Equation of state in 2+1 flavor QCD with improved Wilson quarks by the fixed scale approach

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Equation of state in $2 + 1$ flavor QCD with improved Wilson quarks by the fixed scale approach

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We study the equation of state in $2 + 1$ flavor QCD with nonperturbatively improved Wilson quarks coupled with the renormalization group-improved Iwasaki glue. We apply the $T$-integration method to nonperturbatively calculate the equation of state by the fixed-scale approach. With the fixed-scale approach, we can purely vary the temperature on a line of constant physics without changing the system size and renormalization constants. Unlike the conventional fixed-$N_t$ approach, it is easy to keep scaling violations small at low temperature in the fixed-scale approach. We study $2 + 1$ flavor QCD at light quark mass corresponding to $m_{\pi}/m_{\rho} \approx 0.63$, while the strange quark mass is chosen around the physical point. Although the light quark masses are still heavier than the physical values, our equation of state is roughly consistent with recent results with highly improved staggered quarks at large $N_t$.

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I. INTRODUCTION

The QCD equation of state (EOS) at high temperature plays a key role in understanding the nature of quark gluon plasma (QGP), e.g. as inputs of the hydrodynamical description of QGP space-time evolution in heavy-ion collision experiments [1]. Lattice QCD simulations provide us with the only systematic way to calculate the EOS nonperturbatively without resorting to phenomenological assumptions.

For a quantitatively reliable evaluation of EOS in QCD, it is indispensable to incorporate dynamical up, down, and strange quarks. However, dynamical quarks require a large computational effort on the lattice. Most calculations of EOS have been made in the fixed-$N_t$ approach, in which the temperature $T = (N_t a)^{-1}$ is varied on a lattice with fixed temporal size $N_t$ by varying the lattice spacing $a$ through a variation of coupling parameters on a line of constant physics (LCP). Here, we note that a sizable fraction of the total computational cost is required to systematically carry out zero-temperature simulations to determine the location of the LCP, to get basic information such as the scale and beta functions on the LCP, and to renormalize finite-temperature observables such as the EOS at each simulation point. In QCD with dynamical quarks, such systematic simulations are quite demanding.

We adopt the fixed-scale approach, in which we vary $T$ by varying $N_t$ at a fixed $a$ [2]. In this approach, because all the simulations are done with the same values of the coupling parameters, they are automatically on the same LCP. Furthermore, we need zero-temperature simulation at only one point to renormalize the observables at all $T$'s. Thus, the cost for the zero-temperature simulations can be largely reduced. To take or to confirm the continuum limit, we may repeat the calculations at several values of $a$. As the zero-temperature configurations, we may even borrow high statistic configurations on fine lattices, which were

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generated for spectrum studies at $T = 0$ and are open to the public on the International Lattice Data Grid (ILDG) [3].

The fixed-scale approach is complementary to the conventional fixed-$N_f$ approach in several respects: In the very high $T$ region where $T \simeq O(a^{-1})$, the fixed-scale approach suffers from lattice artifacts due to the coarseness of the lattice in comparison with the typical extent $T^{-1}$ of thermal fluctuations, while in the fixed-$N_f$ approach one can keep $T^{-1}/a = N_f$ finite even in the high temperature limit. In the fixed-scale approach, the spatial volume of the system is kept fixed at all $T$’s with the same spatial lattice size $N_s$, while in the fixed-$N_f$ approach $N_s$ has to be increased to quite large values at high $T$’s to keep the spatial volume. Large spatial volume is important at light quark masses to suppress volume effects in the hadron spectrum and thus in the determination of the scale and LCP. At small $T$’s, typically at $T \leq T_{pc}$, where $T_{pc}$ is the pseudocritical temperature, the fixed-scale approach keeps a small $a$, while the fixed-$N_f$ approach suffers from lattice artifacts due to large $a$. It should be kept in mind here that the fixed-scale approach requires high statistics in the low $T$ region, where we have a severe cancellation in the observables due to the zero-temperature subtraction procedure at large $N_f$. Nevertheless, we think it is worth taking advantage of smaller overall simulation costs with the fixed-scale approach to calculate the EOS in $2 + 1$ flavor QCD with small discretization errors around $T_{pc}$.

Another point of our study is the choice of the quark action on the lattice. Most lattice studies of hot/dense QCD have been done with computationally less expensive staggered-type lattice quarks [4,5]. However, their theoretical basis such as locality and universality are not well established. Therefore, to check the validity of these results it is important to compare the results with those obtained using theoretically sound lattice quarks, such as the Wilson-type quarks. See [6–9] for recent studies of QCD thermodynamics with Wilson-type quarks. A systematic study of the EOS with Wilson-type quarks has been done so far only in the case of two-flavor QCD [10,11]. We extend the study to the more realistic case of $2 + 1$ flavor QCD, using a nonperturbatively improved Wilson quark action coupled to a renormalization group (RG)-improved Iwasaki gauge action.

Thanks to the fixed-scale approach, we can take advantage of using the zero-temperature configurations on the ILDG. Using the same combination of lattice actions as ours, the CP-PACS + JLQCD Collaboration has generated a set of zero-temperature configurations in $2 + 1$ flavor QCD and has studied their hadronic spectrum [12,13]. Another attractive point of the fixed-scale approach in a study with improved Wilson quarks is that, unlike the case of the fixed-$N_f$ approach, we can keep the lattice spacing small at all temperatures and thus can avoid extrapolating the nonperturbative clover coefficient $c_{sw}$ to coarse lattices on which the improvement program is not quite justified.

Choosing a simulation point of the CP-PACS + JLQCD Collaboration, we carry out finite-temperature simulations to perform the first calculation of the EOS in $2 + 1$ flavor QCD with improved Wilson quarks. Although the light quark masses studied are still heavier than their physical values, we find that the EOS obtained is roughly consistent with recent results using highly improved staggered quarks in the fixed-$N_f$ approach at large values of $N_f$.

In the next section, we introduce the $T$-integration method which enables us to calculate the EOS nonperturbatively in the fixed-scale approach. The lattice setup and the simulation parameters are summarized in Sec. III. Results of gauge observables are presented in Sec. IV. In Sec. V the beta functions are evaluated. Our results on the EOS are shown in Sec. VI, and a summary is given in Sec. VII. The Appendix is devoted to a discussion about the choice of the interpolation procedure for the $T$-integration method. Preliminary results of this study have been reported in [14,15].

II. T-INTEGRATION METHOD

In conventional studies of EOS in the fixed-$N_f$ approach, the pressure $p$ is nonperturbatively estimated by the “integration method” [16]:

$$ p = \frac{T}{V} \int_{b_0}^{b} db \cdot \frac{1}{Z} \left( \frac{\partial Z}{\partial b} \right)_{sub} = -\frac{T}{V} \int_{b_0}^{b} db \cdot \left( \frac{\partial S}{\partial b} \right)_{sub}, $$

(1)

where $V$ is the spatial volume of the system, $Z$ is the partition function, $S$ is the lattice action with the coupling parameters $b = (\beta, \kappa_{sub}, \kappa_s, \cdots)$, and $\langle \cdots \rangle_{sub}$ is the thermal average with a zero-temperature contribution subtracted for renormalization. This relation is obtained by differentiating and then integrating the thermodynamic relation $p = (T/V) \ln Z$ in the coupling parameter space of $b$. The initial point $b_0$ is chosen in the low temperature phase such that $p(b_0) = 0$.

This method is inapplicable in the fixed-scale approach because $b$ is fixed in the simulations. To overcome the problem, we have developed the “$T$-integration method” [2]: Using a thermodynamic relation at vanishing chemical potential,

$$ T \frac{\partial}{\partial T} \left( \frac{p}{T^4} \right) = \epsilon - 3p, $$

(2)

where $\epsilon$ is the energy density, we obtain another nonperturbative estimate of the pressure,

$$ \frac{p}{T^4} = \int_{T_0}^{T} dT \frac{\epsilon - 3p}{T^4}, $$

(3)

with the initial temperature $T_0$ chosen such that $p(T_0) = 0$. Here, the trace anomaly $\epsilon - 3p$ is calculated as

$$ \epsilon - 3p = \frac{1}{T^3 V} \int da \left( \frac{\partial S}{\partial b} \right)_{sub}. $$

(4)
where \(a(d\hat{b}/da)\) is a vector of the beta functions which describes the variation of \(\hat{b}\) along the LCP.

When we vary \(T\) along a LCP by varying \(\hat{b}\), the integral in (3) is equivalent to that in (1), with the integration path chosen to be on the same LCP. However, (3) allows us to vary \(T\) without varying \(\hat{b}\). In the fixed-scale approach, we vary \(T\) by varying \(N_t\). Because \(N_t\) is discrete, we have to interpolate the data with respect to \(T\) to carry out the integration of (3). The systematic error from the interpolation should be checked.

In [2], the \(T\)-integration method was tested in quenched QCD and it was shown that the systematic error from the discreteness of \(T\) is under control when \(a\) is chosen sufficiently small, as adopted in spectrum studies. The EOS from the fixed-scale approach was shown to be well consistent with that from the fixed-\(N_t\) approach with large \(N_t\) (\(N_t \geq 8\)), except for the high temperature limit where the fixed-scale approach suffers from lattice discretization errors, as discussed in Sec. I.

### III. LATTICE SETUP

We adopt a nonperturbatively \(O(a)\)-improved Wilson quark action [17] coupled with the RG-improved Iwasaki gauge action [18] to simulate \(2 + 1\) flavor QCD:

\[
S_g = -\beta \sum_x \left\{ \sum_{\mu > \nu} c_0 W_{\mu \nu}^{1/2}(x) + \sum_{\mu, \nu} c_1 W_{\mu \nu}^{1/2}(x) \right\},
\]

\[
S_q = \sum_f \sum_{x, y} \bar{q}_f^x D^{x'y}_{f} q_f^y,
\]

\[
D^{x'y}_{f} = \delta_{x, y} - \kappa \sum_{\mu} \left\{ (1 - \gamma_{\mu}) U_{x, \mu} \delta_{x + \mu, y} + (1 + \gamma_{\mu}) U^+_{x, \mu} \delta_{x - \mu, y} \right\} \delta_{x, y} c_{SW}(\beta) \sum_{\mu > \nu} \sigma_{\mu \nu} F_{\mu \nu}
\]

with \(\kappa_a = \kappa_d = \kappa_{ud}\). The clover coefficient \(c_{SW}(\beta)\) has been evaluated nonperturbatively by the Schrödinger functional method in [12]. Hadronic properties have been systematically studied with this action by the CP-PACS, JLQCD, and PACS-CS Collaborations, down to the physical point [13,19–22].

In this study, we use the zero-temperature configurations by the CP-PACS and JLQCD Collaborations [13], which are open to the public at ILDG/ILDG [3]. The CP-PACS + JLQCD zero-temperature configurations are available at three \(\beta\)’s, five \(\kappa_{ud}\)’s, and two \(\kappa_s\)’s, i.e. at a total of 30 simulation points. Among them, we choose \(\beta = 2.05\), \(\kappa_{ud} = 0.1356\), and \(\kappa_s = 0.1351\), which correspond to the smallest lattice spacing and the lightest \(u\) and \(d\) quark masses (\(m_u/m_d \approx 0.63\)) with \(m_s\) near its physical point (\(m_{\eta_s}/m_\phi \approx 0.74\)). The hadronic radius is estimated to be \(r_0/a = 7.06(3)\) [23]. Setting the lattice scale by \(r_0 = 0.5\) fm, we estimate the scale as \(1/a \approx 2.79\) GeV (\(a \approx 0.07\) fm). The lattice size is \(28^3 \times 56\) (\(N_t a \approx 2\) fm), and the number of thermalized configurations are 650 (6500 trajectories), which are stored every 10 trajectories. Note that the \(u\) and \(d\) quark masses are still much larger than their physical values. We are planning to extend the study down to the physical point [22].

Using the relation between \(T\) and \(N_t\), our range of \(N_t\) corresponds to the range \(T \approx 7469\) MeV at \(\beta = 2.05\), as shown in Fig. 1. Previous studies of the pseudocritical temperature \(T_{pc}\) in two-flavor QCD with improved Wilson quarks at \(N_t \sim 6\) [6,25] suggest \(T_{pc}\) around 200 MeV for \(m_u/m_d \approx 0.63\) in two-flavor QCD. Taking into account the effect of the dynamical \(s\) quark and also our larger values of \(N_t \sim 14\) around the pseudocritical point, we expect a smaller value for \(T_{pc}\). In the succeeding sections, we show

### Table I. Simulation parameters and gauge observables.

The zero-temperature results (\(N_t = 58\)) are taken from [13] by the CP-PACS + JLQCD Collaboration. Temperature \(T\) is determined using \(1/a = 2.79\) GeV (\(a \approx 0.07\) fm) [13]. The Metropolis test is performed every 0.5 trajectories for finite-temperature simulations. \(\delta \tau\) is the molecular dynamics time step, and “Bin size” is the bin size for gauge observables, both in units of trajectories. “Trajectory” is the generated trajectory length after thermalization of “Thermalization” trajectories. “Plaqueette” and “Rectangular” are plaqueette and rectangular loop expectation values. \((L)\) and \(\chi_L\) are the bare Polyakov loop and its susceptibility, respectively.

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<tr>
<th>(N_t)</th>
<th>(T) [MeV]</th>
<th>(\delta \tau)</th>
<th>Trajectory</th>
<th>Thermalization</th>
<th>Bin size</th>
<th>Plaqueette</th>
<th>Rectangular</th>
<th>((L))</th>
<th>(\chi_L)</th>
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<td>1000</td>
<td>500</td>
<td>0.604 033(7)</td>
<td>0.377 087(8)</td>
<td>0.000 213(21)</td>
<td>0.001 820(20)</td>
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<tr>
<td>14</td>
<td>199</td>
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<td>6370</td>
<td>1000</td>
<td>500</td>
<td>0.604 048(2)</td>
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<td>0.001 172(67)</td>
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<tr>
<td>12</td>
<td>232</td>
<td>1/120</td>
<td>6460</td>
<td>1000</td>
<td>300</td>
<td>0.604 178(8)</td>
<td>0.377 314(5)</td>
<td>0.004 911(60)</td>
<td>0.014 25(25)</td>
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<td>10</td>
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<td>1/90</td>
<td>3935</td>
<td>500</td>
<td>200</td>
<td>0.604 262(9)</td>
<td>0.377 446(8)</td>
<td>0.014 70(11)</td>
<td>0.052 58(58)</td>
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<td>8</td>
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<td>2770</td>
<td>500</td>
<td>100</td>
<td>0.604 343(8)</td>
<td>0.377 580(3)</td>
<td>0.040 72(12)</td>
<td>0.115 (13)</td>
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<td>2785</td>
<td>500</td>
<td>50</td>
<td>0.604 590(2)</td>
<td>0.378 018(2)</td>
<td>0.109 81(11)</td>
<td>0.190 (15)</td>
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<td>3510</td>
<td>500</td>
<td>50</td>
<td>0.606 112(9)</td>
<td>0.380 962(1)</td>
<td>0.291 854(74)</td>
<td>0.216 8(92)</td>
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that our data suggest \( T_{pc} \sim 190 \text{ MeV} \) at our simulation point, as shown in Fig. 1 by the vertical shaded line.

The fixed-scale approach is not applicable at very high temperatures, where the lattice spacing \( a \) becomes too coarse to resolve thermal fluctuations [2]. We may estimate a typical length scale of thermal fluctuations by the thermal wavelength \( \lambda \sim 1/E \), where \( E \) is an average energy of massless particles at finite \( T \). We then obtain \( \lambda \sim 1/(3T) \) from \( E \sim 3T \xi(4)/\xi(3) \sim 2.7T \) for the Bose-Einstein distribution and \( E \sim 3T \xi(4)/\xi(3) \times 7/6 \sim 3.15T \) for the Fermi-Dirac distribution. Thus, data at \( T \approx 1/(3a) \) should be taken with care [26]. On the present lattice, the data at \( T \approx 700 \text{ MeV} \) may suffer from some lattice artifacts.

IV. GAUGE OBSERVABLES

The expectation values of gauge observables are measured every 0.5 trajectories. The results of basic observables are summarized in Table I. The time history of the Polyakov loop defined by

\[
L = \frac{1}{V} \sum_{x} \frac{1}{3} \text{Tr} \prod_{\tau=1}^{N_{t}} U_{(\tau-2/3, x)}^{4}
\]

is shown in Fig. 2. The gauge configurations are stored every five trajectories, on which quark observables are measured. By examining the bin-size dependence of the errors, we estimate the statistical errors for gauge observables by the jackknife method with the bin size listed in Table I, while those for quark observables are estimated with the bin size of 25 trajectories after thermalization of 1000 trajectories. Static quark potentials measured on the same configurations are studied in [23,26]. In the following, we disregard the statistical error in \( T \) from that of the lattice scale \( a \), which is about 0.5%. Note that, because the scale is common for all \( T \)'s in the fixed-scale approach, a shift in the scale \( a \) just causes an overall shift of \( T \).

The left panel of Fig. 3 shows the results of the Polyakov loop expectation value \( \langle L \rangle \) and its susceptibility \( \chi_L = N_{t}^{2}(\langle L^{2} \rangle - \langle L \rangle^{2}) \) as functions of \( T \). We find that \( \langle L \rangle \) starts deviating from zero at \( T \approx 180–200 \text{ MeV} \), suggesting the pseudocritical point around there.

For a comparison with the results of previous studies in the fixed-\( N_{t} \) approach, we have to renormalize \( \langle L \rangle \). Although the additive renormalization constant for free energies is independent of \( T \) and thus is common for all \( T \)'s in the fixed-scale approach, the Polyakov loop \( \langle L \rangle \sim e^{-F/T} \) does receive a \( T \)-dependent renormalization. To enable a direct comparison with the results of staggered-type quarks, we adopt the renormalization scheme proposed in [27]; i.e. we renormalize \( L \) such that the singlet free energy from \( L_{\text{ren}} = (Z_{\text{ren}})^{N_{t}} L \) becomes the Lüscher’s universal bosonic-string potential \( -\pi/(12r) + \sigma r \) at \( r = 1.5r_{0} \) [28], where \( \sigma \) is the string tension at \( T = 0 \). Using our potential data at \( T = 0 \) [23], we obtain \( Z_{\text{ren}} = 1.4801(90) \). Our results for \( \langle L_{\text{ren}} \rangle \) and the corresponding susceptibility \( \chi_{L_{\text{ren}}} \) are plotted in the right panel of Fig. 3. We note that the dependences on \( T \) in these quantities are largely influenced by the renormalization factor. In spite of the heavier light quark mass in our study, our results for \( \langle L_{\text{ren}} \rangle \) agree well with a result from the p4 staggered quark action in the fixed-\( N_{t} \) approach at \( N_{t} = 8 \) [29] (see the right panel of Fig. 3). Similar agreement of \( \langle L_{\text{ren}} \rangle \) between a smeared Wilson-type quark action and a smeared staggered-type quark action is reported in [9].

In Fig. 3, we also show the results of Polyakov loop susceptibilities. In the left panel of Fig. 3, besides a faint
bump at $T \sim 200$ MeV, we do not see a clear signal of a peak in $\chi_L$ at the two discrete simulation points in the range 180–200 MeV where $T_{pc}$ is expected. In the right panel of Fig. 3, the existence of a peak of $\chi_{ren}$ around these temperatures is not excluded, but it is due to the large errors there. The origin of the large errors will be discussed in Secs. VI and VII. This is in contrast with the case of our previous study in quenched QCD adopting the fixed-scale approach [30], in which we observe a clear peak of $\chi_L$, and also with the cases of full QCD studies adopting the fixed-$N_t$ approach with staggered-type (see e.g. [31]) and Wilson-type [6,25] quarks. As a possible cause of the absence of a clear peak in this study, we note that the resolution in $T$ is lower than that in our previous quenched study. We may have missed the peak between the simulation points. We also note the following: (i) We probably have a crossover in full QCD around the simulated quark masses instead of the first-order deconfining transition in quenched QCD. (ii) Our previous experience with improved Wilson quarks suggests that the peak becomes milder with increasing $N_t$. Our $N_t \sim 14$ around the crossover point is larger than those adopted in previous studies with the fixed-$N_t$ approach. (iii) The aspect ratio $N_s/N_t$ is not large at low temperatures in this study. All of these will make the peak milder and thus more difficult to detect when the resolution in $T$ is not fine enough.

V. BETA FUNCTIONS

To evaluate the trace anomaly according to (4), we need the beta functions $a(d\beta/da)$ and $a(d\kappa_f/da)$ ($f=ud$ and $s$). In this study, we define LCP’s by $m_q/m_{ps}$ and $m_{\eta_s}/m_\phi$ at $T=0$. The beta functions are determined nonperturbatively through the coupling parameter dependence of zero-temperature observables. We use the data of $am_p$, $m_q/m_p$, and $m_{\eta_s}/m_\phi$ at 30 simulation points of the CP-PACS + JLQCD zero-temperature configurations [13] to extract the beta functions. From a previous experience of two-flavor QCD with improved Wilson quarks in the fixed-$N_t$ approach [11], we expect that, although $a(d\kappa_f/da)$’s are much smaller than $a(d\beta/da)$, in the trace anomaly, the overall magnitude of the quark contribution proportional to $a(d\kappa_f/da)$ is comparable with that of the gauge part proportional to $a(d\beta/da)$, but with opposite sign. Therefore, evaluation of the quark contribution is important.

In our previous attempt [14], we have tried to evaluate the beta functions by the inverse matrix method, which was successful in the case of two-flavor QCD [11]. In 2 + 1 flavor QCD, we fitted the data of $am_p$, $m_q/m_p$, and $m_{\eta_s}/m_\phi$ as functions of three coupling parameters ($\beta$, $\kappa_{ud}$, $\kappa_s$), and inverted the matrix of the slopes of the former in terms of the latter to obtain the beta functions. However, it turned out that errors in $a(d\kappa_f/da)$ are too large to calculate the quark part EOS reliably, although the magnitude of the beta functions and the result for the gauge part of the trace anomaly are consistent with an expectation from the two-flavor case [14]. The situation is also similar when we use the data of pseudoscalar decay constants instead of $m_p$. We find that the large errors in $a(d\kappa_f/da)$ are mainly due to the matrix inversion procedure, through which all components of the inverse matrix get errors of similar magnitude. Because $a(d\kappa_f/da)$ are much smaller than $a(d\beta/da)$, we need more precise values of the slopes to suppress the errors in $a(d\kappa_f/da)$. In the present case of 2 + 1 flavor QCD, the data points of zero-temperature
configurations around the simulation point are not dense enough to achieve the required precision of the slopes.

To avoid the matrix inversion procedure, we now adopt an alternative method, the direct fit method [11]: We fit the coupling parameters, $\beta$, $\kappa_{ud}$, and $\kappa_s$, as a function of three observables, $am_\rho$, $m_\pi/m_\rho$, and $m_{\eta_c}/m_\phi$. Consulting the overall quality of the fits, we adopt the following third order polynomial function of the observables in this study:

$$
\begin{pmatrix}
\beta \\
\kappa_{ud} \\
\kappa_s
\end{pmatrix} = \tilde{c}_0 + \tilde{c}_1 (am_\rho) + \tilde{c}_2 (am_\rho)^2 + \tilde{c}_3 \left( \frac{m_\pi}{m_\rho} \right) + \tilde{c}_4 \left( \frac{m_\pi}{m_\rho} \right)^2 + \tilde{c}_5 (am_\rho) + \tilde{c}_6 \left( \frac{m_{\eta_c}}{m_\phi} \right) + \tilde{c}_7 \left( \frac{m_{\eta_c}}{m_\phi} \right)^2 + \tilde{c}_8 (am_\rho)
\times \left( \frac{m_{\eta_c}}{m_\phi} \right) + \tilde{c}_9 \left( \frac{m_\pi}{m_\rho} \right) + \tilde{c}_{10} (am_\rho)^3 + \tilde{c}_{11} \left( \frac{m_\pi}{m_\rho} \right)^3 + \tilde{c}_{12} \left( \frac{m_{\eta_c}}{m_\phi} \right)^3 + \tilde{c}_{13} (am_\rho) \left( \frac{m_\pi}{m_\rho} \right)^2 + \tilde{c}_{14} (am_\rho)^2 \left( \frac{m_\pi}{m_\rho} \right)
+ \tilde{c}_{15} (am_\rho) \left( \frac{m_{\eta_c}}{m_\phi} \right)^2 + \tilde{c}_{16} (am_\rho)^2 \left( \frac{m_{\eta_c}}{m_\phi} \right) + \tilde{c}_{17} \left( \frac{m_\pi}{m_\rho} \right) \left( \frac{m_{\eta_c}}{m_\phi} \right)^2 + \tilde{c}_{18} \left( \frac{m_\pi}{m_\rho} \right)^2 \left( \frac{m_{\eta_c}}{m_\phi} \right) + \tilde{c}_{19} (am_\rho) \left( \frac{m_\pi}{m_\rho} \right) \left( \frac{m_{\eta_c}}{m_\phi} \right).
$$

Note that the fits for the three coupling parameters are independent of each other. Figures 4 and 5 show the results of the global fit (9) as functions of $m_\pi/\rho$. The fits with $\text{dof} = 10$ lead to reasonable $\chi^2/\text{dof}$ ($\approx 1.63$, $1.08$, and $1.69$ for the fit of $\beta$, $\kappa_{ud}$, and $\kappa_s$, respectively), where the standard deviation of each coupling parameter is estimated by the error propagation rule using the errors of the observables and the partial derivatives of the resulting fitting function, Eq. (9), with respect to the observables, neglecting the covariance among the observables.

We define the LCP by fixing $m_\pi/m_\rho$ and $m_{\eta_c}/m_\phi$. Then, the beta functions are calculated as $\text{d}\beta/\text{d}a = (am_\rho)\partial\beta/\partial (am_\rho)$, etc., in terms of the coefficients $\tilde{c}_1$, $\tilde{c}_2$, $\tilde{c}_5$, $\tilde{c}_8$, $\tilde{c}_{10}$, etc. in (9). The resulting beta functions for our LCP ($m_\pi/m_\rho = 0.6337$, $m_{\eta_c}/m_\phi = 0.7377$) are shown in Fig. 6 as functions of $\beta$. Beta functions at other light quark masses are shown in Fig. 7. As the variable to set the scale, we may alternatively adopt $am_\pi$, $am_K$, or $am_K^*$ instead of $am_\rho$ in (9). Results of the beta functions, at our simulation point ($\beta = 2.05$ on our LCP), adopting various scale setting variables are listed in Table II. Taking the results from $am_\rho$ as the central value, we obtain

$$
\begin{align*}
\frac{\text{d} \beta}{\text{d} a} & = -0.279(24)(^{+40}_{-64}), \\
\frac{\text{d} \kappa_{ud}}{\text{d} a} & = 0.00123(41)(^{+86}_{-64}), \\
\frac{\text{d} \kappa_s}{\text{d} a} & = 0.00046(26)(^{+42}_{-44})
\end{align*}
$$

at our simulation point, where the first brackets are for statistical errors, and the second brackets are for systematic errors estimated by the variation of the scale setting.

**VI. EQUATION OF STATE**

With our lattice action (5) and (6), the trace anomaly $(\epsilon - 3p)/T^4$ is given by

$$
\begin{align*}
\frac{\epsilon - 3p}{T^4} = \frac{N_c^3}{N_f^3} \left( \frac{\text{d} \beta}{\text{d} a} \left( \frac{\partial S}{\partial \beta} \right)_\text{sub} + \frac{\text{d} \kappa_{ud}}{\text{d} a} \left( \frac{\partial S}{\partial \kappa_{ud}} \right)_\text{sub} \right) + \frac{\text{d} \kappa_s}{\text{d} a} \left( \frac{\partial S}{\partial \kappa_s} \right)_\text{sub} \\
& \quad + \frac{\partial S}{\partial \beta} \sum_{\mu > \nu} c_0 W^{1\times 1}_{\mu \nu}(x) + \sum_{\mu > \nu} c_1 W^{1\times 2}_{\mu \nu}(x)
\end{align*}
$$

with

$$
\left( \frac{\partial S}{\partial \beta} \right)_\text{sub} = -\left( \sum_{\mu > \nu} c_0 W^{1\times 1}_{\mu \nu}(x) + \sum_{\mu > \nu} c_1 W^{1\times 2}_{\mu \nu}(x) \right)
+ \frac{\partial c_{SW}}{\partial \beta} \sum_{f = u,d,s} \kappa_f \left( \sum_{\mu > \nu} \left( \sum_{j, \delta = u,d} T^{(f,c)}_{\mu \nu} \delta_{\mu \nu} F_{\mu \nu} (D_f)^{-1} \right)_{\text{sub}} \right).
$$

![Graph](https://example.com/graph.png)

**FIG. 4** (color online). The global fit for coupling parameters, $\beta$, as a function of $m_\pi/\rho$. Square symbols show coupling parameters in the CP-PACS/LQCD study. The solid lines show the global fit results for each simulation point with corresponding $m_\pi/m_\rho$ and $m_{\eta_c}/m_\phi$. To avoid a plot that is too busy, only half of the data points are shown ($\kappa_s = 0.1371, 0.1358$, and $0.1351$ at $\beta = 1.83, 1.90$, and $2.05$, respectively).
where $N_f = 2$ for $f = ud$ and 1 for $f = s$. We evaluate the traces in (12) and (13) by the random noise method with complex U(1) random numbers [25]. The number of noise vectors is 1 for each of the color and spinor indices. Results of the quark contributions in (12) and (13) are summarized in Table III.

In Fig. 8, the results of the trace anomaly (11) are shown by the solid curve. The curve is drawn by the Akima spline interpolation [32]. The central values are the results using the beta functions with the scale setting variable $am_\pi/C26 = 0.6337$ and $m_\eta_{ss}/m_\phi = 0.7377$, as functions of $\beta$. The scale setting is made with $am_\pi/C26$. Beta functions for $ud$ and $s$ are magnified by a factor of 100. Horizontal and vertical bars at each data point represent statistical errors.

FIG. 5 (color online). The same as Fig. 4 but for $\kappa_{ud}$ and $\kappa_s$.

In Fig. 8, the results of the trace anomaly (11) are shown by the solid curve. The curve is drawn by the Akima spline interpolation [32]. The central values are the results using the beta functions with the scale setting variable $am_\pi/C26 = 0.6337$ and $m_\eta_{ss}/m_\phi = 0.7377$, as functions of $\beta$. The scale setting is made with $am_\pi/C26$. Beta functions for $ud$ and $s$ are magnified by a factor of 100. Horizontal and vertical bars at each data point represent statistical errors.
cancel with each other among the contributions from different beta functions, leading to a smaller variation in total. Resulting systematic errors are shown by thick vertical bars in Fig. 8. The systematic errors thus estimated are smaller than the statistical errors in this study.

We find that \((e - 3p)/T^4\) is small at \(T = 174\) MeV but shows a peak at \(T = 199\) MeV and decreases towards higher \(T\). We note that the peak height of about 7 at \(T = 199\) MeV \((N_t = 14)\) is roughly consistent with recent results of highly improved staggered quarks (obtained at \(N_t = 6-12\)) in the fixed-\(N_t\) approach [4,5]. The shape of \((e - 3p)/T^4\) suggests that \(T_{pc}\) is located between 174 and 199 MeV.

Carrying out the \(T^4\) integration (3) using the Akima spline interpolation for the trace anomaly, we obtain the pressure \(p/T^4\) shown in Fig. 8. Here, we have chosen the starting point of the integration to be at \(N_t = 16\), where the trace anomaly vanishes within the statistical error. The energy density \(e/T^4\) is calculated by \(p/T^4\) and \((e - 3p)/T^4\). To our knowledge, this is the first result for EOS in 2 + 1 flavor QCD with dynamical Wilson-type quarks.

In our previous test in quenched QCD, we confirmed that the choice of the interpolation procedure has only minor effects on the EOS [2]. Because the resolution in \(T\) is coarser in the present study, we need to reexamine the influence of the interpolation procedures on the final values of the EOS. The results are summarized in Appendix B. We find that the systematic errors due to the choice of the interpolation procedure are small in the EOS in comparison with the present statistical errors.

The overall large errors in \(p/T^4\) and \(e/T^4\) are mainly due to the large statistic error in \((e - 3p)/T^4\) at \(T = 200\) MeV—they propagate to higher \(T\)'s through the numerical integration. The large statistic error in \((e - 3p)/T^4\) at \(T \leq 200\) MeV is caused by the enhancement factor \(N_t^4\) in (11) (note that \(S\) is proportional to \(N_t^2\)). Although the central value is largely canceled by the zero-temperature subtraction procedure, the errors are magnified. We find that the statistical fluctuation is much larger in the gauge part than in the quark parts. Note that the same difficulty exists also in the fixed-\(N_t\) approach when we increase \(N_t\) toward the continuum limit. In the fixed-scale approach, because high statistics is required at very low temperatures only, the overall numerical cost will still be lower than that in the fixed-\(N_t\) approach when we try to keep a similar magnitude of discretization errors around the transition temperature. In the present test, however, we stop at the current statistics and leave the task for the future investigation at the physical point.

An additional source of errors in Fig. 8 is the spacing of the data points in \(T\): Because our lattice spacing \(a\) is coarser than that of our previous study in quenched QCD [2], and also because \(N_t\) is restricted to be even due to the CPS simulation code with the even-odd preconditioning, we cannot have the resolution as achieved in our previous study. To improve the resolution in \(T\), we need to develop a simulation code for odd \(N_t\)'s. An alternative way out may be to combine results at different lattice spacing \(a\). Note that we can choose small values of \(a\) in the fixed-scale

<table>
<thead>
<tr>
<th>(N_t)</th>
<th>(T[\text{MeV}])</th>
<th>(N_{\text{conf}})</th>
<th>(\langle S^{\text{epp}}\rangle)</th>
<th>(\langle S^{\text{diag}}\rangle)</th>
<th>(\langle S^{\text{epp}}\rangle)</th>
<th>(\langle S^{\text{diag}}\rangle)</th>
</tr>
</thead>
<tbody>
<tr>
<td>58</td>
<td>⋯</td>
<td>390</td>
<td>-4.904 87(46)</td>
<td>1.904 649(79)</td>
<td>-4.748 78(44)</td>
<td>1.909 956(75)</td>
</tr>
<tr>
<td>16</td>
<td>174</td>
<td>447</td>
<td>-0.003 80(82)</td>
<td>-0.000 65(13)</td>
<td>-0.002 71(80)</td>
<td>-0.000 52(12)</td>
</tr>
<tr>
<td>14</td>
<td>199</td>
<td>447</td>
<td>-0.012 15(10)</td>
<td>-0.001 82(17)</td>
<td>-0.010 79(93)</td>
<td>-0.001 53(16)</td>
</tr>
<tr>
<td>12</td>
<td>232</td>
<td>495</td>
<td>-0.029 87(88)</td>
<td>-0.004 83(14)</td>
<td>-0.025 90(86)</td>
<td>-0.003 94(14)</td>
</tr>
<tr>
<td>10</td>
<td>279</td>
<td>287</td>
<td>-0.044 18(11)</td>
<td>-0.006 47(18)</td>
<td>-0.042 22(12)</td>
<td>-0.006 46(18)</td>
</tr>
<tr>
<td>8</td>
<td>348</td>
<td>319</td>
<td>-0.057 11(15)</td>
<td>-0.008 85(15)</td>
<td>-0.059 11(15)</td>
<td>-0.008 98(15)</td>
</tr>
<tr>
<td>6</td>
<td>464</td>
<td>159</td>
<td>-0.085 00(15)</td>
<td>-0.013 94(21)</td>
<td>-0.094 17(15)</td>
<td>-0.015 00(21)</td>
</tr>
<tr>
<td>4</td>
<td>697</td>
<td>95</td>
<td>-0.321 26(24)</td>
<td>-0.049 66(38)</td>
<td>-0.350 21(23)</td>
<td>-0.052 66(35)</td>
</tr>
</tbody>
</table>

TABLE II. Beta functions at our simulation point determined by the global fit (9) or with alternative scale setting variables. Values of \(\chi^2/\text{dof}\) for the fits are also given.

<table>
<thead>
<tr>
<th>Scale setting</th>
<th>(a)</th>
<th>(\chi^2/\text{dof})</th>
<th>(a)</th>
<th>(\chi^2/\text{dof})</th>
<th>(a)</th>
<th>(\chi^2/\text{dof})</th>
</tr>
</thead>
<tbody>
<tr>
<td>(am_p)</td>
<td>-0.279(24)</td>
<td>1.6</td>
<td>0.001 23(41)</td>
<td>1.1</td>
<td>0.000 46(26)</td>
<td>1.7</td>
</tr>
<tr>
<td>(am_{\pi})</td>
<td>-0.319(21)</td>
<td>1.2</td>
<td>0.001 79(38)</td>
<td>0.8</td>
<td>0.000 88(22)</td>
<td>1.3</td>
</tr>
<tr>
<td>(am_K)</td>
<td>-0.252(25)</td>
<td>1.0</td>
<td>0.001 05(44)</td>
<td>1.0</td>
<td>0.000 43(32)</td>
<td>1.3</td>
</tr>
<tr>
<td>(am_{K^*})</td>
<td>-0.215(28)</td>
<td>1.1</td>
<td>0.000 55(47)</td>
<td>1.2</td>
<td>0.000 02(36)</td>
<td>1.8</td>
</tr>
</tbody>
</table>
approach. When $a$’s are well in the scaling region, results for physical observables as functions of $T$ should lie on the same curves for these $a$’s, but at different discrete points. After confirming the insensitivity to a variation of $a$, we may combine the results at different $a$’s to more smoothly interpolate the data in $T$. We leave the application of these methods to future studies of EOS at the physical point.

Besides the large errors, our EOS looks roughly consistent with recent results with highly improved staggered quarks near the physical point: The peak of the trace anomaly from the stout quarks is located at $T = 190–200$ MeV with a peak height of about 4.0 [4]. A preliminary result from the highly improved staggered quarks gives a peak height of about 5.6 at $T = 200–220$ MeV [5]. We recall that our light quark masses are heavier than their physical values. The experience with improved staggered quarks suggests that the peak becomes higher as the light quark masses are increased (see, e.g., [4]).

# VII. SUMMARY

We calculated the EOS in $2 + 1$ flavor QCD with improved Wilson quarks by adopting the fixed-scale approach [2], with which we vary $T$ without varying the system volume on a fine lattice. As the first step towards the EOS with Wilson-type quarks in $2 + 1$ flavor QCD, we made simulations at $m_\pi/m_\rho \approx 0.63$, taking advantage of the fixed-scale approach to make use of high-resolution configurations by the CP-PACS + JLQCD Collaboration at $T = 0$ [13]. Although the light quark masses are still heavier than their physical values, our EOS looks roughly consistent with recent results with highly improved staggered quarks near the physical point [4,5].

To extend the study towards the physical point, however, we found a couple of issues that need to be solved: To obtain statistically accurate EOS at low temperatures, we need a large statistics proportional to $N_t^4$ (a power of $N_t$ is reduced due to the average over the lattice sites). This is, however, an unavoidable step to calculate observables suppressing discretization errors. Another source of systematic errors in EOS is the limited resolution in $T$ due to the discrete variation of $N_t$ in the fixed-scale approach. In the present study, because the lattice spacing $a$ is coarser than our previous quenched study, and because $N_t$ is limited to be even due to the simulation program set we have adopted, this seems to be non-negligible. To improve the resolution in $T$, we need simulations at odd values of $N_t$ and a finer lattice spacing $a$. An alternative way will be to combine results at different $a$’s, since we can choose fine $a$’s with the fixed-scale approach; thus, after confirming that the discretization effects are sufficiently small in the observables under study, we may combine the results at different $a$’s to more smoothly interpolate in $T$. We leave these trials to a forthcoming study with much lighter quarks, adopting the on-the-physical-point configurations by the PACS-CS Collaboration [22].

# ACKNOWLEDGMENTS

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# APPENDIX: COMPARISON OF INTERPOLATION PROCEDURES

To carry out the $T$ integration given by (3), we need to interpolate the data of the trace anomaly at discrete values of $T$ corresponding to the discrete values of $N_t$. In this appendix, we examine the interpolation procedures and their influences on the EOS with our data.

In the left panel of Fig. 9, we apply three different interpolation procedures to our data of the trace anomaly. Beta functions with the scale setting variable $am_\rho$ are adopted. The long-dotted line, dotted line, and solid line represent the results of straight line, cubic spline, and
Akima spline [32] interpolations, respectively. In our previous study in quenched QCD, we have adopted the cubic spline interpolation [2]. With our present data, however, we find the oscillatory interpolation curve by the cubic spline interpolation. This is due to the coarseness of the present data points—data are available only at even values of $N_t$. The cubic spline is not stable for data sets with sharp variations. In such cases, the Akima spline interpolation [32] is widely adopted. The Akima spline is a combination of local cubic polynomials and is known to suppress such oscillatory behavior around sharp variations. From Fig. 9, we find that the Akima spline leads to a more natural curve smoothly following the data points. Therefore, we adopt the Akima spline interpolation in this study.

To estimate the systematic error due to the choice of the interpolation procedure in the EOS, we perform the $T$ integration with these interpolations. The results for the pressure are shown in the right panel of Fig. 9. The strong oscillation of the interpolation curve from the cubic spline is averaged over through the integration, and the results of $p/T^4$ are well consistent for all three interpolations. We thus conclude that the systematic error in the EOS due to the choice of the interpolation procedure is much smaller than the statistical errors.

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