Wigner surmise for Hermitian and non-Hermitian chiral random matrices

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We use the idea of a Wigner surmise to compute approximate distributions of the first eigenvalue in chiral random matrix theory, for both real and complex eigenvalues. Testing against known results for zero and maximal non-Hermiticity in the microscopic large-*N* limit, we find an excellent agreement valid for a small number of exact zero eigenvalues. Compact expressions are derived for real eigenvalues in the orthogonal and symplectic classes and at intermediate non-Hermiticity for the unitary and symplectic classes. Such individual Dirac eigenvalue distributions are a useful tool in lattice gauge theory, and we illustrate this by showing that our results can describe data from two-color quantum chromodynamics simulations with chemical potential in the symplectic class.

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I. MOTIVATION

Probably one of the most used predictions of random matrix theory (RMT) is the so-called Wigner surmise (WS) describing the universal repulsion of energy levels in many systems in nature, including neutron scattering, quantum billiards, and elastomechanical modes in crystals [1]. For large matrices, the nearest-neighbor (nn) spacing distribution $p^{(\beta)}(s)$ is universal and only depends on the repulsion strength which takes discrete values $\beta=1,2,4$ for the three classical Wigner-Dyson (WD) ensembles. It can be computed with surprising accuracy using 2×2 matrices, which is the WS. Although simple arguments discussed in [2] lead to this rule for $\beta=1$, such an approximation is by no means obvious.

The extension from WD to non-Hermitian RMT introduced long ago by Ginibre [3] has become a very active field in the past decade, in particular due to applications in open quantum systems, see [4] for references and other applications. Here the spacing is known only for the class with broken time reversal (β =2) and has been applied in lattice gauge theory (LGT) [5]. However, a simple surmise based on 2×2 matrices does not work here.

In this Rapid Communication we investigate the existence of a surmise for the smallest eigenvalue in chiral RMT and its non-Hermitian extensions. These have become relevant due to applications in quantum chromodynamics (QCD) initiated by [6] and extended to non-Hermitian QCD at finite quark chemical potential μ [7]. QCD at strong coupling is a notoriously difficult theory, and the chiral RMT approach has become an important tool for LGT with exact chiral fermions [8,9]. For non-Hermitian QCD the complex action hampers a straightforward LGT approach, see [10] for a recent discussion and references. Here RMT predictions remain possible for various quantities [11–13].

In this Rapid Communication we will show that an excellent approximation for the first nonzero eigenvalue is possible using a simple $2 \times (2+\nu)$ matrix calculation, capturing the repulsion of a small number ν of zero eigenvalues. Being localized and nonoscillatory the first eigenvalue is much more suitable for LGT than the spectral density, compare,

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e.g., [9,14]. Our surmise fills some gaps in predictions for real eigenvalues in the orthogonal and symplectic classes $(\beta=1,4)$ [15], where until very recently numerically generated RMT had to be used for comparison [16]. We also provide predictions for intermediate non-Hermiticity and test them against QCD-like LGT data from [17]. This further completes the picture, compared to previous approximations [14] $(\beta=2)$ based on a Fredholm determinant expansion [18] and exact results at maximal non-Hermiticity [19] $(\beta=2,4)$.

II. LEVEL SPACING IN THE WD CLASS

We recall here the success of a WS for Hermitian and its failure for non-Hermitian WD ensembles. The WD partition function for an $N \times N$ Hermitian matrix H with real, complex or quaternion real entries is given terms of eigenvalues by

$$\mathcal{Z}_{WD}^{(\beta)} = \int dH e^{-\mathrm{Tr} \, H H^{\dagger}} \sim \int_{\mathbb{R}} \prod_{j=1}^{N} d\lambda_{j} e^{-\lambda_{j}^{2}} |\Delta_{N}(\lambda)|^{\beta}.$$
(1)

The Jacobians of the corresponding ensembles which are called GOE, GUE and GSE ($\beta = 1, 2, \text{ and } 4$) include the Vandermonde determinant, $\Delta_N(\lambda) \equiv \prod_{k>l}^N (\lambda_k - \lambda_l)$.

The large-*N* nn spacing in the bulk of the spectrum can be computed approximately from N=2 (WS) by inserting $\delta(|\lambda_1 - \lambda_2| - s)$ in $\mathcal{Z}_{WD}^{(\beta)}$:

$$p_{WS}^{(\beta)}(s) = a_{\beta}s^{\beta} \exp[-b_{\beta}s^{2}].$$
 (2)

The constants a_{β} and b_{β} follow from fixing the norm and first moment to unity (see, e.g., in [1]). The latter can always be achieved from $\int_0^{\infty} dss \hat{p}(s) = m$ by rescaling $p^{(\beta)}(s) = m\hat{p}^{(\beta)}(ms)$. This fixes the scale compared with $N = \infty$.

The exact result $p^{(\beta)}(s)$ is cumbersome, given in terms of an infinite product of eigenvalues of spheroidal functions (e.g., in [2]), the fifth Painlevé transcendent [2], or combining a Taylor series with coefficients given by sums over permutations and Dyson's asymptotic expansion in a Padé approximation [20]. This is compared to the surmise Eq. (2) in Fig. 1 left. In Table I we give the root of the integrated square deviation for later comparison,

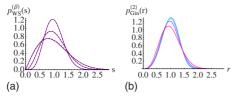


FIG. 1. (Color online) (a) Surmise $p_{WS}^{(\beta)}(s)$ (red) vs exact result [20] (dashed blue) from $\beta=1,2,4$ (bottom to top). (b) $p_{Gin}^{(2)}(r)$ for N=2,3,4,20 in red to blue (bottom to top).

$$\delta = \left\{ \int_0^\infty ds [p^{(\beta)}(s) - p^{(\beta)}_{surmise}(s)]^2 \right\}^{1/2}.$$
 (3)

The non-Hermitian WD ensembles are defined by dropping the Hermiticity constraint in Eq. (1) left [3]. We only display the complex eigenvalue representation for $\beta=2$ and 4 and their Jacobians $\mathcal{J}_{\beta}(z)$ computed in [3]:

$$\mathcal{Z}_{Gin}^{(\beta)} = \int_{\mathbb{C}} \prod_{j=1}^{N} d^{2} z_{j} e^{-|z_{j}|^{2}} \mathcal{J}_{\beta}(z),$$
$$\mathcal{J}_{2}(z) = |\Delta_{N}(z)|^{2}, \quad \mathcal{J}_{4}(z) = \Delta_{2N}(z, z^{*}) \prod_{j=1}^{N} (z_{j} - z_{j}^{*}).$$
(4)

For $\beta=2$ the spacing is obtained from an N=2 surmise by inserting $\delta(|z_1-z_2|-r)$ in $\mathcal{Z}_{Gin}^{(2)}$ and putting one eigenvalue at the origin. The exact spacing for any *N* obtained in [21] uses translational invariance in the bulk

$$\hat{p}_{Gin}^{(2)}(r) = -\frac{\partial E_{Gin}^{(2)}(r)}{\partial r}, \quad E_{Gin}^{(2)}(r) = \prod_{j=1}^{N-1} e^{-r^2} \sum_{k=0}^{j} \frac{r^{2k}}{k!}.$$
 (5)

In Fig. 1 right we compare N=2 with increasing N, all curves having norm and first moment 1. Clearly a surmise *does not work* for the $\beta=2$ Ginibre ensemble ($\delta \approx 0.18$), as previously noted in [21]. For $\beta=4$ and 1 the spacing is currently unknown.

III. FIRST EIGENVALUE IN CHIRAL RMT

The chiral ensembles with real eigenvalues called chGOE, chGUE, and chGSE are defined in terms of $N \times (N + \nu)$ rectangular matrices W with real, complex or quaternion real elements without further symmetry restrictions. Switching to positive eigenvalues $\lambda_j \ge 0$ of the Hermitian Wishart (or covariance) matrix WW^{\dagger} , we obtain

TABLE I. Deviation Eq. (3) in units 10^{-3} between approximate N=2 and exact large-N results (δ_{WS} from [20]).

	δ_{WS}		$\delta_{\mu=0}^{\nu=0}$	$\delta_{\mu=0}^{\nu=1}$	$\delta_{\mu=0}^{\nu=2}$	$\delta_{\mu=1}^{\nu=0}$	$\delta_{\mu=1}^{\nu=1}$	$\delta_{\mu=1}^{\nu=2}$
GUE	0.04	chGUE	0	3.8	7.7	8.0	12.3	14.8
GSE	0.015	chGSE	1.7	6.1	10.6	1.8	3.3	4.4
GOE	0.16	chGOE	3.6	0				

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$$\mathcal{Z}_{\nu}^{(\beta)} = \int_{0}^{\infty} \prod_{j=1}^{N} d\lambda_{j} \lambda_{j}^{d} e^{-\lambda_{j}} |\Delta_{N}(\lambda)|^{\beta}, \quad d \equiv \frac{\beta(\nu+1)}{2} - 1.$$
(6)

Here N_f massless flavors can be added by shifting $d \rightarrow d + N_f$. The gap probability $E^{(\beta)}(s)$ that the interval (0,s) is void follows by integrating in Eq. (6) from s to ∞ . For N=2 we obtain

$$E_{\nu}^{(\beta)}(s) \sim \int_{0}^{\infty} dx dy [(x+s)(y+x+s)]^{d} e^{-2(s+x)-y} y^{\beta}, \quad (7)$$

after shifting variables. The nested integrals can easily be evaluated. Note that d=0 for $\beta=2$, $\nu=0$, and $\beta=1=\nu$. These gap probabilities can be computed exactly for any *N*, and our surmise gives the *exact* result after rescaling.

To compare with Dirac operator eigenvalues we have to switch variables $\lambda_j \rightarrow y_j^2$, coming in eigenvalue pairs $\pm y_j$, and thus to $s \rightarrow s^2$. The distribution of the first positive Dirac eigenvalue follows: $p_{\nu}^{(\beta)}(s) = -\partial_s [E_{\nu}^{(\beta)}(s^2)]$.

We first list all its known $N_f=0$ results in the universal microscopic limit for $\nu \in \mathbb{N}$ in Eqs. (8)–(10): the chGUE for all ν [15,22], the chGOE for $\nu=0$ [23] and odd ν [15], and the chGSE for $\nu=0$ [24] and $\nu>0$ [25]. For the latter, only a convergent Taylor series is known with coefficients $a_j(\nu)$ given by sums over partitions [see Eq. (8) in [25]], much alike for the WS in the WD class,

$$p_{\nu}^{(2)}(s) = s e^{-s^2/4} \det_{i,j=1,\dots,\nu} [I_{i-j+2}(s)]/2, \qquad (8)$$

$$p_{\nu=0}^{(1)}(s) = [(2+s)e^{-s^2/8-s/2}]/4,$$

$$\hat{p}_{\nu=2n+1}^{(1)}(s) \sim s^{(3-\nu)/2}e^{-s^2/8} \Pr[(i-j)I_{i+j+3}(s)], \qquad (9)$$

$$i,j=-n+\frac{1}{2},...,n-\frac{1}{2}$$

$$p_{\nu=0}^{(4)}(s) = (\pi/2)^{1/2}s^{3/2}e^{-s^2/2}I_{3/2}(s),$$

$$\hat{p}_{\nu>0}^{(4)}(s) \sim s^{4\nu+3}e^{-s^2/2} \left(1+\sum_{i=1}^{\infty}a_j(\nu)s^j\right). \qquad (10)$$

Next, we give examples following our surmise Eq. (7) where $p_{\nu}^{\beta}(s)$ is *not* known in elementary form, filling the gaps in Eqs. (8)–(10) for the first two values of $\nu > 0$:

$$\hat{p}_{\nu=2}^{(1)}(s) \sim 3s^3 e^{-s^2/8} + \left(6s^2 - \frac{s^4}{4}\right) e^{-s^2/16} \sqrt{\pi} \operatorname{Erfc}\left[\frac{s}{4}\right],$$
(11)

$$\hat{p}_{\nu=4}^{(1)}(s) \sim \left(s^5 + \frac{s^7}{60}\right) e^{-s^{2}/8} + \left(2s^4 - \frac{s^6}{20}\right) e^{-s^{2}/16} \sqrt{\pi} \text{Erfc}\left[\frac{s}{4}\right],$$

$$\hat{p}_{\nu=1}^{(4)}(s) \sim s^7 (13\ 440 + 1440s^2 + 60s^4 + s^6) e^{-s^{2}/2},$$

$$\hat{p}_{\nu=2}^{(4)}(s) \sim s^{11} (15\ 482\ 880 + 2\ 150\ 400s^2 + 134\ 400s^4 + 4800s^6 + 100s^8 + s^{10}) e^{-s^{2}/2}.$$
(12)

The normalization constants suppressed above easily follow.

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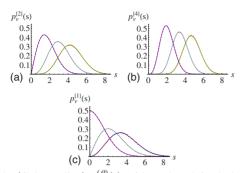


FIG. 2. (Color online) $p_{\nu}^{(\beta)}(s)$ with $\nu=0,1,2$ in dashed blue to green (left to right) for (a) $\beta=2$, (b) $\beta=4$, and (c) $\beta=1$ vs N=2 surmise is in red. Our result Eq. (11) is compared to a numerical simulation at N=20 (black dots).

However, we cannot set the first moment to one as in the WD class. The position of $p_{\nu}^{(\beta)}(s)$ measures the repulsion by ν exact zero eigenvalues, containing important information. Thus we fix the N=2 scale by setting the 1st moment equal to the exact one. Without exact ($\beta=1$, even ν) or concise ($\beta=4, \nu>0$) results, we instead fit to the increasing slope of the known microscopic density $\rho_{\nu}^{(\beta)}(s)$, being the first term in the Fredholm expansion of the first eigenvalue [18] [see also Eq. (18)]. In Fig. 2 we compare approximate to exact first eigenvalues for small topology $\nu=0,1,2$ and all β . The deviation measured by Eq. (3) in Table I increases with ν , becoming visible only for $\nu=2$ (see Fig. 2). This has to be compared to the statistical error in data, see, e.g., Fig. 5.

Note that in chiral RMT the nn spacing also obeys Eq. (2) but *does not follow* from an N=2 surmise [26].

The non-Hermitian chiral ensembles with $\mu \neq 0$ are given in terms of a two-matrix model [11,27]. We only focus on $\beta = 2,4$ here, with their complex eigenvalue representations for $N_f=0$ reading [11,27]

$$\mathcal{Z}_{\nu\mathbb{C}}^{(\beta)} = \int_{\mathbb{C}} \prod_{j=1}^{N} d^2 z_j |z_j|^{\beta\nu+2} K_{\beta\nu/2}(a|z_j|^2) e^{b\Re e z_j^2} \mathcal{J}_{\beta}(z^2).$$
(13)

The weight w(z) depends on $a \equiv (1+\mu^2)/2\mu^2 > b \equiv (1-\mu^2)/2\mu^2 \ge 0$, with $\mu \in (0,1]$. The limit $\mu \to 0$ leads back to real eigenvalues, and at $\mu=1$ non-Hermiticity is maximal. The definition of a gap probability on C is not unique [14,19]. For *radial* ordering it reads

$$E^{(\beta)}(r) \sim \prod_{j=1}^{N} \int_{r}^{\infty} dr_{j} r_{j} \int_{0}^{2\pi} d\theta_{j} w(z_{j}) \mathcal{J}_{\beta}(z).$$
(14)

Differentiation yields $\partial_r E^{(\beta)}(r) = \int_0^{2\pi} d\theta p_{\nu}^{(\beta)}(re^{i\theta}) \equiv P_{\nu}^{(\beta)}(r)$, the integrated first eigenvalue. For $\beta = 2$ ($\beta = 4$) Eq. (14) is given by a Fredholm determinant (Pfaffian) [19]

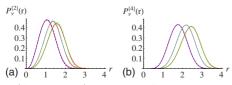


FIG. 3. (Color online) Integrated first eigenvalue $P_{\nu}^{(\beta)}(r)$ at $\mu=1$ for $\nu=0,1,2$ in blue to green dashes (left to right) vs N=2 (red): (a) $\beta=2$ and (b) $\beta=4$.

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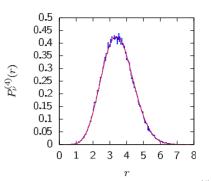


FIG. 4. (Color online) Integrated first eigenvalue $P_{\nu}^{(4)}(r)$ (red) vs lattice data [17] (blue histograms) with volume $V=4^4$, gauge coupling 1.3, $\mu_{\text{Lat}}=0.2$ and mass 20 in lattice units, using a very large number 10^5 configurations.

$$E^{(2)}(r) \sim \det_{1,\dots,N} \left[\int_{r^2}^{\infty} dt t^{k+j+\nu-1} K_{\nu}(at) I_{k+j-2}(bt) \right].$$
(15)

Its matrix elements $A_{jk}^{(\nu)}$ can be computed recursively for any ν by differentiating the following matrix element [19]:

$$A_{11}^{(0)} = \frac{br^2 I_1(br^2) K_0(ar^2) + ar^2 I_0(br^2) K_1(ar^2)}{a^2 - b^2}.$$
 (16)

This leads to a $\beta \times \beta$ determinant (Pfaffian) representation for our N=2 surmise valid for any μ . At $\mu=1$ all Fredholm eigenvalues $1-\lambda_{k=0,\dots,N-1}^{(\beta)}$ are explicitly known [19], providing an exact result for any N as in Eq. (5). It contains incomplete Bessel function series $I_{\nu}^{[k]}(x)$ truncated at power $k \ (\equiv 0$ for k < 0)

$$[1 - \lambda_{k}^{(2)}] = \frac{r^{2(2k+\nu+1)}}{2^{2k+\nu}(k+\nu)!k!} K_{\nu+1}(r^{2}) + r^{2}[I_{\nu+2}^{(k-2)}(r^{2})K_{\nu+1}(r^{2}) + I_{\nu+1}^{(k-1)}(r^{2})K_{\nu+2}(r^{2})].$$
(17)

For $\beta=4$ we have the relation $\lambda_k^{(4)} = \lambda_{2k+1}^{(2)}$ with $\nu \rightarrow 2\nu$ [19]. In Fig. 3 we compare our surmise to this result, truncated at N=8 because of rapid convergence. Here it works better for $\beta=4$ than $\beta=2$, in contrast to $\mu=0$. Due to angular integration only one scale has to be fixed after normalization, which can be done as in the real case.

Next we give a surmise for $p_{\nu}^{(\beta)}(re^{i\theta})$. In Eq. (14) we skip the integration over θ_1 and differentiate wrt r_1 . For N=2 we obtain an exact Fredholm expansion

$$p_{\nu}^{(\beta)}(z) = R_{1,\nu}^{(\beta)}(z) - \int_{0}^{r_{1}} dtt \int_{0}^{2\pi} d\varphi R_{2,\nu}^{(\beta)}(z,te^{i\varphi}), \quad (18)$$

with $z=r_1e^{i\theta_1}=x+iy$. The one- and two-point spectral densities are expressed through the kernel of orthogonal Laguerre polynomials of norm h_j (see [11] for details)

$$R_{1,\nu}^{(2)}(z) = K_N^{(2)}(z, z^*) = w(z) \sum_{j=0}^{N-1} \frac{\left| L_j^{(\nu)} \left(\frac{z^2}{1 - \mu^2} \right) \right|^2}{h_j}, \quad (19)$$

and $R_{2,\nu}^{(2)}(z,u) = R_{1,\nu}^{(2)}(z)R_{1,\nu}^{(2)}(u) - |K_N^{(2)}(z,u^*)|^2$. For $\beta = 4$ we have a Pfaffian of a matrix kernel instead [27]. An example for $p_{\nu=0}^{(4)}(z)$ is shown in Fig. 5 top right. Here two scales have

to be fixed: for z we fit to the increase in the known microscopic density in the x direction and for rescaling $2N\mu^2 \equiv \alpha^2$ to its decrease in the y direction. Since $\alpha \le 2$ for N=2, we conclude that at large-N for $\alpha > 2 p_{\nu}^{(\beta)}(z)$ must become symmetric wrt rotation ($\beta=2$) or reflections wrt the bisector of each quadrant ($\beta=4$). We have checked this, as well as distributions for $0 < \mu < 1$ by generating ensembles of large random matrices.

IV. LATTICE DATA

In [17] two-color QCD was compared to the β =4 microscopic spectral density in the complex plane from chiral RMT [27]. We use the same data here but with higher statistics and refer to [17] for all simulation details. Because unimproved staggered fermions are used, we are in the β =4 class at ν =0. Our N_f =2 data are effectively quenched for the smallest eigenvalues due to a large mass. In Fig. 4 we compare to the first integrated eigenvalue, with α =1.352 being close to maximal non-Hermiticity. No further fits compared to [17] are made.

In Fig. 5 we compare LGT data at intermediate $\mu_{\text{Lat}}=0.1$ to the angle-dependent surmise Eq. (18) by taking cuts. Here two scales are fitted to the data, finding an excellent agreement for $\alpha=0.65$. An alternative to Eq. (18) is to truncate the Fredholm expansion in the large-*N* limit [18]. This was successfully applied to the $\beta=2$ class [14], but higher-order terms are cumbersome.

V. CONCLUSIONS

Conceptually it is possible within chiral RMT to approximate the first eigenvalue distribution using a $2 \times (2+\nu)$ matrix calculation for both real and complex eigenvalues. It is

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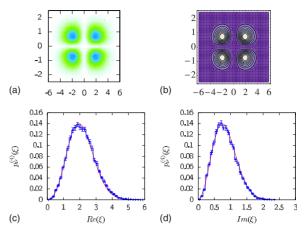


FIG. 5. (Color online) Top: contour plots for Lattice data as in Fig. 4 but with (a) $\mu_{\text{Lat}}=0.1$ vs (b) surmise Eq. (18). Bottom: a single peak cut in (c) x and (d) y directions.

remarkable that this surmise works and captures the repulsion of ν zero eigenvalues. We derived compact expressions for $\beta = 1$ and 4 with real eigenvalues for $\nu > 0$. Second, we have shown that our surmise for $\beta = 4$ successfully describes SU(2) lattice data, in an intermediate regime for $\mu \neq 0$ where no results were previously known. It would be very interesting to extend our results to the $\beta = 1$ non-Hermitian chiral class, having both real and complex eigenvalues.

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