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Unquenched complex Dirac spectra at nonzero chemical potential: two-colour QCD lattice data versus matrix model

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We compare analytic predictions of non-Hermitian chiral random matrix theory with the complex Dirac operator eigenvalue spectrum of two-colour lattice gauge theory with dynamical fermions at nonzero chemical potential. The Dirac eigenvalues come in complex conjugate pairs, making the action of this theory real, and positive for our choice of two staggered flavours. This enables us to use standard Monte-Carlo in testing the influence of chemical potential and quark mass on complex eigenvalues close to the origin. We find an excellent agreement between the analytic predictions and our data for two different volumes over a range of chemical potentials below the chiral phase transition. In particular we detect the effect of unquenching when going to very small quark masses.

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Non-Hermitian operators appear in many different areas of Physics: in S-matrix theory with absorption [1]. in neural network and dissipative quantum dynamics [2], disordered systems with imaginary vector potential [3], Quantum Chromodynamics (QCD) with a non-zero chemical potential μ [4], or θ -vacuum term [5]. Studying the influence of these non-Hermiticity parameters may serve to a better understanding of fundamental problems in nature, such as the deconfinement transition observed in heavy-ion collisions, or the strong CP problem. In the present work we use an analytically solvable random matrix model (MM) as an effective theory based on symmetries, and focus on the applications to QCD at $\mu \neq 0$. Given the wide range of complex MM applications [1] we expect our approach to be relevant to other problems in the same symmetry class.

Lattice Gauge Theory (LGT) is one of the most prominent methods to study nonperturbative QCD. However, the introduction of a chemical potential poses a difficult problem: the Dirac operator and thus the action becomes complex non-Hermitian. In general this invalidates standard Monte-Carlo techniques using importance sampling. the so-called sign problem. The quenched approximation, eliminating the complex fermion determinant, has been explained to fail using a MM approach [4]. It maps QCD to a theory with conjugate quarks, with chiral symmetry unbroken in the massless limit (see [6] for a recent discussion, including phase quenched MM).

Since the role of dynamical fermions is crucial our goal is to show that MM can predict unquenched lattice data at $\mu \neq 0$. Although analytical MM results exist for unquenched QCD [7], a comparison to LGT remains an open challenge, due to a complex valued density with highly oscillatory regions [6, 7]. The various ways to avoid the sign problem including reweighting, Taylor expansion, imaginary chemical potential or combinations of these reviewed in [8] rely essentially on being close to the chiral phase transition at high temperature $T \approx T_c$.

These techniques do not apply easily at the region of our interest at small $T \approx 0$ and μ , where MM are conjectured to coincide with a particular limit of the underlying effective chiral Lagrangian, the ϵ -regime ($\epsilon \chi PT$). MM are a different realisation of the group integral over constant Goldstone modes at $\mu \neq 0$ for QCD [9, 10], and they clearly distinguish the three different chiral symmetry breaking (χSB) patterns at $\mu \neq 0$ [11]. Thus they may serve as a test for algorithms solving the sign problem. In order to access in this region with dynamical fermions we follow the suggestion [12]: gauge group SU(2) or the adjoint representation lead to a real action as complex Dirac eigenvalues come in conjugated pairs. We have chosen SU(2) with 2 staggered fermions because in this symmetry class MM predictions are available [13]. The difficulty is that large masses m effectively quench the Dirac eigenvalues close to the origin. Unquenching is seen only for small m depending on the particular size of the system. Our preliminary results for SU(2) have been presented in [14], compared to SU(2) at $\mu = 0$ [15] and quenched QCD at $\mu \neq 0$ [16, 17].

We start by introducing the MM [13], the complex extension of the chiral Gaussian Symplectic Ensemble, and its predictions. For an arbitrary number of N_f quark flavours with masses m_f it is defined as:

$$\mathcal{Z}_{\nu}(\{m_f\}; \mu) = \int_{\mathbb{H}^{(N+\nu)\times N}} d\Phi d\Psi \prod_{f=1}^{N_f} \det \begin{bmatrix} m_f \mathbf{1} & i\Phi + \mu\Psi \\ i\Phi^{\dagger} + \mu\Psi^{\dagger} & m_f \mathbf{1} \end{bmatrix} \\
\times \exp \left[-N \operatorname{Tr}(\Phi^{\dagger}\Phi + \Psi^{\dagger}\Psi) \right]. \tag{1}$$

The matrices Φ and Ψ of size $(N+\nu)\times N$ contain quaternion real matrix elements without being self-dual, 1 is the quaternion unity element. The Gaussian average of variance $\sigma = 1/\sqrt{2}$ replaces the gauge average, where we have fixed topology or equivalently the number of exact zero eigenvalues $\nu > 0$. The total number of eigenvalues $2N \sim V$ relates to the volume. For $\Psi = \mathbf{1}$ this MM was shown [11] to be in the symmetry class of the

adjoint representation, or gauge group SU(2) with staggered fermions, as in our case. In [11], the qualitative differences between the 3 χ SB classes were compared numerically. Assuming universality, which is true for the MM of QCD [7, 9, 10], we choose the μ -part $\Psi \neq \mathbf{1}$, with the same symmetry as the kinetic part. This choice in eq. (1) permits an eigenvalue representation with Jacobian $\sim \prod_{k>l}^N |z_k^2 - z_l^2|^2 |z_k^2 - z_l^{*2}|^2 \prod_{h=1}^N |z_h^2 - z_h^{*2}|^2$ [13], and mass term $\prod_{f=1}^{N} |z_j^2 + m_f^2|^2$ (here z = x + iy is real at $\mu = 0$). Due to the Jacobian complex eigenvalues are repelled both from the x- and y-axis, the signature of this χ SB class. The QCD MM [7] has no such repulsion.

In [13] all complex eigenvalue correlations were derived in a closed form as the Pfaffian of a matrix kernel of skew orthogonal polynomials on \mathbb{C} . We only give the density for $N_f = 2$ of equal mass needed here. For a discussion of matching MM and staggered flavours we refer to the first of Ref. [15]. No 2nd or 4th root is taken here. Taking the large-N limit at weak non-Hermiticity [1] the eigenvalues, masses $and \mu$ are rescaled as

$$z \cdot 2N\sigma \equiv \xi, \ m \cdot 2N\sigma \equiv \eta, \ \lim_{N \to \infty, \ \mu \to 0} \mu^2 \cdot 2N \equiv \alpha^2, \ (2)$$

resulting in the usual MM level spacing π . The ϵ -regime of the corresponding chiral Lagrangian [18] dictates the same scaling with two different constants: $V\Sigma m$ and $F_{\pi}^{2}V\mu^{2}$, the chiral condensate Σ and decay constant F_{π} , respectively. This implies the existence of one free fit parameter in α^{2} in our model. This observation has very recently been proposed to measure F_{π} for QCD using isospin chemical potential [19]. We note that MM predictions at $\mu = 0$ are parameter-free in units of Σ .

In terms of the rescaled variables eq. (2) the spectral density defined as $\rho(z) \equiv \langle \prod_{i=1}^{N} \delta^{(2)}(z-z_i) \rangle$ is given by

$$\rho^{(N_f=0)}(\xi) = (\xi^{*2} - \xi^2) w(\xi, \xi^*) \kappa(\xi, \xi^*)$$
 (3)

for quenched. The rescaled weight and kernel read

$$w(\xi, \xi^*) = \frac{1}{32\alpha^4} |\xi|^2 K_{2\nu} \left(\frac{|\xi|^2}{2\alpha^2}\right) e^{\frac{1}{4\alpha^2}(\xi^2 + \xi^{*2})}, \quad (4)$$

$$\kappa(\xi, \zeta^*) = \int_0^1 ds \int_0^1 \frac{dt}{\sqrt{t}} e^{-2s(1+t)\alpha^2}$$

$$\times (J_{2\nu}(2\sqrt{st}\xi) J_{2\nu}(2\sqrt{s}\zeta^*) - (\xi \leftrightarrow \zeta^*)), \quad (5)$$

containing K- and J-Bessel functions. The weight for the eigenvalues is non-Gaussian due to the non-trivial decoupling of eigenvectors. The unquenched density

$$\rho^{(N_f=2)}(\xi) = \rho^{(0)}(\xi) - (\xi^{*2} - \xi^2)w(\xi, \xi^*) \times \frac{|\kappa(\xi, i\eta)|^2 - |\kappa(\xi^*, i\eta)|^2}{\kappa(i\eta, i\eta)},$$
(6)

with a mass dependent correction term is shown in Fig. 1 left (for more details, see [13]). Compared to $\mu = 0$ in [15] the peaks locating the eigenvalues are split in two,

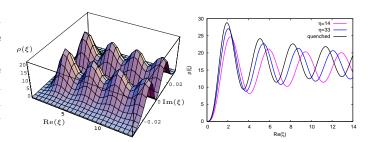


FIG. 1: The spectral density with $N_f = 2$ eq. (6) at $\alpha = 0.012$, $\eta = 8.74$ and $\nu = 0$ (left), and a cut parallel to the x-axis for different masses (right).

the particularity of this χSB class. A cut parallel to the x-axis along the maxima illustrates the effect of dynamical flavours at increasing η . At value $\eta \approx 100$ the smallest eigenvalues are effectively quenched, following eq. (3). Increasing α rapidly damps the oscillations. In

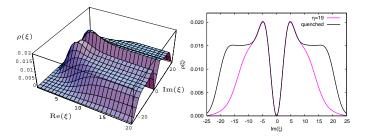


FIG. 2: The density eq. (6) at $\alpha = 3.28$, $\eta = 19$, and $\nu = 0$ (left), and a cut parallel to the y-axis (right) vs. quenched.

Fig. 2 the density is now constant along a stripe, as expected from mean field. The effect of the additional zeros of $\rho^{(N_f=2)}(\xi)$ at $\xi=\pm i\eta$ in the left corners of Fig. 2 are clearly visible. Here we choose a cut in y-direction on the maxima to highlight the difference to quenched. At this value of α the perpendicular cut in x-direction shows no dependence on η , in contrast to the previous figure. In the strong non-Hermiticity limit $\alpha\to\infty$ the plateau extends to ∞ in both x- and y-direction [13]. Away from the edge Fig. 2 is already very close to $\alpha=\infty$, to which we compared our preliminary data [14].

We now turn to the LGT side, describing the details of our simulations summarised in Table I. Our data were generated for gauge group SU(2) with $N_f = 2$ unimproved staggered fermion flavours of equal mass at coupling constant $\beta = 4/g^2 = 1.3$, using the code of [12]. In this setup the fermion determinant remains real and standard Monte Carlo applies. Our choice of parameters was dictated by several constraints: i) Wilson fermions explicitly break chiral symmetry and are already complex at $\mu = 0$, making the effect of $\mu \neq 0$ difficult to disentangle. At $\mu = 0$ Ginsparg-Wilson type fermions have been shown to exactly preserve chiral symmetry and topology (see [20] for references and a comparison to MM). Similar results have been obtained with improved staggered fermions for QCD [21]. Apart from being very expensive it is conceptually not clear so far how to extend these

TABLE I: Summary of simulation parameters at $\beta = 1.3$

\overline{V}	μ	mass ma	level spacing d	η	α	F_{π}^2/Σ	E_{Th}	no. of config.
6^4	$1.0 \cdot 10^{-3}$	0.035	$8.0(2) \cdot 10^{-3}$	13.7	0.0124(3)	0.39(3)	~ 17	45000
6^{4}	$1.0 \cdot 10^{-3}$	0.07	$6.6(2) \cdot 10^{-3}$	33.3	0.0127(3)	0.34(3)	~ 25	15000
6^{4}	$1.0 \cdot 10^{-3}$	20	$5.0(2)\cdot 10^{-3}$	12566	0.0153(3)	0.37(3)	> 100	10000
6^{4}	0.2	0.06	$9.9(1) \cdot 10^{-3}$	19.0	2.21(4)	0.38(1)	~ 10	20000
6^{4}	0.2	20	$4.5(1) \cdot 10^{-3}$	13962	3.28(4)	0.38(2)	> 200	20000
8^4	$5.625 \cdot 10^{-4}$	0.035	$3.0(3) \cdot 10^{-3}$	35.5	0.0107(9)	0.35(5)	~ 20	8 000
8^{4}	$5.625 \cdot 10^{-4}$	0.07	$2.1(1) \cdot 10^{-3}$	104.7	0.0124(3)	0.32(4)	~ 35	5000
8^{4}	$5.625 \cdot 10^{-4}$	20	$1.6(1) \cdot 10^{-3}$	39270	0.0152(3)	0.37(4)	> 35	3000
8^4	0.1125	20	$1.4(1) \cdot 10^{-3}$	44880	3.29(4)	0.38(5)	> 40	6 000

results to $\mu \neq 0$. Thus we use the Dirac operator of standard, unimproved staggered fermions:

$$D_{x,y} = \frac{1}{2a} \Big[\sum_{\gamma = \hat{x}, \hat{y}, \hat{z}} (U_{\gamma}(x) \kappa_{\gamma}(x) \delta_{y,x+\gamma} - \text{h.c.}) \quad (7)$$

$$+ \! \left(U_{\hat{t}}(x) \kappa_{\hat{t}}(x) e^{\mu} \delta_{y,x+\hat{t}} \! - \! U_{\hat{t}}^{\dagger}(y) \kappa_{\hat{t}}(y) e^{-\mu} \delta_{y,x-\hat{t}} \right) \! \Big],$$

with link variables U and staggered phases κ . Consequently our simulations are topology blind, and we set $\nu = 0$ in the following. We note that even at $\mu = 0$ the transition from staggered to the correct continuum symmetry class has not yet been observed. ii) A proper resolution of complex eigenvalues requires of the order of 10⁴ configurations for each parameter set. Even with cheap staggered fermions we are confined to small lattices $V = 6^4$ and 8^4 . iii) We want to make the window where MM apply as large as possible. At the Thouless energy scale [22], $E_{Th} \sim F_{\pi}^2/\Sigma\sqrt{V}$, the Dirac eigenvalues start to feel the propagating Goldstone modes, and leave the ϵ -regime. Increasing β (decreasing F_{π}) at fixed Vfewer Dirac eigenvalues follow MM statistics. This leads us to rather strong coupling. iv) Small masses m considerably slow down the computation. Because of rescaling $\eta = m\Sigma V$ this effects our choice of the volume too.

To compare our data with the prediction eq. (6) we have to determine two parameters from the data. First, we measure the mean level-spacing $d \sim 1/\rho(0)$, by using the Banks-Casher relation at $\mu = 0$, $\pi \rho(0) = \Sigma V$, where $\rho(0)$ is the mean spectral density. We project the complex eigenvalues onto the real axis and measure their distance. Due to $\mu \ll 1$ the 2d geometric distance between eigenvalues agrees within errors. This provides the volume rescaling of lattice eigenvalues, masses and μ^2 :

$$za \cdot \pi/d \equiv \xi$$
, $ma \cdot \pi/d \equiv \eta$, and $\mu^2 C \cdot \pi/d \equiv \alpha^2$. (8)

Second, the remaining constant $C \sim F_\pi^2/\Sigma$ is obtained by a fit to the data. We cut the data rescaled as in eq. (8) parallel to the y-axis on the first maxima, which are easy to identify for all α . Then we fit to the integral of the analytic curve, thus eliminating the choice of bins in this direction (for clarity we show histograms in all Figures). This procedure avoids cumbersome 3d fits and fixes the normalisation of the first maxima.

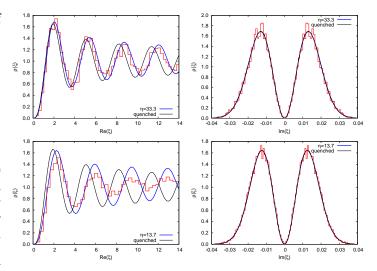


FIG. 3: The effect of dynamical fermions at $\mu=10^{-3}$, $V=6^4$: $\alpha=0.0127$, $\eta=33.3$ (upper), and $\alpha=0.0124$, $\eta=13.7$ (lower).

We start with $\alpha \ll 1$. For mass $\eta = 33.3$ in Fig. 3 (up) the data show an excellent match to the prediction eq. (6) including the 4th maximum, clearly deviating from the quenched curve shown for comparison. The perpendicular cut on the first maxima (up right) is perfectly agreeing, allowing to fit α to an relative error of $\approx 2\%$. The repulsion from the x-axis correctly identifies this χ SB class (QCD has a single peak, see [16]), and we have observed this pattern down to $\mu = 10^{-6}$. Note that fitting α does not move the maxima in the left curves. A similar data set for $V = 8^4$ leads to the same conclusions (see Table 1). Fig. 3 (down) shows the smallest

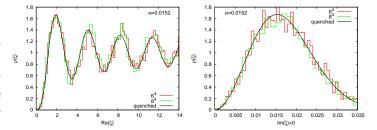


FIG. 4: The scaling $\mu_i^2 V_i = \text{const.}$: $V_1 = 6^4$, $\mu_1 = 10^{-3}$ (green) vs. $V_2 = 8^4$, $\mu_2 = 5,625 \cdot 10^{-4}$ (red), both at ma = 20 for y > 0.

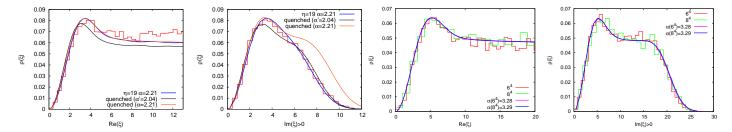


FIG. 5: The effect of dynamical fermions at $\mu = 0.2$, $V = 6^4$: $\alpha = 2.21$, $\eta = 19.0$ for y > 0. Note that $E_{Th} \sim 10$.

mass we could simulate, with $\eta = 13.7$. The data follow the analytic curve including the 2nd minimum, deviating substantially from quenched. The onset of $E_{Th} \sim 17$ is responsible for a smaller window of agreement ($E_{Th} \sim 25$ at $\eta = 33.3$). The perpendicular cut shows again perfect agreement. Despite excessive statistics the fluctuations on the maxima are still visible. To exclude that the observed deviations from quenched are finite volume effects we compare to [15] at $\mu = 0$ and otherwise identical parameters. There finite volume shifts the data to the left of the MM curve. Thus for larger volumes the deviation from guenched would be even more pronounced. Second we check the scaling hypothesis from MM and $\epsilon \chi PT$, choosing $\mu_1^2/\mu_2^2 = 6^4/8^4$ with $\mu_i^2 V_i = \text{const.}$ For simplicity [23] we have chosen a large η to effectively quench the small eigenvalues. In Fig. 4 both sets of data show an excellent agreement to the same curve from eq. (3). Independent fits for α agree within errorbars. Here only y > 0 is shown for a better resolution, the second peak at y < 0 follows from symmetry. The same results are obtained for smaller, unquenched masses (see Table 1).

Next we turn to a larger μ , remaining below the phase transition $\mu_c \approx 0.3$ found in [24] for $V = 6^4$, $\beta = 1.3$. Above the transition at $\mu = 0.4$ the data deviate from the MM as expected. We first determine d, which fixes the scale and η . At these parameter values unquenching is seen in y-direction (Fig. 5 right), which is used to fit α . Here both α and the given η influence the width. We therefore compare the unquenched curve to the quenched one with $\alpha' \neq \alpha$ fitted independently (lower curve). For comparison we also give the quenched curve at the same α (upper curve). Again we display y > 0 only. Clearly the unquenched curve describes the data best both in xand y-direction. The onset of E_{Th} is responsible for the deviation between data and MM for x > 10 in Fig. 5 (left). In Fig. 6 the scaling of constant $\mu_i^2 V_i$ is tested for these μ -values, where we have again chosen $\eta_i \gg 100$ to effectively guench. Both data sets nicely follow the same curve, the fits for α agree within errorbars.

To conclude we have shown that our unquenched SU(2) lattice data are quantitatively very well described by complex MM predictions over a wide range of chemical potentials and masses. We find agreement with $\epsilon\chi PT$ by confirming the scaling $F_\pi^2 V \mu^2$, with a consistent value for $F_\pi^2/\Sigma \sim 0.37$ for all our data. This further indicates

FIG. 6: The scaling of $\mu_i^2 V_i = \text{const.}$: $V_1 = 6^4$, $\mu_1 = 0.2$ (red) vs. $V_2 = 8^4$, $\mu_2 = 0.1125$ (green), both at ma = 20 for y > 0.

a MM- $\epsilon \chi$ PT equivalence for SU(2) at $\mu \neq 0$.

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