Hamiltonian study of improved U(1) lattice gauge theory in three dimensions

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A comprehensive analysis of the Symanzik improved anisotropic three-dimensional U(1) lattice gauge theory in the Hamiltonian limit is made. Monte Carlo techniques are used to obtain numerical results for the static potential, ratio of the renormalized and bare anisotropies, the string tension, lowest glueball masses and the mass ratio. Evidence that rotational symmetry is established more accurately for the Symanzik improved anisotropic action is presented. The discretization errors in the static potential and the renormalization of the bare anisotropy are found to be only a few percent compared to errors of about 20–25 % for the unimproved gauge action. Evidence of scaling in the string tension, antisymmetric mass gap and the mass ratio is observed in the weak coupling region and the behavior is tested against analytic and numerical results obtained in various other Hamiltonian studies of the theory. We find that more accurate determination of the scaling coefficients of the string tension and the antisymmetric mass gap has been achieved, and the agreement with various other Hamiltonian studies of the theory is excellent. The improved action is found to give faster convergence to the continuum limit. Very clear evidence is obtained that in the continuum limit the glueball ratio M_S/M_A approaches exactly 2, as expected in a theory of free, massive bosons.

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I. INTRODUCTION

Lattice gauge theory calculations have demonstrated important qualitative features of QCD, with increasing accuracy. Most lattice gauge theory calculations to date have been performed using Monte Carlo techniques in the Euclidean formulation. Although the Euclidean lattice gauge theory has been a very successful nonperturbative technique to compute the properties of elementary particles over the years, there are areas where progress has been very slow. Examples are QCD at finite temperature, glue thermodynamics, heavy quark spectra, etc. Some of these problems have resisted solution even by the powerful techniques of Euclidean field theory. This suggests that alternative methods should be pursued in parallel with Euclidean lattice gauge theory. A viable alternative that needs to be explored is the Hamiltonian version of QCD. This approach provides a valuable check of the universality of the Euclidean results [1] and has an appealing aspect in reducing lattice gauge theory to a many-body problem. As such the formalism is suited for the application of a host of analytic methods imported from quantum many body theory and condensed matter physics. It has been suggested that Hamiltonian lattice gauge theory could more readily handle finite density QCD [2]. The problems encountered in finite density QCD in the Euclidean formulation have prompted a return to the strong coupling expansions of early Hamiltonian lattice gauge theory [3]. Similar ideas have been pursued recently by Luo et al. [4,5] who propose an alternative Hamiltonian lattice formulation, "the Monte Carlo Hamiltonian," and have already demonstrated its validity and efficiency for the Φ^4 model [6].

Here we attempt to extend the standard Euclidean path integral Monte Carlo techniques to Symanzik improved U(1) gauge theory in three dimensions. Applications of this method have been extremely successful [7,8] and have given rise to great optimism about the possibility of obtaining results relevant to continuum physics from Monte Carlo simulations of lattice versions of the corresponding theory. Such an optimistic view is supported by recent work of Sexton *et al.* [9], Boyd *et al.* [10], Luo *et al.* [11] and Morningstar and Peardon [12] who have attempted to derive masses of the low-lying hadrons from such calculations and report successful results. More recent is the application of path integral Monte Carlo (PIMC) techniques to obtain results in the Hamiltonian limit for the U(1) model in (2+1) dimensions [13] and SU(3) lattice gauge theory in (3+1) dimensions on anisotropic lattices [14].

The use of improved actions [15,16] makes possible accurate Monte Carlo simulations of QCD on coarse lattices with greatly reduced computational effort [12,17,18]. In principle, with an improved action it is possible to achieve lattice volumes large enough to overcome finite-size effects and obtain measurements with good statistical errors. Coupled with tadpole improvement [19], the pursuit of the Symanzik program has led to significant progress in reducing the discretization errors and the renormalization of the anisotropy to the level of a few percent, and makes using anisotropic lattices no more difficult than isotropic ones. At the same time the merits of using an improved anisotropic lattice have been well understood [17,20]. Anisotropic lattices allow us to carry out numerical simulations with a fine temporal resolution while keeping the spatial lattice spacing coarse, i.e., a_t $< a_s$, where a_t and a_s are the lattice spacings in the temporal and spatial directions, respectively. This is especially important for QCD Monte Carlo simulations at finite temperature and heavy particle spectroscopy. But more importantly it should make extrapolations to the continuum limit more reliable.

As mentioned above, our aim in this work is to apply standard Euclidean path integral Monte Carlo techniques to extract the Hamiltonian limit for Symanzik improved

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 $U(1)_{2+1}$ lattice gauge theory on anisotropic lattices. The idea is to measure physical quantities on increasingly anisotropic lattices, and make an extrapolation to the extreme anisotropic limit. The effect of the plaquette improvement is examined by studying the scaling behavior and the sensitivity of scaling coefficients of the string tension and glueball masses in the Hamiltonian limit, $\xi \rightarrow 0$. The rest of the paper is organized as follows: In Sec. II we briefly review the formulation of the Symanzik improved U(1) gauge action in three dimensions on an anisotropic lattice. The details of the simulations and the methods used to extract the observables are described in Sec. III. Here we discuss our techniques for calculating the static quark potential, renormalization of anisotropy, string tension and glueball masses from Wilson loop operators. We present and discuss our results in Sec. IV. Scaling of the string tension, antisymmetric mass gap and the mass ratio in the weak-coupling region are tested against theoretical predictions and compared with the estimates obtained by other studies in the Hamiltonian limit. Our conclusions are given in Sec. V, along with an outline of future work.

II. IMPROVED ANISOTROPIC DISCRETIZATION OF $U(1)_{2+1}$

The Symanzik improved $U(1)_{2+1}$ gauge action on an anisotropic lattice is identical in form to the SU(3) case and is given¹ by [17]

$$S_{g} = \beta_{s} \xi \sum_{x} \sum_{i < j} \left[\frac{5}{3u_{s}^{4}} P_{ij}(x) - \frac{1}{12u_{s}^{6}} [R_{ij}(x) + R_{ji}(x)] \right] \\ + \frac{\beta_{t}}{\xi} \sum_{x} \sum_{i} \left[\frac{4}{3u^{2}u_{t}^{2}} P_{it}(x) - \frac{1}{12u_{s}^{4}u_{t}^{2}} R_{it}(x) \right], \quad (1)$$

where $P_{\mu\nu}$ and $R_{\mu\nu}$ are the 1×1 Wilson loop and 2×1 rectangular loop in the $\mu \times \nu$ plane respectively. At the tree-level the coefficients are chosen so that the action has no $O(a^2)$ discretization corrections. The spatial and temporal square and rectangular loops are given by

$$P_{ij}(x) = [1 - U_i(x)U_j(x+\hat{i})U_i^{\dagger}(x+\hat{j})U_j^{\dagger}(x)]$$
(2)

$$R_{ij}(x) = [1 - U_i(x)U_i(x+\hat{i})U_j(x+2\hat{i})U_i^{\dagger}(x+\hat{i} + \hat{j})U_i^{\dagger}(x+\hat{j})U_j^{\dagger}(x)]$$
(3)

$$P_{it}(x) = [1 - U_i(x)U_t(x+\hat{i})U_i^{\dagger}(x+\hat{t})U_t^{\dagger}(x)]$$
(4)

$$R_{it}(x) = [1 - U_i(x)U_i(x+\hat{i})U_t(x+2\hat{i})U_i^{\dagger}(x+\hat{i} + \hat{t})U_i^{\dagger}(x+\hat{i} + \hat{t})U_i^{\dagger}(x+\hat{i})U_i^{\dagger}(x)],$$
(5)

where x labels the sites of the lattice, *i*, *j* are the spatial indices and $U_i(x)$ is the link variable from site x to $x + \hat{i}$. The rectangular loop that extends two steps in the time direction has not been included in the above action. This has the advantage of eliminating negative residue high energy poles in the gluon propagator [17] but at the same time leaves errors of the order a_t^2 in the action. The $O(a_t^2)$ errors, however, are negligible provided a_t is small compared to a. The bare anisotropy parameter is ξ and is equal to the aspect ratio of the temporal and spatial lattice spacings at the tree level. At higher orders in the perturbative expansion, the bare anisotropy in the simulated action is not the same as the measured value, ξ_{phys} , due to quantum corrections [22]. The couplings β_s and β_t are defined by [23]

$$\beta_s = \frac{1}{g_s^2}, \quad \beta_t = \frac{1}{g_t^2}. \tag{6}$$

The two different couplings in Eq. (1) are necessary in order to ensure that in the continuum limit, physical observables become independent of the kind of lattice regularization chosen. In the case of an asymmetric lattice, this implies that physical quantities have to be independent of the anisotropy factor ξ . To achieve this, one needs to introduce different couplings for spatial and temporal directions, which depend on ξ . The ξ -dependence of the couplings g_s^2 and g_t^2 is due to quantum corrections and leads to an energy sum rule for the quark-antiquark potential, and the glueball mass, which differs in an important way from that which one would expect naively.

In the weak coupling limit of SU(N) lattice gauge theory, the relation between the scales of the couplings in Euclidean and Hamiltonian formulations has been determined analytically from the effective actions [23–26], using the background field method on the lattice. For small values of g_E , the couplings g_s and g_t can be expanded as

$$\frac{1}{g_t^2} = \frac{1}{g_E^2} + c_t(\xi) + O(g_E^2)$$
(7)

$$\frac{1}{g_s^2} = \frac{1}{g_E^2} + c_s(\xi) + O(g_E^2)$$
(8)

where g_E is the Euclidean coupling. For $\xi = 1$, one recovers the usual Euclidean lattice gauge theory, where $g_s = g_t$ $= g_E$. In the limit $\xi \rightarrow 0$, Eqs. (7) and (8) reduce to their Hamiltonian values and one obtains the relation between the Euclidean coupling g_E and its Hamiltonian counterpart g_H . Similar calculations have been performed to determine the anisotropic coefficients c_s and c_t at arbitrary anisotropy for a class of improved actions [20,27–29]. These coefficients become an important tool in the analysis of glue thermodynamics [30], the quark-gluon plasma [31,32] and for the determination of spectral functions at finite temperature [33]. Similar calculations have not yet been done for the U(1) theory, however, as far as we are aware.

¹The notation used here differs slightly from that used in Ref. [21], where the prefactors were absorbed into the definitions of β and ξ . We follow the notation introduced in [17].

Tadpole improvement [19] is introduced by renormalizing the link variables: $U_i(x) \rightarrow U_i(x)/u_s$, and $U_t(x) \rightarrow U_t(x)/u_t$, where the mean fields u_s and u_t can be defined by using the measured values of the average plaquettes in a simulation. In the plaquette mean-link formulation, the mean fields are determined self-consistently and are defined by

$$u_s = \langle P_{ij} \rangle^{1/4}, \quad u_t^2 u_s^2 = \langle P_{it} \rangle. \tag{9}$$

For $a_t \ll a$, the mean temporal link u_t is expected to be very close to unity. For simplicity we use a convenient and gauge invariant definition for u_s in terms of the mean spatial plaquette given by $u_s = \langle \operatorname{Re} \operatorname{Tr} P_{ij} \rangle^{1/4}$, and compute u_t from the temporal plaquette P_{it} in Eq. (9).

To obtain the Hamiltonian estimates from anisotropic lattices, a naive extrapolation procedure is followed. In this procedure we assume classical values of the couplings, i.e., $\beta = \beta_s = \beta_t$ in Eq. (1) and extrapolate the physical quantities to the extreme anisotropic limit, $\xi \rightarrow 0$ at constant β . Such a procedure is not strictly correct, however, at the quantum level because $\beta_s \neq \beta_t \neq \beta$ due to renormalization.²

As an example of the application of PIMC to the Symanzik improved action, we consider the case of compact U(1)gauge theory in three dimensions. The relevance of the model to QCD at finite temperature [35] has made it a standard proving ground for Hamiltonian lattice numerical methods. The model has two essential features in common with QCD, confinement [36–39] and chiral symmetry breaking [40]. Other common features are the existence of a mass gap and a confinement-deconfinement phase transition at some nonzero temperature. These similarities suggest that a comparison of the respective mass spectra should be informative. In the continuum limit of theory, the mass gap M is found to behave as [37]

$$M^2 a^2 \approx \operatorname{const} \beta \exp[-2\pi^2 \beta v(0)] = M_D^2 \qquad (10)$$

where $v(0) \approx 0.2527$ is the lattice Coulomb Green's function at zero separation. In Hamiltonian theory in which the space dimensions are discretized, $v_H(0) \approx 0.3214$ is the analogous Green's function [38]. It has been shown analytically that a linear confinement persists for all nonvanishing couplings, no matter how weak [37,38]. The string tension as a function of coupling also scales exponentially and obeys a lower bound

$$K \ge \operatorname{const} M_D \beta^{-1}.$$
 (11)

An interesting feature to explore in this context is whether the coupling to matter fields will change the permanent confinement status in (2+1) dimensions [41].

III. METHOD

A. Simulation details

To extract estimates of the string tension and the glueball masses using the Symanzik improved action in Eq. (1), a set of simulations are performed on lattices of size $N_s^2 \times N_t$ (N_s = 16 and N_t = 16-64), where N_s and N_t are the number of lattice sites in the spatial and temporal directions respectively. The lattice size in the time direction is adjusted according to the anisotropy used in order to keep the physical length in the spatial and temporal directions equal. To analyze the behavior in the strong and weak coupling regions, gauge configurations are generated using the METROPOLIS algorithm, for a range of couplings $\beta = 1 - 2.5$. The details of the algorithm are discussed elsewhere [13]. Starting from an arbitrary initial gauge configuration, 50 000 sweeps are performed for the equilibration of the configurations and the self-consistent determination of the mean-field parameters. A Fourier acceleration procedure [42,43] is used to overcome the stiffness against variations in the temporal plaquettes for high anisotropies. About 50% of the ordinary METROPOLIS updates are replaced with Fourier updates for $\xi \leq 0.4$ and about 100 000 further sweeps are performed to allow the system to equilibrate.

After thermalization, configurations are stored every 300 sweeps; 1200 stored gauge configurations are used in the measurement of the static quark potential and string tension and 1500 configurations for the glueball masses at each β . Measurements made on the stored configurations are binned into five blocks with each block containing an average of 250 measurements. The mean and standard deviation of the final observables are estimated simply by averaging over the block averages. The simulation parameters used for each configuration set are shown in Table I. It should be noted that Table I does not show all of the used values of parameters in the estimation of the physical quantities in Sec. IV.

B. The interquark potential and string tension

The static quark potential $V(\mathbf{r})$ is extracted from the expectation values of the timelike Wilson loops $W(\mathbf{r},t)$, which are expected to behave as

$$W(\mathbf{r},t) \approx \sum_{i} Z_{i}(\mathbf{r}) \exp[-tV_{i}(\mathbf{r})], \qquad (12)$$

where the summation runs over the excited state contribution to the expectation value, and i=1 corresponds to the lowest energy contribution. To obtain the optimal overlap of Wilson loop (and glueball) operators with the lowest state, it is necessary to suppress the contamination from excited states. This is done by using a simple APE smearing method [17,44,45] which is implemented by the iterative replacement of the original spatial link variables by a smeared link. Following the single-link smearing procedure, every spacelike link variable $U_i(x)$ on the lattice is replaced by

²One-loop calculations of the renormalization of the anisotropy and the gauge coupling in spatial and temporal directions for the improved Abelian lattice gauge theory are currently underway and will be reported elsewhere [34].

Volume	β	ξ	<i>u</i> _t	<i>u</i> _s	$(u_{s})^{4}$	$\langle P \rangle$ at u_s
$16^2 \times 32$	1.0	0.5	0.9991(3)	0.9218(2)	0.7220	0.7221(4)
$16^{2} \times 32$	1.45	0.5	0.9991(2)	0.9468(2)	0.8038	0.8039(3)
$16^{2} \times 32$	1.75	0.5	0.9997(3)	0.9551(3)	0.8324	0.8325(6)
$16^{2} \times 32$	2.0	0.5	1.0	0.9589(2)	0.8458	0.8459(4)
$16^{2} \times 32$	2.5	0.5	1.0	0.9614(4)	0.8543	0.8544(3)
$16^{2} \times 40$	1.0	0.4	1.0	0.9182(3)	0.7110	0.7111(4)
$16^{2} \times 40$	1.45	0.4	1.0	0.9432(3)	0.7917	0.7918(7)
$16^{2} \times 40$	1.75	0.4	1.0	0.9497(3)	0.8138	0.8139(5)
$16^{2} \times 40$	2.0	0.4	1.0	0.9524(5)	0.8228	0.8229(5)
$16^{2} \times 40$	2.5	0.4	1.0	0.9561(2)	0.8358	0.8359(2)
$16^{2} \times 48$	1.0	0.333	1.0	0.9152(3)	0.7016	0.7017(5)
$16^{2} \times 48$	1.45	0.333	1.0	0.9356(3)	0.7663	0.7664(4)
$16^{2} \times 48$	1.75	0.333	1.0	0.9401(2)	0.7815	0.7816(6)
$16^{2} \times 48$	2.0	0.333	1.0	0.9408(4)	0.7836	0.7837(4)
$16^{2} \times 48$	2.5	0.333	1.0	0.9454(4)	0.7991	0.7992(3)
$16^{2} \times 64$	1.0	0.25	1.0	0.9112(6)	0.6895	0.6892(3)
$16^{2} \times 64$	1.45	0.25	1.0	0.9287(3)	0.7440	0.7439(4)
$16^{2} \times 64$	1.75	0.25	1.0	0.9357(5)	0.7665	0.7663(5)
$16^{2} \times 64$	2.0	0.25	1.0	0.9381(4)	0.7745	0.7738(6)
$16^{2} \times 64$	2.5	0.25	1.0	0.9401(4)	0.7812	0.7811(3)

TABLE I. Simulation parameters at various β and ξ values.

$$U_{i}(x) \rightarrow P_{U(1)} \left[\alpha U_{i}(x) + \sum_{j \neq i} \sum_{\pm} U_{j}(x) U_{i}(x \pm \hat{j}) U^{\dagger}(x \pm \hat{i}) \right]$$
(13)

where $U_{-i} \equiv U^{\dagger}(x-\hat{i})$ and *i* and *j* are purely spatial indices. $P_{U(1)}$ denotes the projection onto U(1) and α is the smearing parameter. Operators constructed out of smeared links dramatically reduced the mixing with high frequency modes, thus removing the excited-state contamination in the correlation functions. The smearing fraction is fixed to $\alpha = 0.7$ and ten iterations of the smearing process are used. To reduce the variance, we use the technique of *thermal averaging* [13,46], which amounts to replacing the timelike link variables U_t by their local averages. This technique was applied to all temporal links except those adjacent to the spatial legs of loops, which are not independent. The technique has a dramatic effect in reducing the statistical noise.

The values of the effective potential are measured from the logarithmic ratio of successive Wilson loops

$$V_t(\mathbf{r}) = -\ln\left[\frac{W(\mathbf{r},t+1)}{W(\mathbf{r},t)}\right]$$
(14)

which is expected to be independent of t for t > 0. A plot of the effective potential is shown in Fig. 1 for $\beta = 1.70$ and $\xi = 0.4$ for various separations **r**. The dashed lines indicate the plateau values at various separations. As a result of heavy smearing, a good plateau behavior is seen at small t values for r=1 through 7. For $r \ge 6$, we see that the signal is dominated by noise for t > 6. We fix the fitting range to be, in most cases, t = 2 to 6.

The string tension is then extracted from the Wilson loops by establishing the linear behavior for the static quark potential at large separation \mathbf{r} . We have chosen to fit our results for $V(\mathbf{r})$ to the form [38]

$$V(\mathbf{r}) = a + K\mathbf{r} + b \ln(\mathbf{r}), \qquad (15)$$

where the linear term dominates the behavior at large separations and a logarithmic Coulomb term dominates at small separations.

C. Renormalization of anisotropy

Since the anisotropy ratio, η , is important in QCD simulations on the anisotropic lattices, we study its behavior by numerical simulations. Measurements of renormalization of anisotropy [47] have been made by comparing the static quark potential extracted from correlations along the different spatial and temporal directions. On an anisotropic lattice there are two different potentials $V_t(r)$ and $V_s(r)$ extracted from two different types of loops: timelike W_{xt} and spacelike W_{xy} Wilson loops. The two potentials differ by a factor of ξ_{phys} and by an additive constant, since the self-energy corrections to the static quark potential are different if the quark and antiquark propagate along the temporal or a spatial direction. The natural way to proceed is to build ratios of the Wilson loops,



FIG. 1. Effective potential as a function of Euclidean time *t*. From bottom up the horizontal lines correspond to r=1 through 7 and indicate the plateau values in the range $2 \le t \le 6$.

$$R_{t}(x,t) = \frac{W_{xt}(x,t+1)}{W_{xt}(x,t)}$$
$$R_{s}(x,y) = \frac{W_{xy}(x,y+1)}{W_{xy}(x,y)}.$$
(16)

Asymptotically for large τ and y, the ratios R_t and R_s approach

$$R_{t}(x,t) \approx Z_{x\tau} e^{-\tau V_{t}} + (\text{excited state contr.})$$

$$R_{s}(x,y) \approx Z_{xy} e^{-yV_{s}} + (\text{excited state contr.}). \quad (17)$$

The physical anisotropy is then determined from the ratio of the potentials $V_t(r)$ and $V_s(r)$ estimated from R_t and R_s , respectively. The lattice potentials defined by Eq. (17) contain contributions from the self-energy terms. The potential is simply parametrized as

$$V_{s}(\xi, r) = V_{s}^{0}(\xi) + V_{s}^{f}(\xi, r), \qquad (18)$$

where V_s^t is the lattice potential free of self-energy contributions. The timelike potential V_t is treated similarly. To eliminate the effect of the self-energy term V^0 in the potentials, we define a subtracted potential

$$V_{s}^{sub}(\xi, r, r_{0}) = V_{s}^{f}(\xi, r) - V_{s}^{f}(\xi, r_{0})$$
$$V_{f}^{sub}(\xi, t, t_{0}) = V_{t}^{f}(\xi, t) - V_{t}^{f}(\xi, t_{0})$$
(19)

where the subtraction points r_0 and t_0 are chosen to satisfy $t_0 = \xi r_0$ and the matching of the potential $V_t^f(t_0 = \xi r_0) = V_s^f(r_0)$ should be satisfied there. The subtraction radius r_0 should be chosen to be as small as possible so that fluctuations of the potential do not increase, in which case simulations with high statistics on a larger lattice are required. The renormalized anisotropy is determined from the ratio

$$\xi_{phys} = \frac{V_t^{sub}(\xi, r, r_0)}{V^{sub}(\xi, r, r_0)}$$
(20)

and the measured anisotropy parameter, η , is then given by

$$\eta = \frac{\xi_{phys}}{\xi}.$$
 (21)

An alternative approach to make the comparison is to fit the measured potentials with the forms [48]

$$a_{s}V_{s}(x) = a + \sigma a_{s}^{2}x + c \ln x$$

$$a_{s}V_{t}(t) = a + \sigma a_{s}a_{t}t + c \ln \left[\frac{a_{s}}{a_{t}t}\right].$$
(22)

The renormalized anisotropy ξ_{phys} is then determined from the ratio of the coefficients of the linear terms in the two cases. It is advantageous to use the potential at smaller *r*, where the statistical errors are smaller, and determine ξ_{phys} from the ratio of the coefficients of the Coulomb terms; however, such an estimate depends on short distance effects and is more sensitive to possible discretization errors of $O(a^4/r^4)$ [47].

D. Glueball masses

The numerical analysis of the mass of a glueball having a given J^{PC} proceeds through a study of the timelike correlations between spacelike Wilson loop operators $\Phi_i(t)$

$$C(t) = \langle \bar{\Phi}_i^{\dagger}(t) \bar{\Phi}_i(t) \rangle, \qquad (23)$$

where

$$\Phi_i(t) = \Phi_i(t) - \langle 0 | \Phi_i(t) | 0 \rangle \tag{24}$$



FIG. 2. Effective mass plot for the scalar glueball at β =2.0 and ξ =0.50.

is a gauge invariant vacuum subtracted operator capable of creating a glueball out of the vacuum. It is necessary to subtract the vacuum contribution for the scalar glueball with $J^{PC}=0^{++}$, because in the large Euclidean time limit, C(t) becomes dominated by the lowest energy state carrying the quantum numbers of Φ and these quantum numbers may coincide with that of the vacuum. The vacuum contribution is averaged over the whole ensemble before subtracting from the correlator. The glueball mass of interest is then extracted by studying the exponential decay of the correlation function for large Euclidean times, which is expected to behave as

$$C(t) = c_i [\exp(-m_i t) + \exp(-m_i (T-t))]$$

where m_i is the mass of the lowest-lying glueball which can be created by $\overline{\Phi}_i(t)$, and $T = N_t a_t$ is the extent of the periodic lattice in the time direction. Here, only the lowest "symmetric" (PC = ++) and "antisymmetric" (PC= --) glueball states are studied. The measured values of C(t) are expected to fall on a simple exponential curve assuming that the lattice is fine enough for the glueball mass to exhibit scaling behavior according to the theoretical predictions.

On a finite lattice with lattice spacing a, the operator $\Phi_i(t)$ has a small overlap with the glueball ground state, and the mass extracted from C(t) may be too large owing to the excited-state contamination. The overlap gets worse as the lattice spacing is reduced and nears the continuum limit, a



FIG. 3. Static potential, V(r), as a function of separation r at $\beta = 1.55$ and $\xi = 0.4$. The solid line is a fit to the form V(r) = a + br $+ c \ln(r)$ in the range $3 \le r \le 8$.

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(25)

 $\rightarrow 0$. This is obvious since the physical extension of the glueball remains fixed, while the operator Φ , constructed from small loops, probes an ever smaller region of the glueball wave function as the lattice spacing is decreased. Hence it becomes important to use an improved glueball operator so

as to have approximately the same size as the physical size

of the glueball. For such an operator, the overlap with the

glueball of interest is strong at small lattice spacing and

Following the variational technique of Morningstar and Peardon [17] and the smearing procedure of Teper [50], an optimized operator was found as a linear combination of the

signal-to-noise ratio is also optimal [49].

basic operators ϕ ,



FIG. 4. A close up of the static quark potential, V(r), at small r, using the Symanzik improved action. The solid line is a fit of Eq. (15) to on-axis points r=4 to 8. This plot involves measurements at $\beta=1.55$ for $\xi=0.4$ with 10 smearing sweeps at smearing parameter $\alpha=0.7$.

$$\Phi(t) = \sum_{\alpha} v_{i\alpha} \phi_{i\alpha}(t)$$
(26)

where the index α runs over the rectangular Wilson loops with dimensions $l_x = [n-1,n+1]$, $l_y = [n-1,n+1]$ and smearing $n_s = [m-1,m+1]$. The correlation function C(t)is then computed from the optimized glueball operator $\overline{\Phi}_i(t)$

$$C_{i}(t) = \sum_{t_{0}} \langle 0 | \bar{\Phi}_{i}(t+t_{0}) \bar{\Phi}_{i}(t_{0}) | 0 \rangle.$$
 (27)

 $\beta = 1.4142$ 0.8 $\xi = 0.4$ 0.7 0.6 V(r)0.5 0.4 0.3 0.2 1.5 2 2.5 3 3.5 r

FIG. 5. A close up of the static quark potential, V(r), at small r. The solid line is a fit of Eq. (15) to on-axis points r=4 to 8. This plot involves measurements from the standard Wilson action [13] at β = 1.4142 for ξ =0.4 with 10 smearing sweeps at smearing parameter α =0.7.

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Figure 2 shows the effective mass plot for the scalar glueball for the measurements at $\beta = 2.0$ and $\xi = 0.50$. The signal is seen out to the time slice 6 and reaches a plateau region for $1 \le t \le 5$. The data are noisy for t > 5.

We choose to fit the correlation function C(t) with the simple form

$$C_i(t) = c_1 \cosh m_i [T/2 - t] \tag{28}$$

to determine the glueball mass estimates.

IV. RESULTS AND DISCUSSION

A. Static quark potential and rotational symmetry

A plot of the static quark potential V(r) as a function of r at $\beta = 1.55$ and $\xi = 0.4$ is shown in Fig. 3. The data in this plot were obtained by looking for a plateau in the effective potential.

Because we are concerned to make long distance behavior consistent in both fine and coarse directions, it is advantageous to use Wilson loops of the largest possible spatial extent. However, in practice, the statistical errors in large Wilson loops grow exponentially with separation. We use Wilson loops of size 8×8 and fit the data by Eq. (15). We fit in the range $3 \le r \le 8$ so that we are not sensitive to the Coulomb term and the discretization errors associated with it. We see that the data are fitted very well giving the string tension, $K(=\sigma a^2)=0.044(3)$.

One of the main features of the improved discretization is the improved rotational invariance [21]. Discretization errors in the gluon action affect the extent to which continuum symmetries, such as rotational symmetry, are restored. To explore the extent to which rotational invariance is improved, we measure the potential at off-axis as well as onaxis separations. Thus for improved rotational invariance, the static potential, for example, at $\mathbf{r} = (4,3)$ should agree exactly with that at $\mathbf{r} = (5,0)$. TABLE II. Comparison between the measured anisotropy ξ_{phys} and the input anisotropy ξ for the Symanzik improved and the standard Wilson actions [13]. The measured differences in the offaxis potential at $\mathbf{r} = (1,1)$ are also shown.

Action	β	ξ	ξ_{phys}		$\Delta V(\sqrt{2})$
			$r_0 = 2$	$r_0 = \sqrt{2}$	
Improved	1.35	0.50	0.496(5)	0.493(2)	0.03(1)
action		0.40	0.392(4)	0.390(6)	0.03(2)
		0.333	0.328(3)	0.320(5)	0.04(4)
	1.45	0.444	0.442(6)	0.441(4)	0.04(2)
		0.333	0.326(3)	0.321(6)	0.04(3)
		0.25	0.246(4)	0.240(7)	0.05(4)
	1.55	0.40	0.402(2)	0.398(6)	0.04(2)
		0.25	0.241(5)	0.239(7)	0.06(4)
	1.65	0.333	0.334(5)	0.332(5)	0.02(1)
	1.75	0.25	0.252(4)	0.249(6)	0.03(1)
	2.0	0.333	0.335(6)	0.334(2)	0.03(1)
		0.25	0.242(4)	0.246(1)	0.06(2)
Wilson	1.35	0.444	0.418(5)	0.419(7)	0.12(1)
action	1.55	0.40	0.379(3)	0.374(6)	0.10(1)
	1.70	0.333	0.286(7)	0.289(5)	0.08(1)
	2.0	0.333	0.281(6)	0.292(2)	0.13(2)

For a fixed number of configurations and constant physical volume, we show the results from the Symanzik improved action and standard Wilson action in Figs. 4 and 5 respectively. We see that the off axis points for the improved lattice are excellently fitted by the rotationally invariant fitting curve (15) through r=3 to 8. The data from the Wilson action lie rather less close to the line of best fit. However, at large separations, the standard Wilson action does just as well as the improved action in extracting the static quark potential.



FIG. 6. Subtracted spatial and temporal potentials against the separation *r*. This plot involves measurements at $\beta = 1.75$ and $\xi = 0.333$ for the subtraction point $r_0 = \sqrt{2}$. The temporal potential V_t^{sub} has been rescaled by the input anisotropy.



FIG. 7. Unsubtracted spatial and temporal potentials against the separation *r* at $\beta = 1.75$ and $\xi = 0.333$. The temporal potential $V_t(r)$ has been rescaled by the input anisotropy.

As a quantitive measurement of the improvement, the potential measured in the simulation from nonplanar Wilson loops is compared with an interpolation to the on-axis data [47],

$$\Delta V(r) \equiv \frac{V_{\rm sim}(r) - V_{\rm fit}(r)}{\sigma r}.$$
 (29)

Results for $\mathbf{r} = (1,1)$ are given in Table II. With the mean-field inspired Symanzik improvement, the difference is only a few percent compared to a difference of about 10-20 % for the Wilson action [13].

B. Numerical determination of renormalized anisotropy

We choose the points $r_0 = 2$ and $\sqrt{2}$ to compute the subtraction potentials of Eq. (19) and use them to obtain the ratio in Eq. (20). The subtracted spatial and temporal potentials at the subtraction point $r_0 = \sqrt{2}$ are shown in Fig. 6. The anisotropies measured at these subtraction points for the improved and unimproved actions at different β values are compared in Table II. The two determinations of the anisotropies are in excellent agreement. These results show that the input anisotropy is normalized by less than a few percent for the improved action. This is in contrast with the



FIG. 8. Extrapolation of the string tension to the Hamiltonian limit $\xi \rightarrow 0$, for various β . Solid lines show the quadratic fits in ξ^2 to the data.



FIG. 9. The string tension as a function of inverse coupling. Our present estimates are shown as open triangles. Earlier results from standard Wilson action [13], *t*-expansion [52], Green's Function Monte Carlo simulations [53] and an exact linked cluster expansion [54] are shown as solid triangles, solid circles, open circles and open squares, respectively.

standard Wilson action [13], where the measured anisotropy is found to be about 20-30 % lower than the input anisotropy ξ .

Figure 7 shows the potentials computed from spatial and temporal Wilson loops without subtracting the self-energy terms. We find that the difference between the estimates computed from subtracted and unsubtracted potentials is less than 1% for the Symanzik improved action. We conclude that a few percent renormalization in the anisotropy is sufficiently small that it is unlikely to represent the dominant effect in the final estimates and it is safe to use the bare anisotropy for the tadpole improved Symanzik action.

C. String tension

To obtain estimates of the string tension in the Hamiltonian limit, an extrapolation is performed by a simple quadratic fit in powers of ξ^2 for each β value. The simulations run over a range of anisotropies, $\xi = 1 - 0.25$, thus enabling reliable extrapolation to the Hamiltonian limit. The errors for the extrapolation may be obtained by the "linear, quadratic, cubic" extrapolant method [51]. Figure 8 shows our estimates of the string tension as a function of the anisotropy ξ^2 for various fixed β values. Except at $\beta = 1.35$, a fairly smooth variation of string tension with ξ^2 for various couplings is seen. The curvature in the extrapolation at β



FIG. 10. The logarithm of the string tension as a function of inverse coupling. Our present estimates are shown as open triangles. The solid line is the result of fitting to form (6) for $1.45 \le \beta \le 1.95$. The dash-dot line is the β^4 order strong coupling expansion [57]. Our previous estimates [13] are shown as solid triangles. The dashed line represents the finite size behavior [58].

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=1.35 suggests that our estimate may be somewhat too high there.

Our extrapolated results for the string tension, $K = a^2 \sigma$, together with the earlier Hamiltonian estimates obtained from the *t*-expansion [52], Green's Function Monte Carlo simulations [53] and the Exact Linked Cluster Expansion (ELCE) [54] are plotted as a function of inverse coupling in Fig. 9. We see that the string tension displays an exponential behavior at weak coupling in accordance with the theoretical prediction. It has rigorously been shown that the string tension in $U(1)_{(2+1)}$ undergoes a "roughening transition" [55,56] at some intermediate coupling estimated to be near $\beta \approx 0.8$. Beyond the transition point, the different estimates of the string tension are expected to agree. The on-axis strong coupling series approximants fail to converge beyond $\beta = 0.8$, which prevents the analytic continuation of the series expansion beyond the roughening transition. The t-expansion results, however, do not suffer from this difficulty. A comparison with the GFMC [53] and an exact linked-cluster expansion [54] shows that our estimates are in good agreement with earlier estimates. The *t*-expansion estimates [52] are a little high, but still reasonably accurate. Of course, we do not expect that the results for the improved action should match exactly at finite coupling with other estimates which were computed for the unimproved action. We believe our PIMC estimates are more reliable and accurate and are also clearly consistent with the behavior predicted by Polyakov [37] and Göpfert and Mack [38].

Figure 10 shows the scaling behavior of the string tension together with results obtained using the standard Wilson action [13], as a function of β . The dashed-dot line is the strong-coupling expansion to order β^4 [57] and the solid line represents a fit to the weak-coupling asymptotic form (6). It can be seen that our present estimates appear to match nicely onto the strong and weak coupling expansions in their re-

TABLE III. Comparison of the Hamiltonian results obtained for the symmetric and antisymmetric scalar glueball masses (M_S, M_A) for the Symanzik improved and Wilson actions.

β	Improve	ed action	Wilson action		
	M_{S}	M_A	M_{S}	M_A	
1.0	1.68(1)	1.510(5)	1.9(1)	1.45(7)	
1.35	1.207(9)	0.811(5)	1.4(1)	0.9(1)	
1.55	0.75(1)	0.444(5)	0.8(2)	0.4(1)	
1.70	0.69(2)	0.361(6)	0.5(3)	0.24(9)	
1.90	0.352(6)	0.177(3)	0.3(1)	0.17(9)	
2.0	0.272(4)	0.136(2)	0.2(1)	0.10(6)	

spective limits. The fourth order strong coupling series expansion, obtained from integrated differential approximants, diverges beyond $\beta = 1.30$. In the weak-coupling region, the string tension is consistent with the predicted scaling behavior [38]. An unconstrained fit of the form (6) represents the data rather well in the interval $1.45 \le \beta \le 1.95$. The fit to the data gives a scaling slope of 3.04 ± 0.13 and an intercept of 1.42 ± 0.21 . The intercept of the scaling curve is roughly two times larger in magnitude than the theory predicts, compared to our previous results with the standard Wilson action [13] which were higher than theory by a factor of 5-6. Also in contrast with the Wilson action, a significant reduction in the errors is clearly apparent with the tadpole-improved Symanzik action (see Table III).

In summary, it appears that the overall exponential scaling behavior is the same for both actions, but the constant coefficient is lower for the Symanzik action by a factor of 2 to 3, and closer to the theoretical weak-coupling estimate. It seems highly plausible that a different action should give a different renormalization for the constant coefficient, although no analytic calculation of this effect has been done.



FIG. 11. Extrapolation of the antisymmetric glueball mass to the Hamiltonian limit $\xi \rightarrow 0$. Dashed lines show the quadratic fits in ξ^2 to the data.



FIG. 12. Glueball masses as functions of β . Our present results for the symmetric and antisymmetric glueball masses (M_S , M_A) are shown as open circles and triangles, respectively. Earlier results obtained from the Wilson action [13] are shown as solid circles and triangles, respectively.

D. Antisymmetric mass gap

The weak-coupling behavior of the mass gap is not exactly known. The rigorous analysis of Göpfert and Mack [38] showed that in the continuum limit $U(1)_{2+1}$ reduces to a massive scalar free field theory, with a mass gap M which is expected to decrease exponentially as the lattice spacing goes to zero. They showed that the lattice photon mass in the Villain action on a 3-dimensional Euclidean lattice is given by Eq. (6). It is often claimed in the literature that the Villain action is a high- β approximation of the Wilson action so that Eq. (6) should also hold in the weak-coupling limit of the Wilson model.

The extrapolation of the glueball masses to the Hamiltonian limit is shown in Fig. 11. The extrapolation is performed by using a simple quadratic fit in powers of ξ^2 . Again we see a smooth dependence on ξ^2 , for all β values analyzed here.

Our extrapolated results for the mass gap together with earlier Hamiltonian estimates obtained from the Wilson action [13] are plotted as a function of β in Fig. 12. Unlike the string tension, the strong-coupling expansion of the mass gaps is believed to be analytic near the roughening point. Comparison with earlier Hamiltonian estimates shows that our present data follow quite closely the strong-coupling expansion estimates, obtained by the method of integrated differential approximants, in the strong and weak-coupling regions. The *t*-expansion estimates [52], obtained from Padé approximants, are found to be substantially less accurate than those from the strong coupling expansion. Beyond $\beta=2$, the approximants for the t-expansion do not converge well so that no reliable estimates can be obtained beyond that coupling value.

The asymptotic scaling behavior of the antisymmetric mass gap is shown in Fig. 13. The solid line is the result of fitting for $1.4 \le \beta \le 2.25$ to the form (10) to find the scaling slope and the intercept of the scaling curve. Our results for

these coefficients are shown and compared with previous studies in Table IV. These results are obtained by fitting to the form $M^2 = \beta \exp[-f_0\beta + f_1]$ in the weak-coupling region. It can be seen that the agreement with the earlier results is remarkable. We find the constant coefficient (intercept of the scaling curve) is approximately 1.5 times larger in magnitude than the theory predicts. This is an improvement over our previous estimate using the Wilson action [13], where the constant coefficients were estimated a factor 5-6 times larger. The scaling slope is a little less than the theoretical prediction but in agreement with the estimates obtained in other numerical and analytic calculations (see Table IV). Several studies have provided evidence that the antisymmetric mass gap in the Wilson model of $U(1)_{2+1}$ Hamiltonian lattice gauge theory does not fall in the weak-coupling limit in the same manner as the periodic Gaussian model.³ This may be a signal of nonuniversality for the Abelian U(1)theory.

E. Mass gap ratio

The quantity of interest here is the dimensionless ratio R_M between the symmetric and antisymmetric mass gap in the large β limit. In this limit the ratio $R_M = M_S/M_A$ is expected to tend smoothly to its continuum value. In practice, this limiting value is found by increasing β from strong-coupling until the mass ratio levels off in the weak coupling region. Earlier studies of the photon mass [52] showed that the scaling of the mass ratio sets in for $\beta > 1$. If the continuum

³Based on the fact that periodic Gaussian models are special forms of Wannier-fuction expansions, Suranyi [59] has argued that a natural series of models, beginning with periodic Gaussian and approximating the Wilson model with arbitrary precision, does not exist.



FIG. 13. The logarithm of the antisymmetric mass gap as a function of β . Open triangles show our present estimates. The solid line is the fit to the data for $1 \le \beta \le 2$. Previous estimates from the Wilson action [13] are shown as solid triangles.

theory admits a stable bound state of two photons (a glueball), then the weak-coupling limit of the mass ratio will lie between 1 and 2. If the continuum theory is simply a freefield theory of massive scalar photons, as in the Villain model or if the glueball remains in the continuum theory only as a resonance, then the ratio $R_M = 2$ should be observed as β becomes large. Lana [64] argued for a sharp "transition point" at $\beta = 1.40$ where the symmetric bound state ceases to be stable, and crosses the level consisting of two free axial particles, so that mass ratio suddenly levels out at value 2.

In general, the most accurately calculated physical quantity is the string tension *K*. Therefore the first quantity that we shall extrapolate to the continuum limit will be of the form M/K, where *M* is the glueball mass. The leading corrections to such a ratio are known to be of the order $O(1/\zeta)^2$ [50], where ζ is some length scale. In the weak-coupling limit,

TABLE IV. Results for the coefficients f_0 (slope of the scaling curve) and f_1 (the intercept of the scaling curve) in the weak-coupling limit for the antisymmetric mass gap.

Source	f_0	f_1
Villain (Hamiltonian) [38]	6.345	4.369
Morningstar [52]	5.23	5.94
Hamer and Irving [60]	5.30	6.15
Hamer, Oitmaa and Weihong [61]	5.42	6.27
Plaquette Expansion [62]	5.01	5.82
Heys and Stump [63]	4.97	6.21
Lana [64]	4.10	4.98
Xiyan, Jinming and Shuohong [65]	5.0	5.90
Dabringhaus, Ristig and Clark [66]	4.80	6.26
Darooneh and Modarres [67]	4.40	5.78
Present Work	5.39(9)	6.3(1)

$$M_A/K \rightarrow \pi^2/2$$
 as $\beta \rightarrow \infty$,

in contrast to QCD, where M/\sqrt{K} is constant in the continuum limit. Figure 14 displays the estimates for this quantity, as a function of β . The data certainly appear to level out below $a_{eff}^2 \approx 0.006$ at a value 6. This is quite close to the predicted value for the Villain model.

Figure 15 shows the behavior of the dimensionless mass ratio $R_M = M_S/M_A$ as a function of effective lattice spacing squared, a_{eff}^2 , where a_{eff} is defined from Eq. (10) as [13]

$$a_{eff} = \sqrt{8\pi^2\beta}e^{-\pi^2\beta v(0)}$$

The plot shows that the mass ratio approaches very closely to the expected value of 2 in the large β limit. A linear fit to the data from $0.00026 \le a_{eff}^2 \le 0.0066$ gives $R_M = 2.007 \pm 0.01$, which is consistent with a continuum limit value $R_M = 2$. There is no sign of a stable, scalar glueball bound state, nor of any sharp "break" in the mass ratio. In comparison with previous results using the unimproved action, there are two notable features. First, the results are much more accurate. Second, the corrections to the continuum limit appear to be linear in a_{eff}^2 , whereas those for the unimproved action were linear in a_{eff} [13]. This provides impressive evidence that the improved action has achieved what it was designed to do: produce faster convergence to the continuum limit (see Table V).

V. SUMMARY AND CONCLUSIONS

In this work, we have applied the standard Euclidean path integral Monte Carlo method to obtain results in the Hamiltonian limit of the Symanzik improved $U(1)_{2+1}$ lattice gauge theory on anisotropic lattices. Monte Carlo results were obtained for the static quark potential, renormalized anisotropy, the string tension, and the lowest-lying glueball masses. The interquark potential with the improved action



FIG. 14. The dimensionless ratio $M_A/\beta K$ as a function of effective lattice spacing, a_{eff} . Open triangles show our present PIMC estimates. Our previous estimates using the Wilson action [13] are shown as solid triangles.

exhibits good rotational symmetry. We found that both the improved and standard Wilson action do equally well in extracting the static potential at large separations. The improved discretization allows substantially more accurate estimates of the string tension and the glueball masses. The extrapolations to the Hamiltonian limit were performed by simple quadratic fits.

In this limit the string tension displays an asymptotic behavior which is in excellent agreement with the behavior predicted by the weak-coupling approximation of Göpfert and Mack [38]. The scaling coefficient of the scaling curve for the mass gap was estimated to be roughly twice as large in magnitude as the weak-coupling prediction. This is an improvement over our previous unimproved results which were estimated to be larger by a factor of 5-6. We believe that these estimates can be further improved by taking the renormalization of the couplings into account, i.e., computing the action beyond the tree-level.

The weak-coupling behavior of the antisymmetric mass gap was found to agree with the theoretical expectations. The parameters agree quite well with the results obtained from previous numerical and analytic calculations. The mass ratio $R_M = M_S/M_A$ was observed to scale to 2.007 ± 0.01 in the continuum limit. No sign of a glueball bound state or a sharp crossover between the levels was seen. This is in excellent agreement with the statement of Göpfert and Mack [38], that the continuum limit corresponds to a theory of free bosons, where $R_M = 2$ exactly. It also shows very clearly the value of



FIG. 15. Mass ratio, R_M , as a function of effective lattice spacing, a_{eff} . Our present estimates are shown as the open triangles. The solid line is the fit to the data in the range $0.00026 \le a_{eff}^2 \le 0.0066$. The solid triangles show our earlier estimates obtained from the Wilson action [13].

TABLE V. Comparison of the Hamiltonian estimates for dimensionless mass ratio R_M obtained using the improved and standard Wilson actions.

β	R _M			
	Improved action	Wilson action		
1.0	1.11(2)	1.32(1)		
1.35	1.4(1)	1.5(2)		
1.55	1.7(1)	2.1(1.1)		
1.70	1.9(2)	2.2(1.5)		
1.90	2.0(3)	2.2(1.5)		
2.0	2.0(5)	2.4(1.6)		

using an improved action, in giving more rapid convergence to the continuum limit, as well as improved accuracy.

Taking advantage of the improved discretization on anisotropic lattices, we aimed to apply the Euclidean path integral Monte Carlo approach to examine the Hamiltonian limit of U(1) theory. The results obtained here clearly demonstrate that PIMC is a more reliable technique than other quantum Monte Carlo methods, such as Green's Function Monte Carlo, which gave only qualitative estimates of the string tension and the mass gap (see Table VI).

Although suffering from the disadvantage that it requires an extrapolation to the anisotropic limit, PIMC is the preferred Monte Carlo technique for obtaining reliable results in the Hamiltonian formulation, just as in the Euclidean formulation. In order to make the PIMC method a more valuable tool in Hamiltonian lattice gauge theories, it will be crucial to show that it allows one to treat eventually matter fields as well as gauge fields especially in the non-Abelian case.

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TABLE VI. PIMC results for the string tension, K, symmetric and antisymmetric glueball masses M_S , M_A and the mass ratio R_M .

β	a_{eff}	K	M_S/K	M_A/K	R_M
1.0	0.3724	0.329(2)	5.10(5)	4.58(4)	1.11(2)
1.15	0.2481	0.189(3)	7.8(1)	6.1(1)	1.28(3)
1.25	0.1884	0.141(2)	8.9(2)	6.4(1)	1.38(4)
1.35	0.1425	0.112(5)	10.8(5)	7.3(4)	1.4(1)
1.40	0.1239	0.096(5)	10.4(5)	6.7(3)	1.6(1)
1.45	0.1076	0.067(4)	13.9(9)	8.0(5)	1.7(2)
1.55	0.0810	0.045(3)	16.6(1.0)	9.8(6)	1.7(1)
1.60	0.0702	0.041(2)	18.1(1.1)	10.1(6)	1.8(1)
1.65	0.0608	0.034(2)	19.3(1.3)	10.4(9)	1.8(2)
1.70	0.0527	0.028(2)	24.1(2.1)	12.6(1.0)	1.9(2)
1.75	0.0425	0.023(2)	22.0(2.2)	11.6(1.1)	1.9(3)
1.80	0.0395	0.021(2)	21.6(2.7)	11.1(1.3)	1.9(3)
1.85	0.0341	0.016(2)	23.4(2.5)	12.3(1.3)	1.9(3)
1.90	0.0295	0.013(2)	25.9(3.2)	13.0(1.6)	1.6(3)
2.0	0.0220	0.010(2)	26.0(4.5)	12.9(2.2)	2.0(5)
2.10	0.0164	0.009(2)	23.4(5.0)	11.5(2.5)	2.0(6)
2.20	0.0122	0.008(2)	18.9(8.7)	9.6(2.4)	

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