 2, where k_L^2 is the usual (squared) lattice momentum $k_L^2 = 4\sum_i \sin^2(k_i/2)$. It follows that $L^{-3}dS_P/d(a^2) = -Ak_L^2 + B$, and further simulations find that the constants A, B are independent of P_0 . This information, together with the knowledge of the potential, allows us to deduce the action S_P (cf. [5] for details), and compare with the strong-coupling result:

$$S_{P} = \begin{cases} .0292(8) \sum_{\mathbf{x}} \sum_{i=1}^{3} P_{\mathbf{x}} P_{\mathbf{x}+\hat{i}} \text{ (relative weights method)} \\ .0285 \sum_{\mathbf{x}} \sum_{i=1}^{3} P_{\mathbf{x}} P_{\mathbf{x}+\hat{i}} \text{ (strong-coupling expansion)} \end{cases}$$
(3.7)

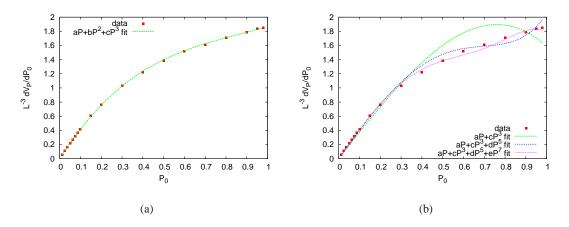


Figure 3: Derivatives of the potential. Subfigure (a) shows the best fit to the data by a polynomial $aP + bP^2 + cP^3$, while subfigure (b) shows a best fit by polynomials with two, three, and four odd powers of *P*, which are forms that might be expected from unbroken center symmetry.

4. The potential term at weaker couplings

Our convention is that the potential term is local, while the kinetic term vanishes for spatially constant configurations U. In that case

$$V_P = \sum_{\mathbf{x}} \mathscr{V}(U_{\mathbf{x}}) \text{ and } \mathscr{V}(U) = \frac{1}{L^3} S_P(U)$$
 (4.1)

and $K_P = S_P - V_P$ by definition. To compute the potential defined in this way, we only need to compute the variation of S_P with respect to the amplitude of spatially constant configurations. As at strong coupling, we choose the set (3.4) and compute $L^{-3}dS_P/da = L^{-3}dV_P/P_0$, but this time we work at a weaker coupling $\beta = 2.2, N_t = 4$, which is still in the confined phase, and fit the result to a polynomial in P_0 . The result is shown in Fig. 3(a), and the data is nicely fit by a cubic polynomial $c_1P_0 + c_2P_0^2 + c_3P_0^3$.

At first sight this result would seem to violate center symmetry, which for SU(2) requires that $V_P(P) = V_P(-P)$, and therefore the derivative should be an odd function of P, which would rule out the quadratic $c_2P_0^2$ term. On the other hand, fits to the data using only odd powers of P_0 are not very good, as we see in Fig. 3(b). The fit improves as more terms are added; it is, however, a little like fitting a step function with a truncated Fourier series. But in fact the quadratic term does not violate center symmetry, because simulations at $P_0 < 0$ find that the quadratic term is actually $c_2P_0^2 \times \text{sign}(P_0)$, and therefore, upon integration we find a center symmetric, albeit non-analytic, potential

$$V_P = \sum_{\mathbf{x}} \left(\frac{1}{2} c_1 P_{\mathbf{x}}^2 + \frac{1}{3} c_2 |P_{\mathbf{x}}|^3 + \frac{1}{4} c_3 P_{\mathbf{x}}^4 \right)$$
(4.2)

5. Towards the kinetic term

We consider two different sets of configurations. Type I will be small plane wave deformations on a constant background (3.6), as in the last section. Type II will simply be plane wave configurations

$$U_{\mathbf{x}}^{(n)} = P_{\mathbf{x}}^{(n)} \mathbb{1} + i\sqrt{1 - (P_{\mathbf{x}}^{(n)})^2} \sigma_3$$

$$P_{\mathbf{x}}^{(n)} = A_n \cos(\mathbf{k} \cdot \mathbf{x})$$

$$A_n = A_0 + \left(n - \frac{1}{2}(M+1)\right) \Delta A , \quad n = 1, 2, ..., M$$
(5.1)

For Type I the *a*-dependence begins at $O(a^2)$ and we again compute $dS_P/d(a^2)$, but unlike the strong-coupling case the result does depend on P_0 . For Type II configurations, we compute dS/dA.

It turns out that an expression for S_P which seems to fit the data from Type I and Type II configurations has the form

$$S_{P} = 2c \left\{ \sum_{\mathbf{xy}} P_{\mathbf{x}} \left(\sqrt{-\nabla_{L}^{2} + gP_{av}^{2} + g'\Delta P^{2}} \right)_{\mathbf{xy}} P_{\mathbf{y}} - \sum_{\mathbf{x}} \sqrt{gP_{av}^{2} + g'\Delta P^{2}} P_{\mathbf{x}}^{2} \right\} + \sum_{\mathbf{x}} \left(\frac{1}{2} c_{1} P_{\mathbf{x}}^{2} + \frac{1}{3} c_{2} |P_{\mathbf{x}}^{3}| + \frac{1}{4} c_{3} P_{\mathbf{x}}^{4} \right)$$
(5.2)

where

$$P_{av} = \frac{1}{L^3} \sum_{x} P_x$$
 and $\Delta P^2 = \frac{1}{L^3} \sum_{x} (P_x - P_{av})^2$ (5.3)

The data (red) and best fits (green) to the data for Type I and Type II configurations are shown in Figs. 4(a) and 4(b) respectively. The constants c_1, c_2, c_3 match the constants already determined for the potential. Although this action has been derived from data at very special (Type I and II) configurations, it is natural to conjecture, based on the success of this strategy at strong couplings, that (5.2) holds throughout the field configuration space.

6. Conclusions

The relative weights method for deriving the Polyakov line action has been tested at strong lattice couplings, where the answer known, and the method works well in that case. At a weaker coupling $\beta = 2.2$, $N_t = 4$ for pure SU(2) lattice gauge theory, where the Polyakov line action is not really known, a conjecture for the full action has been presented.

However, this conjecture comes with a strong caveat. The relative weights technique computes path derivatives of the Polyakov line action at any point along a path in the field configuration space, and the action S_P has to be deduced from that information. In the present case these derivatives have been computed only at very special points in configuration space, and these points are quite atypical, when compared to the thermalized Polyakov line configurations that would be generated in the course of a standard lattice Monte Carlo finite temperature simulation. The approximate form that the action takes at such special points cannot be guaranteed to hold in the more general situation. Therefore it is crucial to test the proposed action in other, more typical, regions of field configuration space. If the form (5.2) of the action survives this test, then the next check would be to compute an observable such as the Polyakov line correlator, to see if it agrees with the same quantity computed in the underlying lattice gauge theory. I hope to report on the results of these studies in the near future.

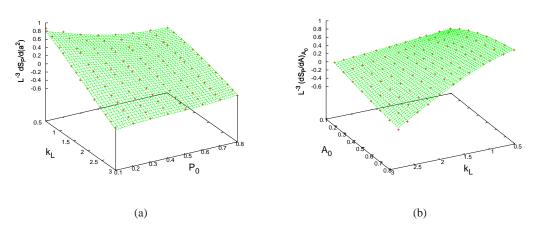


Figure 4: Derivatives of S_p for Type I and Type II configurations. Red crosses are data points, and the green surface is a best fit to the data derived from the analytic form (5.2). (a) Type I configurations: $L^{-3}dS/d(a^2)$ vs. lattice momentum k_L and Polyakov line P_0 . (b) Type II configurations: Variation of the effective action with plane wave amplitude, $L^{-3}dS_P/dA$ evaluated at $A = A_0$, for Polyakov line configurations proportional to plane waves $P_{\mathbf{x}} = A \cos(\mathbf{k} \cdot \mathbf{x})$, as a function of A_0 and lattice momentum k_L .

References

- [1] Y. D. Mercado and C. Gattringer, (2012), arXiv:1204.6074.
- [2] M. Fromm, J. Langelage, S. Lottini, and O. Philipsen, JHEP 1201, 042 (2012), arXiv:1111.4953.
- [3] G. Aarts and F. A. James, JHEP 1201, 118 (2012), arXiv:1112.4655.
- [4] J. Greensite and K. Splittorff, Phys.Rev. D86, 074501 (2012), arXiv:1206.1159.
- [5] J. Greensite, (2012), arXiv:1209.5697.
- [6] L. Dittmann, T. Heinzl, and A. Wipf, JHEP 0406, 005 (2004), arXiv:hep-lat/0306032.
- [7] T. Heinzl, T. Kaestner, and A. Wipf, Phys.Rev. D72, 065005 (2005), arXiv:hep-lat/0502013.