

Individual eigenvalue distributions of crossover chiral random matrices and low-energy constants of $SU(2) \times U(1)$ lattice gauge theory

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We compute individual distributions of low-lying eigenvalues of a chiral random matrix ensemble interpolating symplectic and unitary symmetry classes by the Nyström-type method of evaluating the Fredholm Pfaffian and resolvents of the quaternion kernel. The one-parameter family of these distributions is shown to fit excellently the Dirac spectra of $SU(2)$ lattice gauge theory with a constant $U(1)$ background or dynamically fluctuating $U(1)$ gauge field, which weakly breaks the pseudoreality of the unperturbed $SU(2)$ Dirac operator. The observed linear dependence of the crossover parameter with the strength of the $U(1)$ perturbations leads to precise determination of the pseudo-scalar decay constant, as well as the chiral condensate in the effective chiral Lagrangian of the AI class.
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1. Introduction

The grounds for universality, i.e., insensitivity to details of the system of concern, of local correlation of energy levels of stochastic and quantum-chaotic Hamiltonians have been well uncovered by now, in terms of ten-fold classification of symmetric superspaces on which spectral nonlinear σ models describing spontaneous symmetry breaking reside [1], and of semiclassical equivalence between periodic orbits and the aforementioned σ models [2]. On the other hand, the presence of explicit symmetry breaking perturbations is known to induce a crossover between different universality classes [3], in such a way that is also insensitive to the systems' details. An example of this universality crossover is the Gaussian orthogonal ensemble (GOE)–Gaussian unitary ensemble (GUE) transition in a disordered or chaotic system under a magnetic field [4,5]. In the realm of lattice gauge theory, where Dirac operators play the rôle of stochastic Hamiltonians [6,7], the crossover between the chiral Gaussian unitary ensemble (chGUE) itself and the chGUE–GUE crossover, associated with an imaginary isospin chemical potential [8] and a finite lattice-spacing effect in the Wilson Dirac operator [9], respectively, have been utilized to determine the pion decay constant and the Wilsonian chiral perturbation constants from relatively small lattices. The aim of this work is to apply this spectral approach to the determination of low-energy constants in another setting, namely $SU(2)$ gauge theory under $U(1)$ perturbations, either in the form of a constant imaginary chemical potential [10,11] or a dynamically fluctuating one. In contrast to these preceding works, which used n -level correlation functions or the smallest eigenvalue distribution, our strategy in this paper is to employ *multiple* spectral observables, which allows for precise fitting of lattice data, namely, individual distributions of the k th smallest Dirac eigenvalues [12] inclusively. The practical advantages of our

method will be proved in the precision of the low-energy constants determined, as preliminarily reported in Ref. [13].

This paper is composed mainly of two parts: Sect. 2 is devoted to analytic treatments in random matrix theory, and Sect. 3 to its application to Dirac eigenvalue distributions measured in lattice simulations. In Sect. 2 we start by reviewing the established results of the chiral Gaussian symplectic ensemble (chGSE)–chGUE crossover, namely, the derivation of the quaternion kernel K [14–16]. Then we apply the Nyström-type method [17,18] to that kernel and compute individual eigenvalue distributions in the form of the Fredholm Pfaffian and resolvents of $\hat{K}$ [19]. The relationship between the chiral Lagrangian in the ε regime and the nonlinear σ model from random matrices interpolating chGSE–chGUE leads to identification between parameters in each theory. In Sect. 3 we introduce our models of lattice gauge theory and explain our strategy of fitting the Dirac spectra using individual eigenvalue distributions of chGSE and of the chGSE–chGUE crossover. Optimally fitting parameters (mean level spacings Δ and crossover parameters ρ) will be exhibited in Tables 1–6, leading to very precise determination of the low-energy constants (chiral condensate Σ and pseudo-scalar decay constant F) as presented in Table 7. Our conclusions, including a possible direction of study, will be summarized in Sect. 4.

2. chGSE–chGUE crossover

2.1. Crossover random matrix ensemble

Let N and N' be even positive integers. Let $\mathbf{A}$ and $\mathbf{B}$ be $(N/2) \times (N'/2)$ quaternion matrices, which can be represented as ordinary $N \times N'$ matrices A and B as¹

$$A = \sum_{\mu=0}^3 \left(A_{jk}^{(\mu)} \right) \otimes \sigma_{\mu}, \quad B = \sum_{\mu=0}^3 \left(B_{jk}^{(\mu)} \right) \otimes \sigma_{\mu} \quad (j = 1, \dots, N/2, k = 1, \dots, N'/2). \quad (1)$$

Here the four units of the quaternion field $\mathbb{H}$ are represented by the 2×2 unit matrix and Pauli matrices, $\{\sigma_{\mu}\} = (\mathbb{I}, -i\sigma_1, -i\sigma_2, -i\sigma_3)$. Let these matrix elements belong to

$$A_{jk}^{(\mu)} \in \mathbb{R}, \quad B_{jk}^{(\mu)} \in \mathbb{C}, \quad (2)$$

i.e., $\mathbf{A}$ is quaternion-real and $\mathbf{B}$ is not (i.e., B is a generic $N \times N'$ complex matrix). We consider $A_{jk}^{(\mu)}$, $\text{Re } B_{jk}^{(\mu)}$, and $\text{Im } B_{jk}^{(\mu)}$ to be independent random variables, distributed according to the Gaussian distributions $e^{-\frac{1}{2} \text{tr } AA^{\dagger}}$ and $e^{-\text{tr } BB^{\dagger}}$, respectively. We define an ensemble of $(N + N') \times (N + N')$ Hermitian matrices H of the form

$$H = \begin{bmatrix} 0_{N \times N} & C \\ C^{\dagger} & 0_{N' \times N'} \end{bmatrix}, \quad C = e^{-\tau} A + \sqrt{1 - e^{-2\tau}} B, \quad (3)$$

where τ is a real parameter, initially introduced by Dyson as a fictitious time for the Brownian motion of eigenvalues [3]. This ensemble is called a “chiral” random matrix ensemble because it enjoys the chiral symmetry $\{H, \gamma_5\} = 0$ with $\gamma_5 = \text{diag}(\mathbb{I}_N, -\mathbb{I}_{N'})$. This anticommutation relation implies that the eigenvalues of H consist of $\min(N', N)$ pairs of generically nonzero eigenvalues of equal magnitude and opposite signs (i.e., $(\pm 1) \times$ singular values of C), and $\nu = |N' - N|$

¹ We denote quaternions in bold symbols and their 2×2 complex matrix representatives in the corresponding italic symbols.

zero eigenvalues. The presence of B violates the quaternion-reality of C and the self-duality of H (i.e., $H_{ij}^{(\mu)} \otimes \sigma_\mu = H_{ji}^{(\mu)} \otimes \sigma_\mu^\dagger$, $i, j = 1, \dots, N + N'$), and lifts the Kramers degeneracy of all nonzero eigenvalues of H (i.e., singular values of A alone). Accordingly, this random matrix ensemble interpolates between two limiting cases, the chiral Gaussian symplectic ensemble (chGSE) at $\tau = 0$ and the chiral Gaussian unitary ensemble (chGUE) at $\tau \rightarrow \infty$, depending on a single parameter τ (at a fixed, finite N and N').

2.2. Joint eigenvalue distribution

In order to make the paper self-contained, below we sketch the derivation of the probability distribution of singular values of C , and refer the reader to Refs. [14–16] for rigorous proofs of the relevant formulas. We start from the unnormalized probability measure of the matrix elements of C ,

$$dC \int dA e^{-\frac{1}{2} \text{tr} AA^\dagger} \int dB e^{-\text{tr} BB^\dagger} \delta \left(e^{-\tau} A + \sqrt{1 - e^{-2\tau}} B - C \right) \\ \propto dC \int dA \exp \left(-\frac{1}{2} \text{tr} AA^\dagger - \frac{1}{1 - e^{-2\tau}} \text{tr} (C - e^{-\tau} A)(C - e^{-\tau} A)^\dagger \right), \quad (4)$$

where $dA = \prod_{j,k,\mu} dA_{jk}^{(\mu)}$, $dB = \prod_{j,k,\mu} d^2 B_{jk}^{(\mu)}$, and $dC = \prod_{j,k,\mu} d^2 C_{jk}^{(\mu)}$. Without loss of generality we assume $N \leq N'$. We employ the singular value decomposition $A = SMS'^\dagger$ and $C = U\Lambda U'^\dagger$ with $S \in \text{USp}(N)$, $S' \in \text{USp}(N')$, $U \in \text{U}(N)$, $U' \in \text{U}(N')$ and parametrize the singular values as $M = \begin{bmatrix} \text{diag}(\mu_1, \dots, \mu_N) & 0_{N \times v} \end{bmatrix}$, $\Lambda = \begin{bmatrix} \text{diag}(\lambda_1, \dots, \lambda_N) & 0_{N \times v} \end{bmatrix}$. Kramers-degenerate pairs of singular values of A are ordered such that $\mu_{i+N/2} = \mu_i$ ($i = 1, \dots, N/2$). The measures dA and dB on quaternion-real matrices and complex matrices take the following respective forms (in what follows we suppress all constant factors in the measures):

$$dA = d(S, S') \left(\prod_{i=1}^{N/2} d\mu_i \mu_i^{2v+3} \right) \Delta_{N/2}(\mu^2)^4, \quad dC = d(U, U') \left(\prod_{i=1}^N d\lambda_i \lambda_i^{2v+1} \right) \Delta_N(\lambda^2)^2. \quad (5)$$

Here $d(S, S')$ and $d(U, U')$ denote the invariant measures on the respective angular degrees of freedom, and Δ denote the Vandermonde determinants $\Delta_{N/2}(\mu^2) := \prod_{i>j}^{N/2} (\mu_i^2 - \mu_j^2)$ and $\Delta_N(\lambda^2) := \prod_{i>j}^N (\lambda_i^2 - \lambda_j^2)$. The probability measure of the singular values $\{\lambda_i\}$ of C follows from Eqs. (4) and (5) by integrating out the unitary matrices (U, U') . The integrations over symplectic matrices (S, S') decouple after redefining the unitary matrices by $S^\dagger U \rightarrow U$ and $S'^\dagger U' \rightarrow U'$, leading to the expression

$$\text{Eq. (4)} = \prod_{i=1}^N d\lambda_i \lambda_i^{2v+1} \exp \left(-\frac{\lambda_i^2}{1 - e^{-2\tau}} \right) \Delta_N(\lambda^2)^2 \int_0^\infty \prod_{j=1}^{N/2} d\mu_j \mu_j^{2v+3} \exp \left(-\frac{\mu_j^2}{\tanh \tau} \right) \\ \times \Delta_{N/2}(\mu^2)^4 \int_{\text{U}(N)} dU \int_{\text{U}(N')} dU' \exp \left(\frac{1}{\sinh \tau} \text{Re tr } U \Lambda U'^\dagger M^t \right). \quad (6)$$

We employ the Berezin–Karpelevich formula [20–22] for the integration over (U, U') and take the pairwise confluent limit $\mu_{i+N/2} \rightarrow \mu_i$ for all $i = 1, \dots, N/2$:

$$\begin{aligned} \int_{U(N)} dU \int_{U(N')} dU' e^{\frac{1}{\sinh \tau} \text{Re tr } U \Lambda U'^{\dagger} M^t} &\propto \frac{\det \left[I_{\nu} \left(\frac{\lambda_i \mu_j}{\sinh \tau} \right) \right]_{i,j=1}^N}{\Delta_N(\lambda^2) \Delta_N(\mu^2) \prod_{i=1}^N (\lambda_i \mu_i)^{\nu}} \Bigg|_{\mu_{i+N/2} \rightarrow \mu_i} \\ &\propto \frac{\det \left[I_{\nu} \left(\frac{\lambda_i \mu_j}{\sinh \tau} \right) \frac{\partial}{\partial \mu_j^2} I_{\nu} \left(\frac{\lambda_i \mu_j}{\sinh \tau} \right) \right]_{i=1, \dots, N}^{j=1, \dots, N/2}}{\Delta_N(\lambda^2) \Delta_{N/2}(\mu^2)^4 \prod_{i=1}^N \lambda_i^{\nu} \prod_{j=1}^{N/2} \mu_j^{2\nu}}. \end{aligned} \quad (7)$$

Here I_{ν} denotes the Bessel function of the pure imaginary argument, $I_{\nu}(z) = J_{\nu}(iz)$. By substituting Eq. (7) into Eq. (6) and performing a change of variables from the singular values of rectangular matrices C and A to the eigenvalues of Wishart matrices $CC^{\dagger}$ and $AA^{\dagger}$, $x_i = \lambda_i^2$ and $y_i = \mu_i^2$, the probability measure becomes proportional to

$$\text{Eq. (6)} \propto \prod_{i=1}^N dx_i \sqrt{w(x_i)} \Delta_N(x) \int_0^{\infty} \prod_{j=1}^{N/2} dy_j y_j \det \left[g(x_k, y_{\ell}) \frac{\partial g(x_k, y_{\ell})}{\partial y_{\ell}} \right]_{k,\ell=1, \dots, N/2}^{k=1, \dots, N}. \quad (8)$$

Here we have introduced the Laguerre weight $w(x) := x^{\nu} e^{-x}$, and the symmetric function

$$g(x, y) := \frac{e^{-(2\nu+1)\tau}}{1 - e^{-2\tau}} \exp \left(-\frac{x+y}{2 \tanh \tau} \right) I_{\nu} \left(\frac{\sqrt{xy}}{\sinh \tau} \right). \quad (9)$$

With the multiplicative constant chosen as the above, the function $g(x, y)$ admits an alternative interpretation as a one-particle Green’s function for the Brownian motion at time τ ,

$$g(x, y) = \sqrt{w(x)w(y)} \sum_{k=0}^{\infty} \frac{L_k^{\nu}(x)L_k^{\nu}(y)}{h_k} e^{-\gamma_k \tau}, \quad (10)$$

with the norm given by $h_k = (k + \nu)!/k!$ and the “one-particle energy” by $\gamma_k = 2k + \nu + 1$. Using a lemma by Mehta (A.17 of Ref. [23]), the $(N/2)$ -fold integral of an $N \times N$ determinant in Eq. (8) can be decomposed into an $N \times N$ Pfaffian of single integrals:

$$\text{Eq. (8)} \propto \prod_{i=1}^N dx_i \sqrt{w(x_i)} \Delta_N(x) \text{Pf} [F(x_j, x_k)]_{j,k=1, \dots, N}, \quad (11)$$

$$F(x, x') := \int_0^{\infty} dy y \left\{ g(x, y) \frac{\partial g(x', y)}{\partial y} - \frac{\partial g(x, y)}{\partial y} g(x', y) \right\}. \quad (12)$$

2.3. Quaternion determinant

In this subsection we summarize the procedure presented in Ref. [23], Sect. 14. We introduce a set of arbitrary monic polynomials $\{R_k(x)\}_{k=0,1,\dots}$ and arbitrary positive numbers $\{r_k\}_{k=0,1,\dots}$, and define functions $\{\psi_k(x)\}$ by

$$\psi_{2k}(x) = \frac{\sqrt{w(x)} R_{2k}(x)}{\sqrt{r_k}}, \quad \psi_{2k+1}(x) = \frac{\sqrt{w(x)} R_{2k+1}(x)}{\sqrt{r_k}} \quad (13)$$

and their F -convolutions $\{\phi_k(x)\}$ by

$$\phi_k(x) = - \int_0^\infty dx' F(x, x') \psi_k(x'). \quad (14)$$

Next we introduce functions $D(x, x')$, $S(x, x')$, and $I(x, x')$, which are bilinear combinations of $\{\psi_k(x)\}$ and $\{\phi_k(x)\}$:

$$D(x, x') = \sum_{k=0}^{N/2-1} \left(\psi_{2k}(x) \psi_{2k+1}(x') - \psi_{2k+1}(x) \psi_{2k}(x') \right), \quad (15)$$

$$S(x, x') = \sum_{k=0}^{N/2-1} \left(\phi_{2k}(x) \psi_{2k+1}(x') - \phi_{2k+1}(x) \psi_{2k}(x') \right), \quad (16)$$

$$I(x, x') = - \sum_{k=0}^{N/2-1} \left(\phi_{2k}(x) \phi_{2k+1}(x') - \phi_{2k+1}(x) \phi_{2k}(x') \right), \quad (17)$$

and the corresponding $N \times N$ matrices D_N , S_N , and I_N by

$$D_N = [D(x_i, x_j)]_{i,j=1,\dots,N}, \quad S_N = [S(x_i, x_j)]_{i,j=1,\dots,N}, \quad I_N = [I(x_i, x_j)]_{i,j=1,\dots,N}. \quad (18)$$

Since a $2N \times 2N$ antisymmetric matrix $\begin{bmatrix} D_N & S_N^t \\ -S_N & -I_N \end{bmatrix}$ is a product of two rectangular matrices of size $2N \times N$ and $N \times 2N$:

$$\begin{bmatrix} D_N & S_N^t \\ -S_N & -I_N \end{bmatrix} = \begin{bmatrix} \psi_{2k}(x_i) & \psi_{2k+1}(x_i) \\ -\phi_{2k}(x_i) & -\phi_{2k+1}(x_i) \end{bmatrix}_{k=0,\dots,N/2-1}^{i=1,\dots,N} \begin{bmatrix} \psi_{2k+1}(x_j) & -\phi_{2k+1}(x_j) \\ -\psi_{2k}(x_j) & \phi_{2k}(x_j) \end{bmatrix}_{j=1,\dots,N}^{k=0,\dots,N/2-1},$$

its rank is N at most; and also it is N at least due to the linear independence of $\{\psi_i(x)\}_{i=0,\dots,N-1}$. Accordingly the lower N rows $[-S_N \ -I_N]$ are linear combinations of the upper N rows $[D_N \ S_N^t]$.

Now consider a Pfaffian of another antisymmetric $2N \times 2N$ matrix $\begin{bmatrix} D_N & S_N^t \\ -S_N & -I_N - F_N \end{bmatrix}$ with $F_N = [F(x_i, x_j)]_{i,j=1,\dots,N}$. Since adding $[S_N \ I_N]$ (or minus its transpose) to the lower N rows (the right N columns) does not change the determinants due to the aforementioned linear dependence with the upper N rows (the left N columns), we readily obtain

$$\text{Pf} \begin{bmatrix} D_N & S_N^t \\ -S_N & -I_N - F_N \end{bmatrix} = \text{Pf} \begin{bmatrix} D_N & 0 \\ 0 & -F_N \end{bmatrix} = (-1)^{N/2} \text{Pf } D_N \cdot \text{Pf } F_N. \quad (19)$$

On the other hand,

$$\begin{aligned} \text{Pf } D_N &= \text{Pf} \left[[\psi_{2k}(x_i) \ \psi_{2k+1}(x_i)]_{k=0,\dots,N/2-1}^{i=1,\dots,N} \begin{bmatrix} \psi_{2k+1}(x_j) \\ -\psi_{2k}(x_j) \end{bmatrix}_{j=1,\dots,N}^{k=0,\dots,N/2-1} \right] \\ &= \det [\psi_{k-1}(x_i)]_{i,k=1,\dots,N} = \det [\sqrt{w(x_i)} x_i^{k-1}]_{i,k=1,\dots,N} \propto \prod_{i=1}^N \sqrt{w(x_i)} \cdot \Delta_N(x). \end{aligned} \quad (20)$$

Using Eqs. (19) and (20), the probability measure (11) now reads

$$\text{Eq. (11)} \propto \prod_{i=1}^N dx_i \text{Pf}(ZK_N), \quad (21)$$

$$K_N = \begin{bmatrix} S_N & J_N \\ D_N & S_N^t \end{bmatrix}, \quad J_N = I_N + F_N, \quad Z = \begin{bmatrix} 0 & \mathbb{I}_N \\ -\mathbb{I}_N & 0 \end{bmatrix}.$$

Since its upper diagonal block is the transpose of the lower diagonal block and both off-diagonal blocks are antisymmetric, K_N can be regarded a $2N \times 2N$ complex matrix representative of an $N \times N$ quaternion self-dual matrix, which we call $\mathbf{K}_N = [\mathbf{K}(x_i, x_j)]_{i,j=1,\dots,N}$. Using Dyson's lemma $\text{qdet } \Phi = \text{Pf}(Z\Phi)$ for the quaternion determinant (qdet) [24] of a quaternion self-dual matrix Φ , we finally obtain

$$\text{Eq. (21)} = \prod_{i=1}^N dx_i \text{qdet} \mathbf{K}_N(x_1, \dots, x_N). \quad (22)$$

2.4. Skew-orthogonal polynomials

We would like to choose $\{R_k(x)\}$ and $\{r_k\}$ (which are so far arbitrary) in such a way that the 2×2 matrix $K(x, y) = \begin{bmatrix} S(x, y) & J(x, y) \\ D(x, y) & S(y, x) \end{bmatrix}$ representing the quaternion kernel $\mathbf{K}(x, y)$ enjoys the quasi-projectivity

$$\int_0^\infty dy K(x, y) K(y, z) = K(x, z) \begin{bmatrix} 1 & 0 \\ 0 & 0 \end{bmatrix} + \begin{bmatrix} 0 & 0 \\ 0 & 1 \end{bmatrix} K(x, z), \quad (23)$$

and is correctly normalized:

$$\int_0^\infty dx K(x, x) = N. \quad (24)$$

These two relationships would yield a crucial property that the restricted quaternion matrix $\mathbf{K}_n = [\mathbf{K}(x_i, x_j)]_{i,j=1,\dots,n}$ satisfies the recursion relation

$$\int_0^\infty dx_n \text{qdet} \mathbf{K}_n(x_1, \dots, x_n) = (N - n + 1) \text{qdet} \mathbf{K}_{n-1}(x_1, \dots, x_{n-1}), \quad (25)$$

from which a k -level correlation function is expressed as $\text{qdet} \mathbf{K}_k = \text{Pf}(ZK_k)$. It is well established that the properties (23), (24) are fulfilled by requiring $\{R_k(x)\}$ to be skew-orthogonal with respect to the skew inner product $\langle \cdot, \cdot \rangle$ and $\{r_k\}$ to be their skew-norms,

$$\langle f, g \rangle := \int_0^\infty dx \int_0^\infty dy \sqrt{w(x)w(y)} F(x, y) f(x) g(y) = -\langle g, f \rangle, \quad (26)$$

$$\langle R_{2k}, R_{2k+1} \rangle = -\langle R_{2k+1}, R_{2k} \rangle = r_k, \quad \text{all others} = 0. \quad (27)$$

Under this choice, $F(x, x')$ itself is expressed as

$$F(x, x') = \sum_{k=0}^{\infty} (\phi_{2k}(x) \phi_{2k+1}(x') - \phi_{2k+1}(x) \phi_{2k}(x')). \quad (28)$$

Thus the matrix element of $J_N = [J(x_i, x_j)]_{i,j=1,\dots,N} = [I(x_i, x_j) + F(x_i, x_j)]_{i,j=1,\dots,N}$ takes the form

$$J(x, x') = \sum_{k=N/2}^{\infty} \left(\phi_{2k}(x) \phi_{2k+1}(x') - \phi_{2k+1}(x) \phi_{2k}(x') \right), \quad (29)$$

due to the definition (17).

One can verify that

$$R_{2k}^{(0)}(x) = \sum_{j=0}^k \frac{2^{2k} k! \Gamma(k + (\nu + 1)/2)}{2^{2j} j! \Gamma(j + (\nu + 1)/2)!} (2j)! L_{2j}^{\nu-1}(x) \quad \text{and} \quad R_{2k+1}^{(0)}(x) = -(2k + 1)! L_{2k+1}^{\nu-1}(x) \quad (30)$$

satisfy the skew-orthogonality (27) at $\tau = 0$ (i.e., chGSE, with the skew inner product denoted by $\langle \cdot, \cdot \rangle_{(0)}$) with $r_k^{(0)} = (2k + 1)!(2k + \nu)!$ [25]. Due to a lemma [26],

$$\langle R_m, R_n \rangle = e^{-(\gamma_m + \gamma_n)\tau} \langle R_m^{(0)}, R_n^{(0)} \rangle_{(0)}, \quad (31)$$

which directly follows from the definition (26),

$$R_{2k}(x) = \sum_{j=0}^k \frac{2^{2k} k! \Gamma(k + (\nu + 1)/2)}{2^{2j} j! \Gamma(j + (\nu + 1)/2)!} (2j)! L_{2j}^{\nu-1}(x) e^{(\gamma_{2j} - \gamma_{2k})\tau} \quad \text{and} \quad R_{2k+1}(x) = R_{2k+1}^{(0)}(x) \quad (32)$$

satisfy the skew-orthogonality (27) at $\tau > 0$ with $r_k = r_k^{(0)} e^{-(\gamma_{2k} + \gamma_{2k+1})\tau}$ [14]. Together with the definitions (13)–(16) and (29), the quaternion kernel elements at finite N are completely determined.

2.5. Microscopic crossover scaling limit

Now we concentrate on the case in which the Kramers degeneracy is weakly broken by the small parameter $\tau \ll 1$. Then the spectral density $\rho(\lambda)$ of H in the large- N limit is identical to that of chGSE ($\tau = 0$), i.e., Wigner's semicircle $\rho(\lambda) = \pi^{-1} \sqrt{4N - \lambda^2}$. We magnify the vicinity of the origin by introducing the rescaled variables² $x_i := \lambda_i / \Delta$, which measure the eigenvalues in units of the mean level spacing at the origin, $\Delta = 1/\rho(0) = \pi/\sqrt{4N}$. Moreover, in order to realize a nontrivial crossover behavior we take the triple-scaling limit $N \rightarrow \infty, \lambda \rightarrow 0, \tau \rightarrow 0$ while keeping the combination $\rho = \sqrt{\tau}/\Delta$ and x_i fixed finite. In this limit, sums over k turn

² Here we have abused the notation slightly: Up to Sect. 2.4, $x_i = \lambda_i^2$ denotes squared eigenvalues of H , whereas after Sect. 2.5 $x_i = \lambda_i / \Delta$ denote microscopically rescaled eigenvalues of H . Accordingly the function symbols $K(x, y)$, $S(x, y)$, etc. are also used with two different meanings.

to integrals over $v := k/N$ and Laguerre polynomials reduce to Bessel functions, $\sqrt{w(z)}L_k^v(z) \sim k^{v/2}J_v(2\sqrt{kz})$ as $k \rightarrow \infty$. Accordingly, the quaternion kernel elements (15), (16), (29) reduce to (see footnote 2) [14]

$$S(x, y) = \pi \sqrt{xy} \left\{ \frac{J_v(\pi x) y J_{v-1}(\pi y) - x J_{v-1}(\pi x) J_v(\pi y)}{x^2 - y^2} - \frac{J_v(\pi x)}{2} \pi \int_0^1 dv e^{\pi^2 \rho^2 (v^2 - 1)} J_v(\pi v y) \right\}, \quad (33)$$

$$D(x, y) = \frac{\pi^2 \sqrt{xy}}{2} \int_0^1 dv v \int_0^1 du e^{\pi^2 \rho^2 v^2 (1+u^2)} \{J_v(\pi v u x) J_v(\pi v y) - J_v(\pi v x) J_v(\pi v u y)\}, \quad (34)$$

$$J(x, y) = \frac{\pi^3 \sqrt{xy}}{2} \int_1^\infty dv v^2 e^{-2\pi^2 \rho^2 v^2} \{J_v(\pi v x) y J_{v-1}(\pi v y) - x J_{v-1}(\pi v x) J_v(\pi v y)\}. \quad (35)$$

Thus the correlation function of n positive rescaled eigenvalues $\{x_i\}$ of H in the vicinity of the origin is finally expressed as

$$R_n(x_1, \dots, x_n) = \text{Pf} \left(Z [K(x_i, x_j)]_{i,j=1}^n \right), \quad K(x, y) = \begin{bmatrix} S(x, y) & J(x, y) \\ D(x, y) & S(y, x) \end{bmatrix}. \quad (36)$$

2.6. Individual eigenvalue distributions

It is well known that, for a determinant process in which n -point correlation functions are expressed in terms of a scalar kernel $K(x, y)$ as $R_n(x_1, \dots, x_n) = \det [K(x_i, x_j)]_{i,j=1}^n$, the probability $E_\ell(I)$ for an interval I to contain exactly ℓ points is expressed as a Fredholm determinant (Det) of K over I [23]:

$$E_\ell(I) = \text{Prob}[\#(I) = \ell] = \frac{1}{\ell!} \left(-\frac{\partial}{\partial \xi} \right)^\ell \text{Det}(\mathbb{I} - \xi \hat{K}_I) \Big|_{\xi=1}. \quad (37)$$

Here $\hat{K}_I$ acts on L^2 -functions $f(x)$ over I as $(\hat{K}_I f)(x) = \int_I dy K(x, y) f(y)$. This argument directly carries over to our case of the quaternion determinant process in which n -point correlation functions are expressed in terms of a quaternion kernel $\mathbf{K}(x, y)$ as $R_n(x_1, \dots, x_n) = \text{qdet} \mathbf{K}_n(x_1, \dots, x_n) = \text{Pf} \left(Z [K(x_i, x_j)]_{i,j=1}^n \right)$, leading to

$$E_\ell(I) = \frac{1}{\ell!} \left(-\frac{\partial}{\partial \xi} \right)^\ell \text{Pf}(Z - \xi Z \hat{K}_I) \Big|_{\xi=1} = \frac{1}{\ell!} \left(-\frac{\partial}{\partial \xi} \right)^\ell \text{Det}(\mathbb{I} - \xi \hat{K}_I)^{1/2} \Big|_{\xi=1}. \quad (38)$$

This time, $\hat{K}_I$ acts on 2-component L^2 -functions $F(x)$ over I as $(\hat{K}_I F)(x) = \int_I dy K(x, y) \cdot F(y)$. By differentiating $\text{Det}(\mathbb{I} - \xi \hat{K}_I)^{1/2} = \exp \frac{1}{2} \text{Tr} \log(\mathbb{I} - \xi \hat{K}_I)$ in ξ , the first few $E_\ell(s)$ are expressed in terms of the Fredholm determinant and the resolvents of the operator $\hat{K}_I$,

$$T_n(I) := \frac{1}{2} \text{Tr}(\hat{K}_I (\mathbb{I} - \hat{K}_I)^{-1})^n,$$

as [19]

$$\begin{aligned}
E_0(I) &= \text{Det}(\mathbb{I} - \hat{K}_I)^{1/2} \\
E_1(I) &= E_0 T_1 \\
E_2(I) &= \frac{E_0}{2!} (T_1^2 - T_2) \\
E_3(I) &= \frac{E_0}{3!} (T_1^3 - 3T_1 T_2 + 2T_3) \\
E_4(I) &= \frac{E_0}{4!} (T_1^4 - 6T_1^2 T_2 + 3T_2^2 + 8T_1 T_3 - 6T_4) \\
E_5(I) &= \frac{E_0}{5!} (T_1^5 - 10T_1^3 T_2 + 20T_1^2 T_3 + 15T_1 T_2^2 - 30T_1 T_4 - 20T_2 T_3 + 24T_5) \\
E_6(I) &= \frac{E_0}{6!} \left\{ T_1^6 - 15T_1^4 T_2 + 40T_1^3 T_3 + 45T_1^2 T_2^2 - 90T_1^2 T_4 - 120T_1 T_2 T_3 - 15T_2^3 \right. \\
&\quad \left. + 144T_1 T_5 + 90T_2 T_4 + 40T_3^2 - 120T_6 \right\} \\
E_7(I) &= \frac{E_0}{7!} \left\{ T_1^7 - 21T_1^5 T_2 + 70T_1^4 T_3 + 105T_1^3 T_2^2 - 210T_1^3 T_4 - 420T_1^2 T_2 T_3 - 105T_1 T_2^3 \right. \\
&\quad \left. + 504T_1^2 T_5 + 630T_1 T_2 T_4 + 280T_1 T_3^2 + 210T_2^2 T_3 - 840T_1 T_6 - 504T_2 T_5 \right. \\
&\quad \left. - 420T_3 T_4 + 720T_7 \right\}.
\end{aligned} \tag{39}$$

After specializing to $I = [0, s)$ and abbreviating $E_\ell(s) := E_\ell([0, s))$, the probability distribution $p_k(s)$ of the k th smallest positive eigenvalue is given in terms of $E_0(s), \dots, E_{k-1}(s)$ as

$$p_k(s) = -\frac{d}{ds} \sum_{\ell=0}^{k-1} E_\ell(s). \tag{40}$$

This relationship follows from a simple observation that, for a joint of two intervals $[0, s + ds) = [0, s) \cup [s, s + ds) := I \cup dI$, the probability that the narrower interval dI contains more than one eigenvalue is of order $O(ds^2)$, so to the order $O(ds^1)$ one has

$$E_\ell(s + ds) \simeq \text{Prob}[\#(I) = \ell \cap \#(dI) = 0] + \text{Prob}[\#(I) = \ell - 1 \cap \#(dI) = 1]. \tag{41}$$

Subtracting Eq. (41) from the definition of $E_\ell(s)$ gives

$$E_\ell(s) - E_\ell(s + ds) \simeq \text{Prob}[\#(I) = \ell \cap \#(dI) = 1] - \text{Prob}[\#(I) = \ell - 1 \cap \#(dI) = 1], \tag{42}$$

which is equivalent, in the limit $ds \searrow 0$, to

$$-\frac{d}{ds} E_\ell(s) = p_{\ell+1}(s) - p_\ell(s), \tag{43}$$

with $p_0(s) = 0$ understood. Summing over $\ell = 0, \dots, k-1$ gives Eq. (40).

An efficient way of numerically evaluating the Fredholm determinant of a trace-class operator $\hat{K}_I$ is the Nyström-type discretization [17,18]

$$\text{Det}(\mathbb{I} - \hat{K}_I) \simeq \det(\mathbb{I}_M - \mathcal{K}_I), \text{ where } \mathcal{K}_I = [K(x_i, x_j) \sqrt{w_i w_j}]_{i,j=1}^M \tag{44}$$

is an $M \times M$ matrix evaluated with a quadrature rule consisting of a set of M points $\{x_i\} \in I$ and associated weights $\{w_i\}$ such that $\int_I f(x) dx \simeq \sum_{i=1}^M f(x_i) w_i$. As the order M of the quadrature

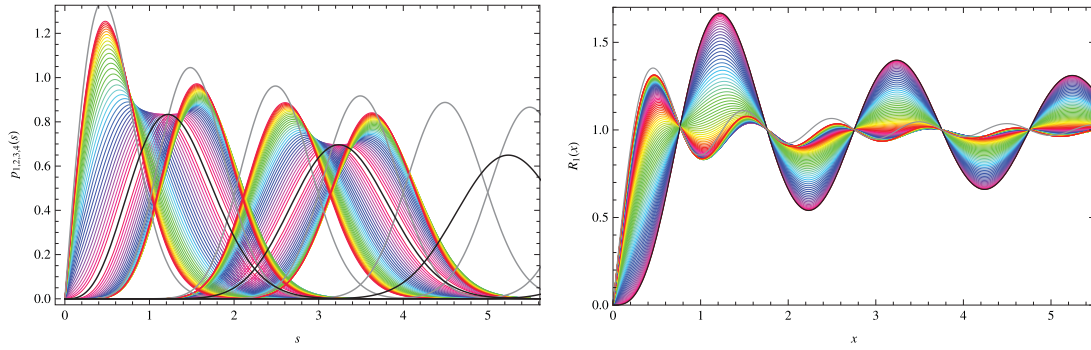


Fig. 1. First four eigenvalue distributions $p_1(s), \dots, p_4(s)$ (left) for $0.04 \leq \rho \leq 0.70$ (step 0.01, purple to red) and the microscopic spectral density $R_1(x)$ (right) for $0.01 \leq \rho \leq 1.00$ (step 0.01, purple to red) for the chGSE (black) to chGUE (gray) crossover, at $\nu = 0$.

increases, the RHS of Eq. (44) is proven to converge to its LHS uniformly and exponentially fast in M [17,18]. We also need to evaluate the resolvents in Eq. (39), which are likewise approximated as

$$\text{Tr}(\hat{K}_I(\mathbb{I} - \hat{K}_I)^{-1})^n \simeq \text{tr}(\mathcal{K}_I(\mathbb{I}_M - \mathcal{K}_I)^{-1})^n. \quad (45)$$

Obviously these formulae hold for a 2×2 -matrix-valued kernel as well. For our purpose we employ the Gauss–Legendre quadrature rule in which $\{x_1, \dots, x_M\}$ are the nodes of the Legendre polynomial $P_M(x)$ on a shifted and rescaled domain $[-1, 1] \mapsto I = [0, s]$. We have applied the Nyström-type method to the kernel (33)–(36) for the chGSE–chGUE crossover at $\nu = 0$, bearing in mind that the topological charge ν is washed away in the staggered Dirac operator that we shall employ in Sect. 3. We have numerically evaluated $p_1(s), \dots, p_4(s)$ with M at least 20 for far-more-than-sufficient precision, and confirmed the stability of the results for increasing M . Plots of $p_1(s), \dots, p_4(s)$ for $0 \leq s \leq 5.5$ are exhibited in Fig. 1 (left).

For comparison, the spectral densities that comprise the former, $R_1(x) = \sum_{k=1}^{\infty} p_k(x) = S(x, x)$ in Eq. (33), are plotted in Fig. 1 (right). The practical advantage of adopting individual eigenvalue distributions over n -level correlation functions (including $R_1(x)$) for fitting is clear from the figures: As the oscillation of the latter consists of overlapping multiple peaks, the characteristic shape of each peak is inevitably smeared, resulting in a rather structureless curve for which an accurate fit is difficult. On the other hand, the shape of the former is clearly distinguishable and is extremely sensitive to the ρ parameter, because the ratio of the two $p_k(s)$ of the chGUE–chGSE crossover at different ρ grows as $\exp(\text{const.}s^2)$ for large s . Therefore, the $p_k(s)$ are expected to admit very precise one-parameter fitting of data by the least-squares method (as will be shown in Tables 3–6 of Sect. 3).

2.7. Effective theory and low-energy constants

Now we shall relate the crossover random matrix ensemble to the nonlinear σ models originating from gauge theory. In continuum, QCD-like theories with N_F flavors of quarks in a real representation have pseudoreal Dirac operators, as does our case of SU(2) lattice gauge theory with fundamental staggered fermions (Eq. (52) in Sect. 3). The low-energy effective Lagrangian of these theories is universally determined by the spontaneous breaking of the Pauli–Gürsey extended flavor group SU($2N_F$) down to its vector subgroup SO(N_F) and takes the form

$$\mathcal{L}_{\text{eff}}(Q) = \frac{1}{2}F^2 \text{tr} \partial_\nu Q^\dagger \partial_\nu Q - \frac{1}{2} \Sigma m \text{Re tr} \hat{M} Q \quad (46)$$

in the leading order of the p -expansion. Here $Q(x)$ is a symmetric $SU(2N_F)$ matrix-valued Nambu–Goldstone field (called a nonlinear σ model of class AI), m the degenerate quark mass, and $\hat{M} = \sigma_1 \otimes \mathbb{I}_{N_F}$. It contains two phenomenological constants: F the pseudo-scalar decay constant and $\Sigma = \langle \bar{\psi}\psi \rangle / N_F$ the chiral condensate (both measured in the chiral and zero-chemical potential limit $m, \mu \rightarrow 0$). The effect of introducing the quark number chemical potential μ to the fundamental theory is unambiguously incorporated in $\mathcal{L}_{\text{eff}}$ through flavor covariantization in the symmetric rank-2 tensor representation [27],

$$\partial_\nu Q \mapsto \nabla_\nu Q = \partial_\nu Q - i\mu\delta_{\nu 0} (\hat{B}Q + Q\hat{B}), \quad (47)$$

with $\hat{B} = \sigma_3 \otimes \mathbb{I}_{N_F}$. If the theory is in a finite volume $V = L^4$ and the Thouless energy defined as $E_c \sim F^2/(\Sigma L^2)$ is much larger than m (called the ε -regime), the path integral is dominated by the zero mode only and takes the form

$$Z = \int dQ \exp \left(V\mu^2 F^2 \text{tr} (\hat{B}Q^\dagger \hat{B}Q + \hat{B}\hat{B}) + \frac{1}{2} V\Sigma m \text{Re tr } \hat{M}Q \right). \quad (48)$$

The above *form* (not the concrete symmetric space on which Q takes values) of action is in fact common for all classes of nonlinear σ models (in even or odd dimensions, with Dyson indices $\beta = 1, 2, 4$) in which the global symmetry is broken by the chemical potential [28].

On the other hand, the characteristic polynomial $\langle \det(\lambda - H)^{N_F} \rangle$ of a random matrix H (3) in the microscopic crossover scaling limit explained in Sect. 2.5 can be evaluated by exponentiating the determinant with N_F flavors of Grassmannian vectors and using a standard technique of Hubbard–Stratonovich transformations (see, e.g., Ref. [29]) for matrices A and B . One can easily show that the outcome is the same nonlinear σ model as Eq. (48) with parameters replaced by

$$VF^2\mu^2 \rightarrow \frac{\pi^2}{2}\rho^2, \quad V\Sigma m \rightarrow i\pi x, \quad (49)$$

respectively. Note that the above identification can also be read off from the exponents in the quaternion kernel elements (33)–(35).

3. Dirac spectrum

In this section we shall fit the probability distributions analytically derived in the previous section to the Dirac operator spectra measured from two types of lattice gauge simulations: (a) $SU(2)$ gauge theory with the imaginary chemical potential, and (b) $SU(2) \times U(1)$ gauge theory. In either case the pseudoreality of the staggered $SU(2)$ Dirac operator is weakly violated by the $U(1)$ field, which is applied as a fixed background [30] or dynamically fluctuating.

3.1. Simulation details

(a) $SU(2)$ gauge theory with the imaginary chemical potential (ICP): The $SU(2)$ variables on temporal links of a hypercubic lattice of size $V = L^4$ are multiplied by a constant phase:

$$\tilde{U}_\nu(x) = U_\nu(x) \times \begin{cases} e^{i2\pi\varphi} & (\nu = 4, x_4 = L - 1) \\ 1 & (\text{else}). \end{cases} \quad (50)$$

The phase $2\pi\varphi$ can be regarded as the Aharonov–Bohm (AB) flux [4] penetrating the temporal circle and is gauge-equivalent to the imaginary chemical potential $\mu = i\mu_I = 2\pi i\varphi/L$, i.e., a fixed

U(1) background $B_\nu = (2\pi\varphi/L)\delta_{\nu,4}$. We chose antiperiodic/periodic boundary conditions in the temporal/spatial directions, respectively, and consider a small twisting along the temporal direction ($\varphi \ll 1$).

(b) $SU(2) \times U(1)$ gauge theory: Following Ref. [31], non-compact U(1) link variables $B_\nu(x)$ are generated under the Coulomb gauge-fixing condition (with an additional constraint for the $B_4(x)$) and at the unit coupling constant, and are multiplied to SU(2) link variables $U_\nu(x)$,

$$\tilde{U}_\nu(x) = U_\nu(x) e^{ieB_\nu(x)}, \quad (51)$$

with e denoting the bare U(1) coupling constant. Fermions are quenched both for the SU(2) and U(1) gauge fields. For the pure SU(2) case $e = 0$ (or $\varphi = 0$), the staggered SU(2) Dirac operator in the fundamental representation

$$D_{x,y} = \sum_{\nu=1}^4 (-1)^{\sum_{i=1}^{\nu-1} x_i} \left(\tilde{U}_\nu(x) \delta_{x,y+\hat{\nu}} - \tilde{U}_\nu^\dagger(x) \delta_{x,y-\hat{\nu}} \right) \quad (52)$$

is pseudoreal, i.e., satisfies

$$\mathcal{T} D \mathcal{T}^{-1} = D \quad \text{with} \quad \mathcal{T}^2 = (\mathcal{Z}\mathcal{C})^2 = -\mathbb{I} \quad (53)$$

($\mathcal{C}$ denotes complex conjugation) for either choice of periodic or antiperiodic boundary condition in each direction, and we impose periodic boundary conditions in all four directions and consider the U(1) part as a small perturbation (i.e., $e \ll 1$).

As the presence of the AB flux φ or the U(1) coupling e breaks the pseudoreality (53), they parametrize antiunitary symmetry breaking in Dirac operators and are anticipated to be the lattice gauge theory counterpart of the crossover parameter ρ in random matrix ensemble interpolating chGSE and chGUE. We note that while the effect of the ICP μ_I on the low-energy effective Lagrangian is completely dictated on the symmetry ground [32,33] and is related to the pseudo-scalar decay constant F as for the real chemical potential [27], the effect of e cannot be directly related to F , because integrating over the dynamical U(1) gauge field in the effective Lagrangian would lead to nonlocal self-couplings of pseudo-scalar mesons.

3.1.1. Simulation setup

We measured low-lying spectra of the naïve staggered Dirac operator (52) on small lattices of volume $V = 4^4$ and 6^4 . In order to examine the validity of our method for the strong-coupling to the near-continuum scaling regions, we chose the bare SU(2) gauge coupling constant $\beta = 4/g^2$ from the range $\beta = 0, 0.25, \dots, 1.75$ (step 0.25) on $V = 4^4$ and $\beta = 0, 0.25, \dots, 2.0$ (step 0.25), 2.1 on $V = 6^4$. The simplest algorithm is employed in generating SU(2) gauge configurations: unimproved plaquette action and the 10-hit heat-bath update combined with over-relaxation. The antiunitary symmetry violation parameters are set to be: (a) $\varphi = 0.01, \dots, 0.06$ (step 0.01) on $V = 4^4$ and $\varphi = 0.01, \dots, 0.05$ (step 0.005) on $V = 6^4$, and (b) $e = 0.002, \dots, 0.006$ (step 0.001), 0.008, 0.0010 on $V = 4^4$ and $e = 0.0004, \dots, 0.0016$ (step 0.0002), 0.0020, 0.0024, 0.0028 on $V = 6^4$. $N_{\text{conf}} = 40\,000$ (10\,000) configurations are generated and diagonalized on $V = 4^4$ (6^4) for each set of parameters (β, φ) or (β, e) .

3.1.2. Fitting Dirac spectra

Our procedure of fitting the Dirac spectra to the individual eigenvalue distributions of the crossover chiral random matrices consists of the following two steps:

- (i) *Determination of the mean level spacing Δ at the origin.* In order to determine the physical scale of Dirac eigenvalues upon which the effects of perturbations are to be evaluated, we measure for N_{conf} independent configurations four low-lying nondegenerate eigenvalues $\lambda_{2i-1} = \lambda_{2i}$ ($i = 1, \dots, 4$) of the pure SU(2) Dirac operator (due to its quaternionic nature, all eigenvalues are doubly degenerate).

For each i , the mean level spacing Δ_i at the spectral origin is determined by best-fitting the histogram of the unfolded Dirac eigenvalue λ_{2i}/Δ_i to the normalized individual eigenvalue distribution $p_i(s)$ of chGSE so that $\chi^2/\text{d.o.f.}$ is minimized by varying Δ_i . In doing so, we discard the left tail $[0, s_{\min}]$ and the right tail $[s_{\max}, \infty)$ of the probability distribution $p_i(s)$ for which $\int_0^{s_{\min}} ds p_i(s) = \int_{s_{\max}}^{\infty} ds p_i(s) \simeq 10^{-3}$, and split the mid-range into B bins of fixed widths $\delta s = 0.1$, $[s_{\min}, s_{\max}] = I_1 \cup \dots \cup I_B$. Then we define χ^2 from the measured frequency $F_b = \#\{\lambda_{2i} \in I_b\}$ and its analytic prediction $f_b = N_{\text{conf}} \int_{I_b} ds p_i(s)$ by $\chi^2 = \sum_{b=1}^B (F_b - f_b)^2 / f_b$. The statistical error $\delta\Delta_i$ is estimated as a deviation from the optimal Δ_i at which $\chi^2/\text{d.o.f.}$ increases by unity. The combined value of the mean level spacing at the origin $\bar{\Delta}$ is obtained as the weighted average of $(\Delta_i, \delta\Delta_i)$, $i = 1, \dots, 4$. We have confirmed that these four data are always mutually consistent, so that their combination helps to improve the statistical error in $\bar{\Delta}$ as compared to the previous method of using the smallest eigenvalue only [10,11].

- (ii) *Determination of the crossover parameter ρ .* Next we switch on the U(1) perturbation (a) or (b) and measure the Dirac spectra $\{\lambda_k\}$ for $N_{\text{conf}} = \mathcal{O}(10^4)$ independent configurations. The effect of such perturbations on Δ (i.e., the chiral condensate) is negligible in the lowest order in the ε -expansion that we are working on. U(1) perturbations split once-Kramers-degenerate pairs of eigenvalues, $\lambda_{2i-1} < \lambda_{2i}$. We take the first two pairs of Dirac eigenvalues $(\lambda_{2i-1}, \lambda_{2i})$, $i = 1, 2$, and define the unfolded eigenvalues as $(s_{2i-1}, s_{2i}) = (\lambda_{2i-1}/\Delta_i, \lambda_{2i}/\Delta_i)$. Then by using the same strategy as in Step (i), their histograms $P_k(s_k)$ are fitted to the analytic results $p_k(s)$ (39), (40) of the crossover random matrices, with the crossover parameter ρ_k being varied. The statistical error $\delta\rho_k$ of the crossover parameter is again estimated as a deviation from the optimal ρ_k at which $\chi^2/\text{d.o.f.}$ increases by unity.

The crossover parameter $\bar{\rho}$ corresponding to a particular choice of φ or e is eventually determined as the weighted average of $(\rho_k, \delta\rho_k)$, $k = 1, \dots, 4$. The advantage of using both once-degenerate eigenvalue pairs (over using, e.g., three low-lying eigenvalues) is now evident: Histograms of s_{2i-1} and s_{2i} always shift in the opposite directions under the Kramers-breaking perturbation, so that the effect of small errors in the determination of the overall scale Δ_i in Step (i) (which shifts both histograms in the same direction) is expected to be canceled in the final value of $\bar{\rho}$ obtained by combining ρ_{2i-1} and ρ_{2i} .

3.1.3. Low-energy constants

Due to the correspondence (49), two low-energy constants contained in the chiral Lagrangian, the chiral condensate Σ and the pseudo-scalar decay constant F , are directly related to Δ and ρ measured in Steps (i) and (ii), the former by the Banks–Casher relation

$$\Sigma = \frac{\pi}{\Delta V}; \quad (54)$$

the latter for (a) the SU(2)+ICP model by

$$F^2 = \frac{\pi^2}{2V} \left(\frac{\rho}{\mu_1} \right)^2 = \frac{\pi^2}{2V} \left(\frac{\rho}{2\pi\varphi/L} \right)^2, \quad (55)$$

and for (b) the SU(2)×U(1) model by

$$\frac{F^2 \mu_1^2}{e^2} = \frac{\pi^2}{2V} \left(\frac{\rho}{e} \right)^2. \quad (56)$$

Note that the combination “ $F^2 \mu_1^2$ ” in the LHS of Eq. (56) is to be regarded as a single coefficient of the $-\text{tr} \hat{B} Q^\dagger \hat{B} Q$ term in the chiral Lagrangian (48). As this quantity should be proportional to e^2 , our aim here is to determine the unknown proportionality constant set by the dynamics. Thus we shall check, for each β in either case of (a) or (b), the stability of the ratio $R(\varphi) = \rho/\varphi$ or $R(e) = \rho/e$ as φ or e is varied, and determine its mean value as a weighted average of $\{R(\varphi), \delta R(\varphi)\}$ or $\{R(e), \delta R(e)\}$.

3.2. Simulation results

3.2.1. Fitting Dirac spectra

In Tables 1 and 2 we exhibit optimal values of the mean level spacings Δ_i of SU(2) Dirac spectra (a) under the antiperiodic boundary condition on the temporal direction (to be used for SU(2)+ICP), and (b) under the periodic boundary conditions on all directions (to be used for SU(2) × U(1)), respectively. At each β , we adopt only the Δ_i that pass the χ^2 test of the fitting: $\chi^2/\text{d.o.f.} < 2.0$. In Fig. 2 (top) we exhibit sample plots of histograms of the four smallest nondegenerate Dirac

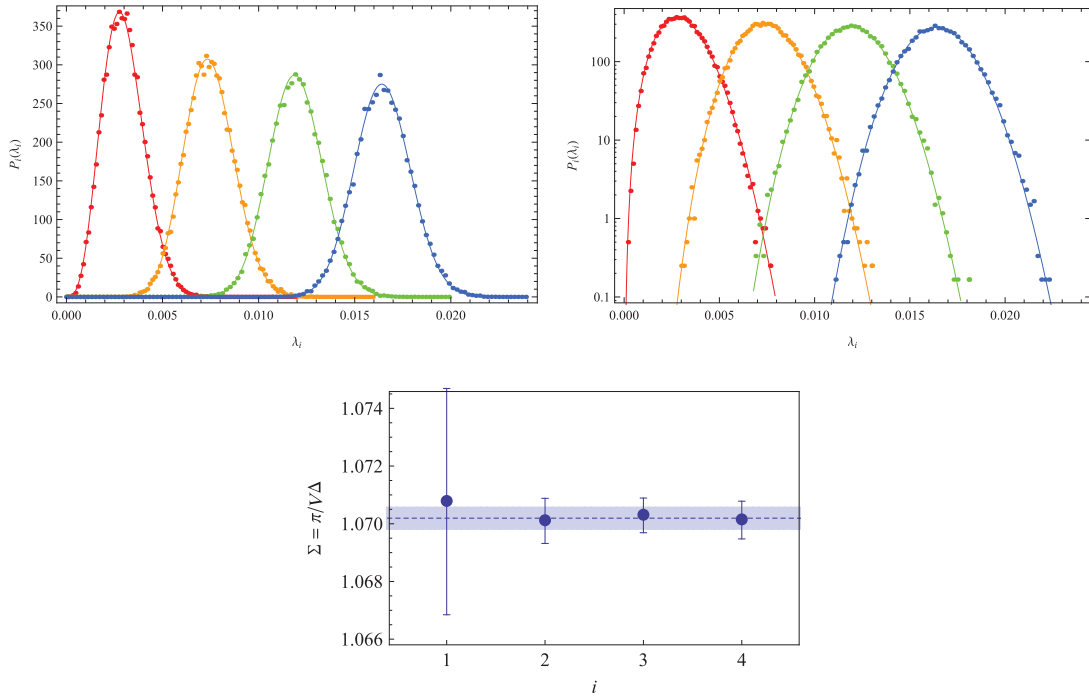


Fig. 2. (Top) Linear and logarithmic plots of histograms of the four smallest nondegenerate Dirac eigenvalues $P_i(\lambda_i)$ ($i = 1, \dots, 4$) (red to blue) of pure SU(2) gauge theory at $\beta = 1.0$, on $V = 6^4$ and $N_{\text{conf}} = 30\,000$, and individual eigenvalue distributions $p_i(s)$ of chGSE, each being optimally rescaled by a constant Δ_i . (Bottom) Values of chiral condensate $\Sigma = \pi/(V\Delta_i)$ (circles), their weighted average (horizontal line), and the combined error (strip).

Table 1. Mean level spacings of the pure SU(2) Dirac spectrum on $V = 4^4$ and 6^4 with the antiperiodic boundary condition on the temporal direction, in units of $10^{-2}a^{-1}$.

V	β	Δ_1	Δ_2	Δ_3	Δ_4	$\bar{\Delta}$	$\chi^2/\text{d.o.f.}$
4^4	0	0.929(2)	0.931 5(7)	0.930 9(5)	0.931 3(4)	0.931 1(3)	0.99–1.00
	0.25	0.969(2)	0.970 8(9)	0.970 9(6)	0.970 3(4)	0.970 5(3)	0.85–1.00
	0.5	1.019(2)	1.019 0(9)	1.018 0(6)	1.017 5(4)	1.017 9(3)	0.74–1.02
	0.75	1.073(2)	1.073 5(8)	1.074 9(6)	1.075 6(5)	1.075 0(3)	1.01–1.36
	1.0	1.151(2)	1.150 0(9)	1.150 4(7)	–	1.150 3(5)	0.68–1.47
	1.25	1.255(2)	1.254(1)	–	–	1.254 4(9)	0.81–1.46
	1.5	1.408(3)	–	–	–	1.408(3)	0.73
	1.75*	1.705(4)	–	–	–	1.705(4)	0.91
6^4	0	0.185 4(7)	0.185 2(3)	0.185 0(2)	0.185 2(2)	0.185 2(1)	0.92–1.02
	0.25	0.191 2(7)	0.192 6(3)	0.193 1(2)	0.192 7(2)	0.192 8(1)	0.98–1.20
	0.5	0.203 0(8)	0.202 7(4)	0.202 0(2)	0.202 3(2)	0.202 3(1)	0.95–0.99
	0.75	0.213 5(7)	0.213 3(2)	0.214 3(2)	0.213 5(2)	0.213 7(1)	0.85–1.01
	1.0	0.229 3(8)	0.227 7(4)	0.227 9(3)	0.227 8(2)	0.227 8(1)	0.93–1.00
	1.25	0.247 5(9)	0.248 0(4)	0.247 9(3)	0.248 1(2)	0.248 0(2)	0.63–0.99
	1.5	0.279 5(9)	0.278 7(5)	0.278 3(3)	0.277 9(2)	0.278 2(2)	0.96–1.00
	1.75	0.334(1)	0.333 1(6)	0.332 9(4)	–	0.333 0(3)	0.73–1.00
	2.0	0.482(2)	–	–	–	0.482(2)	1.12
	2.1*	0.640(2)	–	–	–	0.640(2)	1.49

Table 2. Mean level spacings of the pure SU(2) Dirac spectrum on $V = 4^4$ and 6^4 with periodic boundary conditions on all four directions, in units of $10^{-2}a^{-1}$.

V	β	Δ_1	Δ_2	Δ_3	Δ_4	$\bar{\Delta}$	$\chi^2/\text{d.o.f.}$
4^4	0	0.932(2)	0.930 6(8)	0.931 1(5)	0.931 3(4)	0.931 1(3)	0.62–1.13
	0.25	0.972(2)	0.971 8(8)	0.970 8(6)	0.970 4(4)	0.970 8(3)	0.84–1.13
	0.5	1.020(2)	1.016 0(9)	1.016 2(6)	1.017 0(4)	1.016 7(3)	0.97–1.20
	0.75	1.080(2)	1.077 0(9)	1.077 4(6)	1.076 5(5)	1.076 9(3)	0.76–1.15
	1.0	1.153(2)	1.152(1)	1.152 5(7)	–	1.152 4(5)	1.00–1.14
	1.25	1.249(2)	1.253(1)	–	–	1.252 5(9)	0.99–1.00
	1.5	1.412(2)	–	–	–	1.412(2)	0.93
	1.75*	1.698(3)	–	–	–	1.698(3)	1.35
6^4	0	0.186 5(6)	0.185 4(3)	0.185 3(2)	0.185 4(2)	0.185 4(1)	0.79–1.00
	0.25	0.192 8(7)	0.192 9(3)	0.192 8(2)	0.192 8(2)	0.192 8(1)	0.99–1.00
	0.5	0.202 4(8)	0.202 4(4)	0.202 0(2)	0.202 0(2)	0.202 1(1)	0.77–1.01
	0.75	0.214 6(8)	0.213 7(4)	0.213 8(2)	0.213 6(2)	0.213 7(1)	0.98–1.01
	1.0	0.228 6(8)	0.229 2(4)	0.228 5(3)	0.228 5(2)	0.228 6(1)	0.99–1.00
	1.25	0.248 1(9)	0.247 8(6)	0.248 4(3)	0.247 7(2)	0.248 0(2)	0.97–1.01
	1.5	0.277(1)	0.277 8(5)	0.278 3(3)	0.278 2(2)	0.278 1(2)	0.99–1.00
	1.75	0.333(1)	0.331 8(6)	0.333 1(4)	–	0.332 7(3)	0.96–1.24
	2.0	0.480(2)	–	–	–	0.480(2)	1.66
	2.1*	0.639(3)	–	–	–	0.639(3)	0.99

eigenvalues $P_i(\lambda_i)$ versus the corresponding individual eigenvalue distributions $p_i(s)$ of chGSE, each being optimally rescaled by Δ_i . The mutual consistency of such Δ_i observed in Fig. 2 (bottom) justifies the use of the weighted average, listed in the seventh columns in Tables 1 and 2. This sample figure illustrates that the error in Δ_1 (i.e., in Σ), which is relatively larger than that in the other three Δ_i , is considerably improved by a factor of 5–10 and is down to $\mathcal{O}(10^{-4})$ by the use of the

Table 3. Crossover parameters ρ for the SU(2)+ICP model on $V = 4^4$.

β	ρ	φ						$\chi^2/\text{d.o.f.}$
		0.01	0.02	0.03	0.04	0.05	0.06	
0	ρ_1	0.061(1)	0.122(1)	0.183(1)	0.244(2)	—	—	0.68–1.06
	ρ_2	0.061(2)	0.121(2)	0.183(2)	0.238(3)	—	—	0.95–1.71
	ρ_3	0.060(2)	0.118(2)	0.181(2)	0.243(2)	—	—	0.50–0.69
	ρ_4	0.062(2)	0.125(2)	0.184(2)	0.245(3)	—	—	0.64–1.19
	$\bar{\rho}$	0.061 2(8)	0.121 4(9)	0.183(1)	0.243(1)	—	—	—
0.25	ρ_1	0.060(1)	0.117(1)	0.176(1)	0.234(2)	0.291(2)	—	0.65–1.21
	ρ_2	0.057(2)	0.118(2)	0.174(2)	0.234(3)	0.296(4)	—	0.63–1.22
	ρ_3	0.062(2)	0.117(2)	0.176(2)	0.235(2)	0.291(3)	—	0.80–1.49
	ρ_4	0.056(2)	0.116(2)	0.174(2)	0.231(3)	0.298(4)	—	0.76–1.30
	$\bar{\rho}$	0.058 6(8)	0.116 9(9)	0.175 3(9)	0.234(1)	0.292(1)	—	—
0.5	ρ_1	0.056(1)	0.108(1)	0.168(1)	0.224(2)	0.279(2)	—	0.57–1.56
	ρ_2	0.056(2)	0.117(2)	0.167(2)	0.227(3)	0.282(4)	—	0.35–1.33
	ρ_3	0.056(2)	0.111(2)	0.165(2)	0.228(2)	0.279(2)	—	0.64–1.25
	ρ_4	0.057(2)	0.112(2)	0.171(2)	0.217(3)	0.277(4)	—	0.82–1.34
	$\bar{\rho}$	0.056 1(8)	0.111 3(9)	0.167 7(9)	0.224(1)	0.279(1)	—	—
0.75	ρ_1	0.055(1)	0.108(1)	0.162(1)	0.216(1)	0.270(2)	—	0.63–0.90
	ρ_2	0.053(2)	0.106(2)	0.159(2)	0.213(3)	0.266(4)	—	0.76–1.23
	ρ_3	0.056(2)	0.110(2)	0.164(2)	0.218(2)	0.271(2)	—	0.78–1.39
	ρ_4	0.051(2)	0.104(2)	0.158(2)	0.210(3)	0.262(3)	—	0.98–1.18
	$\bar{\rho}$	0.053 8(8)	0.107 3(9)	0.161 0(9)	0.215(1)	0.269(1)	—	—
1.0	ρ_1	0.052(1)	0.102(1)	0.153(1)	0.203(1)	0.254(2)	—	0.78–1.05
	ρ_2	0.050(2)	0.101(2)	0.151(2)	0.201(2)	0.251(3)	—	0.60–1.16
	ρ_3	0.050(2)	0.101(2)	0.151(2)	0.202(2)	0.252(2)	—	0.78–1.39
	ρ_4	0.051(2)	0.102(2)	0.153(2)	0.204(3)	0.253(3)	—	0.90–1.23
	$\bar{\rho}$	0.050 8(8)	0.101 5(9)	0.152 2(9)	0.203(1)	0.253(1)	—	—
1.25	ρ_1	0.048(1)	0.096(1)	0.143(1)	0.190(1)	0.238(2)	0.285(2)	0.88–1.23
	ρ_2	0.047(2)	0.095(2)	0.142(2)	0.189(2)	0.236(3)	0.283(4)	0.70–1.14
	ρ_3	—	—	—	—	—	—	—
	ρ_4	—	—	—	—	—	—	—
	$\bar{\rho}$	0.048(1)	0.095(1)	0.143(9)	0.190(1)	0.237(1)	0.285(2)	—
1.5	ρ_1	—	0.087(1)	0.130(1)	0.174(1)	0.217(2)	0.261(2)	0.69–0.94
	ρ_2	—	0.086(2)	0.129(2)	0.172(2)	0.214(3)	0.254(3)	0.62–1.03
	ρ_3	—	—	—	—	—	—	—
	ρ_4	—	—	—	—	—	—	—
	$\bar{\rho}$	—	0.087(1)	0.130(1)	0.173(1)	0.216(1)	0.259(1)	—
1.75*	ρ_1	—	0.075(2)	0.113(2)	0.150(2)	0.188(2)	0.226(2)	0.88–1.99
	ρ_2	—	0.074(2)	0.110(2)	0.147(3)	0.180(3)	0.216(4)	0.56–1.30
	ρ_3	—	—	—	—	—	—	—
	ρ_4	—	—	—	—	—	—	—
	$\bar{\rho}$	—	0.075(1)	0.112(2)	0.149(2)	0.186(2)	0.224(2)	—

weighted average of the four. Note that histograms at $\beta = 1.75$ on $V = 4^4$ and at $\beta = 2.1$ on $V = 6^4$ (marked by * in Tables 1 and 2) failed to be fitted into the chGSE predictions in the above criterion as $\chi^2/\text{d.o.f.}$ exceeds 3. Thus we chose to relax it by cutting off the tails of the distributions for which $p_1(s) < 0.2$ from fitting, which leads safely to $\chi^2/\text{d.o.f.} < 2.0$.

Table 4. Crossover parameters ρ for the SU(2)+ICP model on $V = 6^4$.

β	ρ	φ									$\chi^2/\text{d.o.f.}$
		0.01	0.015	0.02	0.025	0.03	0.035	0.04	0.045	0.05	
0	ρ_1	0.093(3)	0.135(3)	0.178(3)	0.228(3)	—	—	—	—	—	0.54–1.18
	ρ_2	0.089(3)	0.138(4)	0.187(5)	0.228(6)	—	—	—	—	—	0.47–1.40
	ρ_3	0.083(4)	0.135(4)	0.173(4)	0.223(4)	—	—	—	—	—	0.56–1.30
	ρ_4	0.099(4)	0.138(4)	0.191(5)	0.236(6)	—	—	—	—	—	0.76–1.09
	$\bar{\rho}$	0.091(2)	0.136(2)	0.180(2)	0.228(2)	—	—	—	—	—	—
0.25	ρ_1	0.088(3)	0.135(3)	0.177(3)	0.223(3)	—	—	—	—	—	0.87–1.32
	ρ_2	0.087(3)	0.130(4)	0.171(4)	0.215(5)	—	—	—	—	—	0.51–1.37
	ρ_3	0.090(4)	0.133(4)	0.179(4)	0.218(4)	—	—	—	—	—	0.66–1.05
	ρ_4	0.087(4)	0.132(4)	0.169(4)	0.229(6)	—	—	—	—	—	0.76–1.19
	$\bar{\rho}$	0.088(2)	0.133(2)	0.175(2)	0.221(2)	—	—	—	—	—	—
0.5	ρ_1	0.083(3)	0.119(3)	0.168(3)	0.217(3)	—	—	—	—	—	0.96–1.25
	ρ_2	0.085(3)	0.135(4)	0.170(4)	0.201(5)	—	—	—	—	—	0.60–0.94
	ρ_3	0.083(4)	0.123(4)	0.166(4)	0.222(4)	—	—	—	—	—	0.77–1.32
	ρ_4	0.086(4)	0.130(4)	0.173(5)	0.200(5)	—	—	—	—	—	0.48–1.23
	$\bar{\rho}$	0.084(2)	0.125(2)	0.169(2)	0.213(2)	—	—	—	—	—	—
0.75	ρ_1	0.082(3)	0.123(3)	0.157(3)	0.202(3)	0.240(3)	—	—	—	—	0.70–1.47
	ρ_2	0.077(3)	0.119(4)	0.171(4)	0.205(5)	0.246(6)	—	—	—	—	0.46–1.12
	ρ_3	0.087(4)	0.124(4)	0.163(4)	0.207(4)	0.241(4)	—	—	—	—	0.60–1.14
	ρ_4	0.076(4)	0.116(4)	0.159(4)	0.200(5)	0.244(6)	—	—	—	—	0.80–1.33
	$\bar{\rho}$	0.081(2)	0.121(2)	0.161(2)	0.203(2)	0.242(2)	—	—	—	—	—
1.0	ρ_1	0.076(3)	0.111(3)	0.153(3)	0.188(3)	0.233(3)	—	—	—	—	0.49–0.97
	ρ_2	0.080(3)	0.117(4)	0.157(4)	0.197(5)	0.228(6)	—	—	—	—	0.71–1.27
	ρ_3	0.078(4)	0.108(4)	0.146(4)	0.186(4)	0.227(4)	—	—	—	—	0.75–1.38
	ρ_4	0.078(4)	0.125(4)	0.165(5)	0.199(5)	0.233(6)	—	—	—	—	0.70–1.13
	$\bar{\rho}$	0.078(2)	0.114(2)	0.154(2)	0.191(2)	0.231(2)	—	—	—	—	—
1.25	ρ_1	0.074(3)	0.107(3)	0.146(3)	0.180(3)	0.213(3)	—	—	—	—	0.54–1.31
	ρ_2	0.073(3)	0.113(4)	0.144(4)	0.180(5)	0.219(5)	—	—	—	—	0.67–1.05
	ρ_3	0.071(4)	0.106(4)	0.152(4)	0.181(4)	0.214(4)	—	—	—	—	0.84–0.98
	ρ_4	0.074(4)	0.111(4)	0.138(4)	0.179(5)	0.222(5)	—	—	—	—	0.66–1.24
	$\bar{\rho}$	0.073(2)	0.109(2)	0.145(2)	0.180(2)	0.216(2)	—	—	—	—	—
1.5	ρ_1	0.068(3)	0.099(3)	0.136(3)	0.169(3)	0.196(3)	0.230(3)	—	—	—	0.61–0.89
	ρ_2	0.067(3)	0.104(4)	0.132(4)	0.158(4)	0.206(5)	0.240(6)	—	—	—	0.60–1.02
	ρ_3	0.069(4)	0.106(4)	0.141(4)	0.172(4)	0.201(4)	0.232(4)	—	—	—	0.80–1.17
	ρ_4	0.063(4)	0.094(4)	0.128(4)	0.157(4)	0.200(5)	0.231(6)	—	—	—	0.84–1.32
	$\bar{\rho}$	0.067(2)	0.100(2)	0.135(2)	0.166(2)	0.199(2)	0.232(2)	—	—	—	—
1.75	ρ_1	0.058(3)	0.090(3)	0.115(3)	0.148(3)	0.172(3)	0.209(3)	—	—	—	0.67–1.21
	ρ_2	0.060(3)	0.084(3)	0.121(4)	0.144(4)	0.181(5)	0.203(5)	—	—	—	0.67–1.60
	ρ_3	0.057(4)	0.091(4)	0.117(4)	0.148(4)	0.178(4)	0.216(4)	—	—	—	0.76–1.54
	ρ_4	0.060(4)	0.083(4)	0.116(4)	0.145(4)	0.172(5)	0.192(5)	—	—	—	0.77–1.68
	$\bar{\rho}$	0.059(2)	0.087(2)	0.117(2)	0.147(2)	0.175(2)	0.207(2)	—	—	—	—
2.0	ρ_1	—	0.072(3)	0.095(3)	0.117(3)	0.140(3)	0.163(3)	0.185(3)	0.207(3)	0.230(3)	0.94–1.48
	ρ_2	—	0.063(3)	0.085(3)	0.108(4)	0.131(4)	0.152(4)	0.173(4)	0.194(5)	0.214(5)	0.90–1.37
	ρ_3	—	—	—	—	—	—	—	—	—	—
	ρ_4	—	—	—	—	—	—	—	—	—	—
	$\bar{\rho}$	—	0.068(2)	0.091(2)	0.114(2)	0.137(2)	0.159(2)	0.181(2)	0.204(3)	0.226(3)	—
2.1*	ρ_1	—	—	0.079(4)	—	0.114(4)	0.133(4)	0.151(4)	0.170(4)	0.189(4)	1.55–1.90
	ρ_2	—	—	0.064(5)	—	0.101(5)	0.119(5)	0.137(5)	0.150(6)	0.167(6)	1.42–1.94
	ρ_3	—	—	—	—	—	—	—	—	—	—
	ρ_4	—	—	—	—	—	—	—	—	—	—
	$\bar{\rho}$	—	—	0.073(3)	—	0.109(3)	0.128(3)	0.147(3)	0.165(3)	0.183(3)	—

Table 5. Crossover parameters ρ for the $SU(2) \times U(1)$ model on $V = 4^4$.

β	ρ	e							$\chi^2/\text{d.o.f.}$
		0.002	0.003	0.004	0.005	0.006	0.008	0.010	
0	ρ_1	0.093(1)	0.136(1)	0.186(1)	0.232(2)	0.276(2)	—	—	0.62–1.52
	ρ_2	0.905(2)	0.142(2)	0.181(2)	0.225(3)	0.277(4)	—	—	0.73–1.38
	ρ_3	0.089(2)	0.136(2)	0.183(2)	0.231(2)	0.273(2)	—	—	0.82–1.20
	ρ_4	0.094(2)	0.139(2)	0.187(2)	0.230(3)	0.280(4)	—	—	0.81–1.08
	$\bar{\rho}$	0.091 8(8)	0.137 8(9)	0.184(1)	0.230(1)	0.276(1)	—	—	—
0.25	ρ_1	0.088(1)	0.136(1)	0.176(1)	0.225(2)	0.265(2)	—	—	0.48–1.14
	ρ_2	0.088(2)	0.128(2)	0.177(2)	0.216(3)	0.264(3)	—	—	0.76–1.75
	ρ_3	0.087(2)	0.137(2)	0.175(2)	0.226(2)	0.262(2)	—	—	0.55–1.04
	ρ_4	0.088(2)	0.128(2)	0.176(2)	0.214(3)	0.261(3)	—	—	0.75–1.53
	$\bar{\rho}$	0.088 0(8)	0.133 1(9)	0.176 0(9)	0.222(1)	0.264(1)	—	—	—
0.5	ρ_1	0.083(1)	0.126(1)	0.167(1)	0.210(2)	0.251(2)	—	—	0.50–0.99
	ρ_2	0.086(2)	0.128(2)	0.170(2)	0.212(3)	0.256(3)	—	—	0.53–0.94
	ρ_3	0.086(2)	0.121(2)	0.170(2)	0.206(2)	0.255(2)	—	—	0.79–1.34
	ρ_4	0.082(2)	0.133(2)	0.166(2)	0.218(3)	0.248(3)	—	—	0.38–1.34
	$\bar{\rho}$	0.084 0(8)	0.126 2(9)	0.168 2(9)	0.210(1)	0.252(1)	—	—	—
0.75	ρ_1	0.078(1)	0.118(1)	0.159(1)	0.199(1)	0.239(2)	—	—	0.68–1.22
	ρ_2	0.082(2)	0.122(2)	0.163(2)	0.202(2)	0.244(3)	—	—	0.48–0.84
	ρ_3	0.081(2)	0.116(2)	0.161(2)	0.196(2)	0.242(2)	—	—	0.72–1.01
	ρ_4	0.080(2)	0.124(2)	0.159(2)	0.204(3)	0.237(3)	—	—	0.62–1.40
	$\bar{\rho}$	0.079 9(8)	0.119 8(9)	0.160 2(9)	0.200(1)	0.240(1)	—	—	—
1.0	ρ_1	0.075(1)	0.114(1)	0.150(1)	0.189(1)	0.225(2)	0.300(2)	—	0.75–1.27
	ρ_2	0.075(2)	0.111(2)	0.150(2)	0.184(2)	0.223(3)	0.296(4)	—	1.00–1.55
	ρ_3	0.076(2)	0.114(2)	0.151(2)	0.189(2)	0.226(2)	0.303(3)	—	0.94–1.62
	ρ_4	0.073(2)	0.111(2)	0.147(2)	0.183(2)	0.218(2)	0.283(4)	—	0.86–1.20
	$\bar{\rho}$	0.074 8(8)	0.112 6(9)	0.149 4(9)	0.187(1)	0.224(1)	0.298(1)	—	—
1.25	ρ_1	0.068(1)	0.103(1)	0.138(1)	0.172(1)	0.208(1)	0.278(2)	—	0.60–1.15
	ρ_2	0.069(2)	0.104(2)	0.138(2)	0.172(2)	0.206(3)	0.272(4)	—	0.53–0.95
	ρ_3	0.069(2)	0.107(2)	0.139(2)	0.177(2)	0.210(2)	0.281(2)	—	1.14–1.57
	ρ_4	0.069(2)	0.097(2)	0.134(2)	0.162(2)	0.201(3)	0.264(3)	—	1.32–1.65
	$\bar{\rho}$	0.068 8(8)	0.103 2(8)	0.137 7(9)	0.171 6(9)	0.207(1)	0.276(1)	—	—
1.5	ρ_1	0.063(1)	0.092(1)	0.125(1)	0.154(1)	0.187(1)	0.250(2)	0.313(2)	0.68–1.06
	ρ_2	0.060(2)	0.093(2)	0.121(2)	0.154(2)	0.180(2)	0.236(3)	0.286(4)	0.61–1.25
	ρ_3	—	—	—	—	—	—	—	—
	ρ_4	—	—	—	—	—	—	—	—
	$\bar{\rho}$	0.062(1)	0.092(1)	0.123(1)	0.154(1)	0.185(1)	0.246(1)	0.308(2)	—
1.75*	ρ_1	0.054(2)	—	0.106(2)	—	0.158(2)	0.210(2)	0.263(2)	1.35–1.63
	ρ_2	0.049(2)	—	0.100(2)	—	0.148(2)	0.195(3)	0.236(3)	1.16–1.77
	ρ_3	—	—	—	—	—	—	—	—
	ρ_4	—	—	—	—	—	—	—	—
	$\bar{\rho}$	0.052(1)	—	0.103(1)	—	0.154(1)	0.206(1)	0.258(2)	—

Next we exhibit the optimal values of the crossover parameters $\rho_k (k = 1, \dots, 4)$, determined by fitting individual Dirac eigenvalue histograms to the chGSE–chGUE predictions, (a) for the

Table 6. Crossover parameters ρ for the $SU(2) \times U(1)$ model on $V = 6^4$.

β	ρ	e										$\chi^2/\text{d.o.f.}$
		0.0004	0.0006	0.0008	0.0010	0.0012	0.0014	0.0016	0.0020	0.0024	0.0028	
0	ρ_1	0.090(3)	0.149(3)	0.192(3)	0.241(3)	—	—	—	—	—	—	0.72–1.03
	ρ_2	0.102(4)	0.134(4)	0.187(5)	0.232(6)	—	—	—	—	—	—	0.97–1.31
	ρ_3	0.098(4)	0.147(4)	0.196(4)	0.236(4)	—	—	—	—	—	—	0.67–1.45
	ρ_4	0.092(4)	0.138(4)	0.183(5)	0.240(6)	—	—	—	—	—	—	0.59–1.09
	$\bar{\rho}$	0.095(2)	0.143(2)	0.190(2)	0.238(2)	—	—	—	—	—	—	—
0.25	ρ_1	0.093(3)	0.134(3)	0.185(3)	0.225(3)	—	—	—	—	—	—	0.49–1.21
	ρ_2	0.089(3)	0.143(4)	0.180(5)	0.238(6)	—	—	—	—	—	—	0.97–1.41
	ρ_3	0.093(4)	0.132(4)	0.184(4)	0.223(4)	—	—	—	—	—	—	0.67–1.31
	ρ_4	0.085(4)	0.147(4)	0.180(5)	0.243(6)	—	—	—	—	—	—	1.01–1.51
	$\bar{\rho}$	0.092(2)	0.138(2)	0.183(2)	0.228(2)	—	—	—	—	—	—	—
0.5	ρ_1	0.090(3)	0.127(3)	0.177(3)	0.215(3)	—	—	—	—	—	—	0.73–1.19
	ρ_2	0.085(3)	0.140(4)	0.173(4)	0.230(6)	—	—	—	—	—	—	0.57–1.20
	ρ_3	0.085(4)	0.128(4)	0.173(4)	0.214(4)	—	—	—	—	—	—	0.64–1.10
	ρ_4	0.091(4)	0.135(4)	0.180(5)	0.222(5)	—	—	—	—	—	—	0.82–0.98
	$\bar{\rho}$	0.088(2)	0.131(2)	0.176(2)	0.218(2)	—	—	—	—	—	—	—
0.75	ρ_1	0.082(3)	0.124(3)	0.166(3)	0.206(3)	0.249(3)	—	—	—	—	—	0.77–0.96
	ρ_2	0.084(3)	0.128(4)	0.167(4)	0.213(5)	0.250(6)	—	—	—	—	—	0.74–0.79
	ρ_3	0.082(4)	0.123(4)	0.166(4)	0.205(4)	0.248(5)	—	—	—	—	—	0.84–1.02
	ρ_4	0.082(4)	0.129(4)	0.163(4)	0.214(5)	0.244(6)	—	—	—	—	—	0.94–1.12
	$\bar{\rho}$	0.082(2)	0.125(2)	0.166(2)	0.208(2)	0.248(2)	—	—	—	—	—	—
1.0	ρ_1	0.081(3)	0.116(3)	0.155(3)	0.193(3)	0.232(3)	—	—	—	—	—	1.02–1.36
	ρ_2	0.073(3)	0.117(4)	0.154(4)	0.194(5)	0.231(6)	—	—	—	—	—	0.75–1.01
	ρ_3	0.083(4)	0.113(4)	0.161(4)	0.191(4)	0.241(5)	—	—	—	—	—	0.36–1.45
	ρ_4	0.072(4)	0.119(4)	0.155(4)	0.196(5)	0.234(6)	—	—	—	—	—	0.75–1.09
	$\bar{\rho}$	0.078(2)	0.116(2)	0.156(2)	0.193(2)	0.234(2)	—	—	—	—	—	—
1.25	ρ_1	0.070(3)	0.108(3)	0.143(3)	0.178(3)	0.216(3)	—	—	—	—	—	0.70–1.13
	ρ_2	0.076(3)	0.110(4)	0.149(4)	0.182(5)	0.222(6)	—	—	—	—	—	0.76–0.95
	ρ_3	0.073(4)	0.111(4)	0.144(4)	0.184(4)	0.217(4)	—	—	—	—	—	0.87–1.23
	ρ_4	0.071(4)	0.108(4)	0.143(4)	0.179(5)	0.213(5)	—	—	—	—	—	0.48–0.94
	$\bar{\rho}$	0.072(2)	0.109(2)	0.144(2)	0.180(2)	0.217(2)	—	—	—	—	—	—
1.5	ρ_1	0.062(3)	0.095(3)	0.130(3)	0.160(3)	0.195(3)	—	—	—	—	—	0.77–1.33
	ρ_2	0.068(3)	0.098(3)	0.132(4)	0.165(4)	0.196(5)	—	—	—	—	—	0.68–1.26
	ρ_3	0.069(4)	0.100(4)	0.129(4)	0.164(4)	0.193(4)	—	—	—	—	—	0.78–1.25
	ρ_4	0.059(4)	0.094(4)	0.127(4)	0.157(4)	0.189(5)	—	—	—	—	—	0.96–0.99
	$\bar{\rho}$	0.064(2)	0.097(2)	0.130(2)	0.161(2)	0.194(2)	—	—	—	—	—	—
1.75	ρ_1	0.052(3)	0.079(3)	0.108(3)	0.133(3)	0.163(3)	0.187(3)	0.217(3)	—	—	—	0.93–1.24
	ρ_2	0.057(3)	0.085(3)	0.113(4)	0.140(4)	0.167(4)	0.194(5)	0.220(5)	—	—	—	0.87–1.20
	ρ_3	0.060(4)	0.075(4)	0.115(4)	0.129(4)	0.170(4)	0.183(4)	0.226(4)	—	—	—	0.66–1.22
	ρ_4	0.049(4)	0.088(4)	0.102(4)	0.141(4)	0.154(4)	0.195(5)	0.204(5)	—	—	—	0.46–1.09
	$\bar{\rho}$	0.055(2)	0.081(2)	0.109(2)	0.135(2)	0.163(2)	0.188(2)	0.217(2)	—	—	—	—
2.0	ρ_1	—	0.057(3)	0.080(3)	0.096(3)	0.118(3)	0.134(3)	0.156(3)	0.194(3)	0.234(3)	0.272(3)	0.95–1.44
	ρ_2	—	0.058(3)	0.073(3)	0.095(3)	0.110(4)	0.132(4)	0.147(4)	0.180(5)	0.215(5)	0.244(6)	0.90–1.94
	ρ_3	—	—	—	—	—	—	—	—	—	—	—
	ρ_4	—	—	—	—	—	—	—	—	—	—	—
	$\bar{\rho}$	—	0.57(2)	0.077(2)	0.096(2)	0.115(2)	0.133(2)	0.153(2)	0.190(2)	0.229(3)	0.266(3)	—
2.1*	ρ_1	—	—	0.062(4)	—	0.091(4)	—	—	0.147(4)	0.176(4)	0.205(4)	1.03–1.56
	ρ_2	—	—	0.059(5)	—	0.087(5)	—	—	0.139(5)	0.162(6)	0.187(7)	1.29–1.76
	ρ_3	—	—	—	—	—	—	—	—	—	—	—
	ρ_4	—	—	—	—	—	—	—	—	—	—	—
	$\bar{\rho}$	—	—	0.061(3)	—	0.089(3)	—	—	0.144(3)	0.172(3)	0.201(3)	—

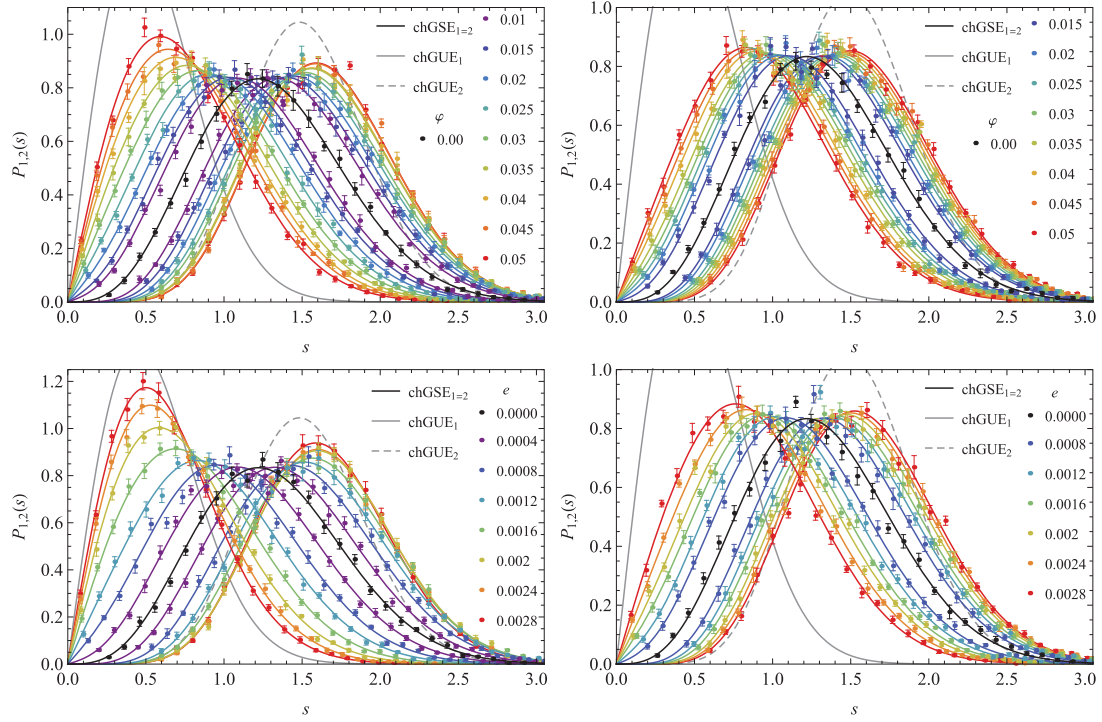


Fig. 3. The first two Dirac eigenvalue distributions $P_{1,2}(s)$ of the SU(2)+ICP model (top) and SU(2) $\times$ U(1) model (bottom) with increasing perturbations φ and e , from black ($\varphi, e = 0$) to purple (smallest φ, e) to red (largest φ, e), on $V = 6^4$ and at $\beta = 1.0$ (left) and at $\beta = 2.0$ (right). Those of chGSE and chGUE are plotted in black and gray, respectively, and the best-fit curves of the chGSE–chGUE crossover are plotted in the same colors as the corresponding lattice data. The error bars of the histograms in Figs. 3–5 are estimated by the jackknife method.

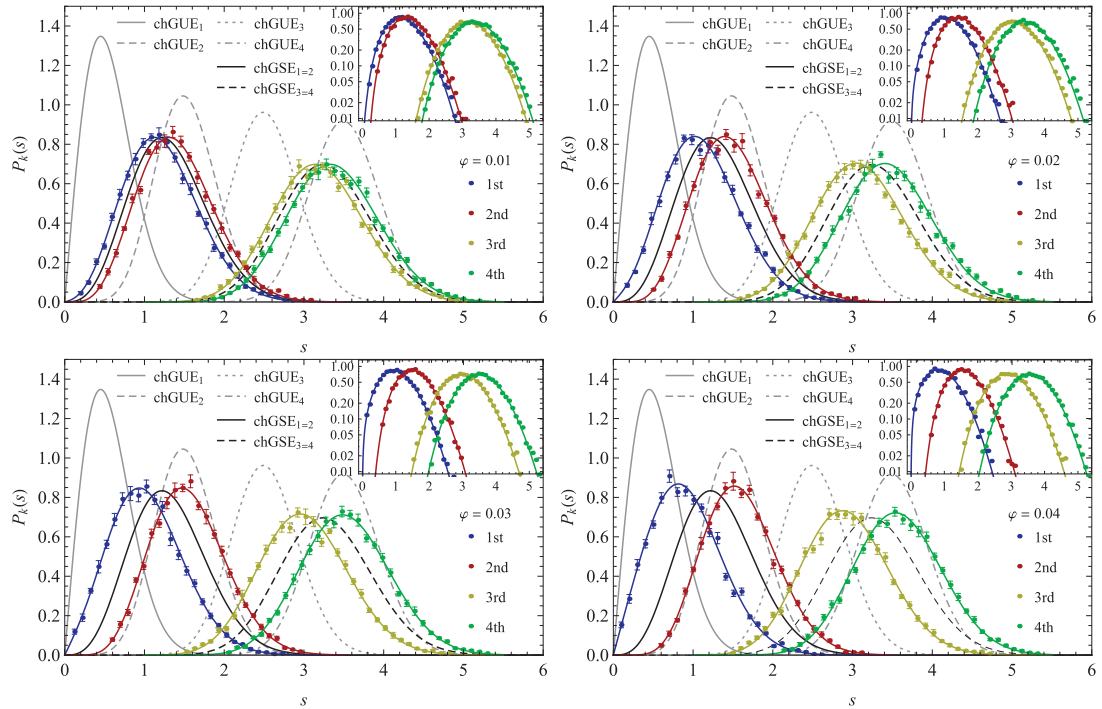


Fig. 4. Linear and logarithmic (inset) plots of the first four Dirac eigenvalue distributions $P_{1,2,3,4}(s)$ (blue, red, yellow, green) of the SU(2)+ICP model on $V = 6^4$ and at $\beta = 1.75$, and best-fit curves of the chGSE–chGUE crossover. Top left: $\varphi = 0.01$, top right: $\varphi = 0.02$, bottom left: $\varphi = 0.03$, bottom right: $\varphi = 0.04$.

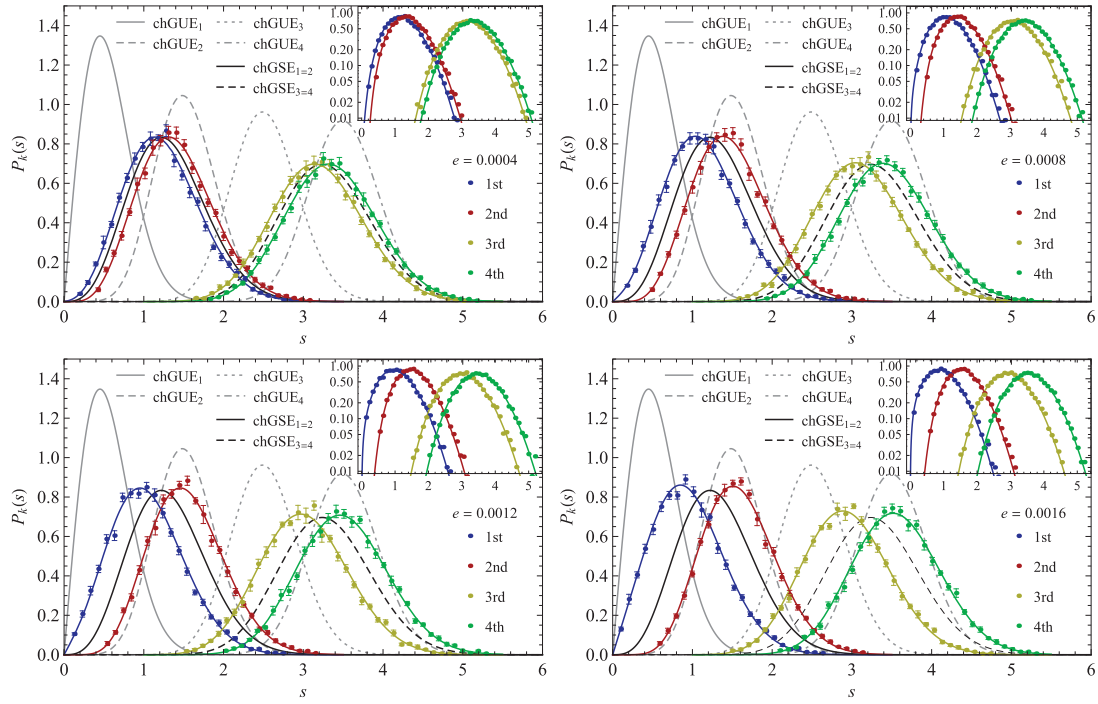


Fig. 5. Linear and logarithmic (inset) plots of the first four Dirac eigenvalue distributions $P_{1,2,3,4}(s)$ (blue, red, yellow, green) of the $SU(2) \times U(1)$ model on $V = 6^4$ and at $\beta = 1.75$, and best-fit curves of the chGSE–chGUE crossover. Top left: $e = 0.0004$, top right: $e = 0.0008$, bottom left: $e = 0.0012$ and bottom right: $e = 0.0016$.

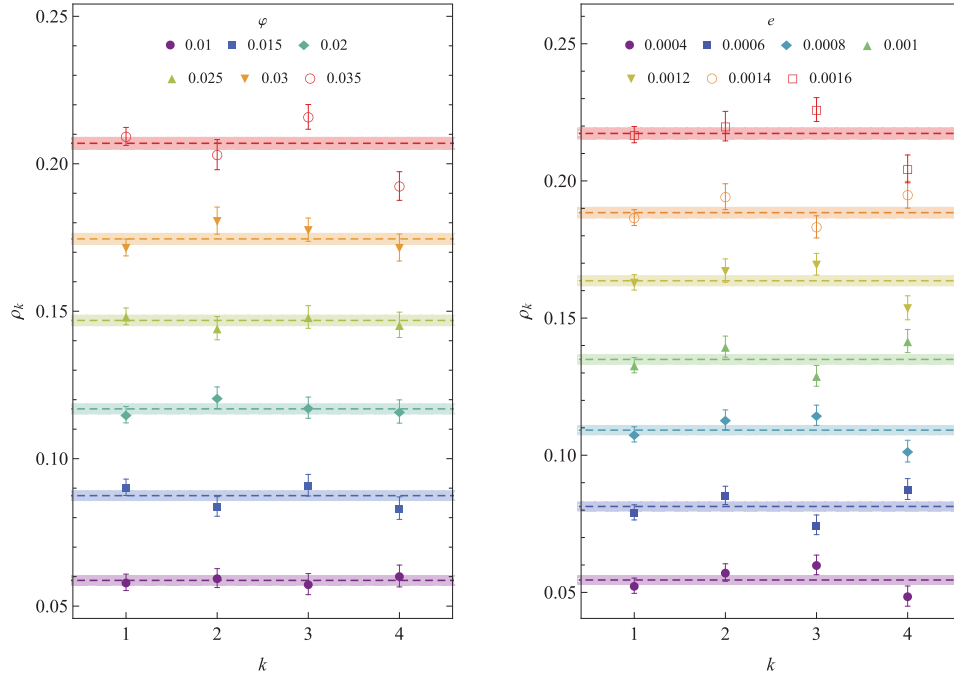


Fig. 6. Crossover parameters $\rho_{1,2,3,4}$ determined from $P_{1,2,3,4}(s)$. Their weighted averages and combined errors are shown in horizontal lines and strips. Left: $SU(2)+ICP$ model with $\varphi = 0.01–0.035$ (purple to red). Right: $SU(2) \times U(1)$ model with $e = 0.0004–0.0016$ (purple to red). Both are at $\beta = 1.75$ and on $V = 6^4$.

SU(2)+ICP model in Tables 3 and 4, and (b) for the SU(2) $\times$ U(1) model in Tables 5 and 6. Again we adopted only the cases that passed the χ^2 test with $\chi^2/\text{d.o.f.} < 2.0$, and the exceptional treatment of cutting off the tails of the smallest eigenvalue distributions for which $p_1(s) < 0.2$ from fitting is applied to $\beta = 1.75$ on $V = 4^4$ and $\beta = 2.1$ on $V = 6^4$ (marked by * in Tables 3–6). Whenever the histogram of the unfolded eigenvalues $s_{2i-1} = s_{2i}$ of the pure SU(2) Dirac operator is well fitted into the chGSE prediction in our criterion, so is the corresponding pair $\{P_{2i-1}(s_{2i-1}), P_{2i}(s_{2i})\}$ of the perturbed Dirac operator to the chGSE–chGUE crossover prediction, as expected.

Figure 3 shows samples of the histograms $\{P_1(s_1), P_2(s_2)\}$ of the perturbed Dirac operator eigenvalues and best-fit distributions of the chGSE–chGUE crossover. As the perturbation φ or e is increased, the eigenvalue distributions are clearly seen to follow the random matrix curves. They respond more rapidly to the perturbation for smaller β (left panels) than for larger β (right panels), indicating that F is a decreasing function of β . For the sake of graphical visibility, we also exhibit samples of histograms of $\{P_1(s_1), \dots, P_4(s_4)\}$ and best-fit distributions of the chGSE–chGUE crossover, at fixed $\beta = 1.75$ and with increasing perturbations φ (Fig. 4) and e (Fig. 5), respectively.

Figure 6 illustrates the crossover parameters ρ_k determined from $P_k(s_k)$ for the SU(2)+ICP model and for the SU(2) $\times$ U(1) model at $\beta = 1.75$ and $V = 6^4$, at each value of perturbation φ or e . We notice that ρ_1 and ρ_2 (ρ_3 and ρ_4) have a tendency to counter-move across the weighted average of the four, as anticipated in Sect. 3.1. Combined use of $P_{2i-1}(s_{2i-1})$ and $P_{2i}(s_{2i})$ is indeed seen to reduce the errors of the best-fit parameters $\bar{\rho}$ in both panels.

3.2.2. Low-energy constants

The chiral condensate Σ is determined by Eq. (54) from $\bar{\Delta}$ summarized in Tables 1 and 2. In Table 7 we list these values in the third column (SU(2)+ICP) and in the fifth column (SU(2) $\times$ U(1)), at each β

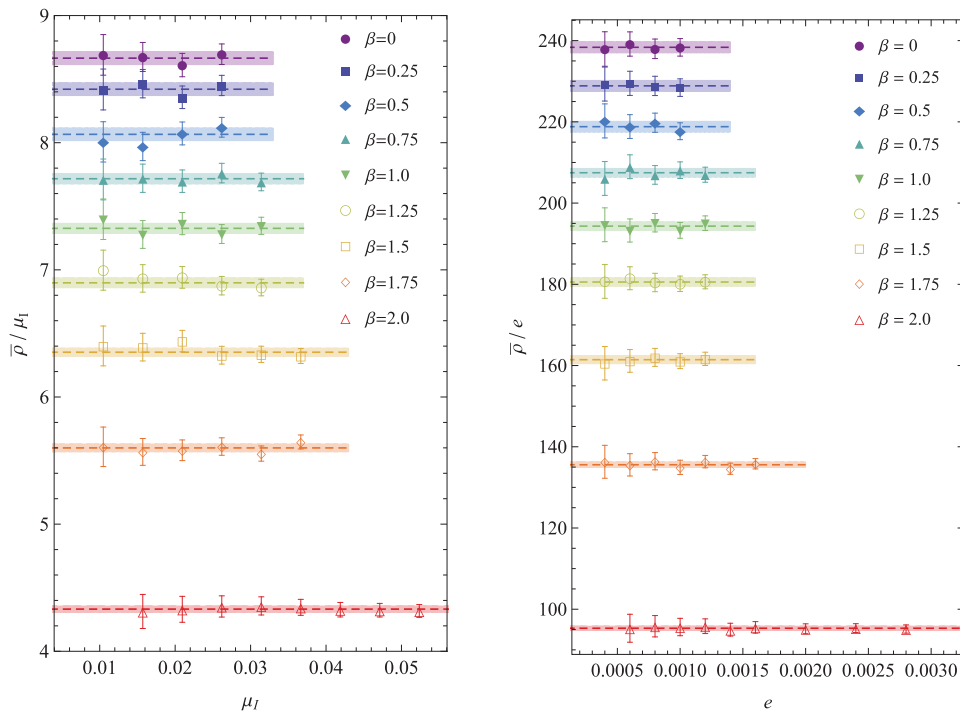


Fig. 7. Ratios between the crossover parameter and the strength of perturbation: $\bar{\rho}/\mu_1$ for the SU(2)+ICP model (left) and $\bar{\rho}/e$ for the SU(2) $\times$ U(1) model (right), on $V = 6^4$ and at various β . Colored dashed lines and strips represent the weighted averages and combined errors of the ratios.

Table 7. Low-energy constants Σ and F derived from the SU(2)+ICP model and SU(2) $\times$ U(1) model on $V = 4^4, 6^6$, and their extrapolation to the thermodynamic limit (TDL).

V	β	SU(2)+ICP		SU(2) $\times$ U(1)		
		Σa^3	$F^2 a^2$	Σa^3	$F^2 \mu_1^2 a^4 / e^2$	$F^2 \mu_1^2 a^4 / (e^2 V)$
4^4	0	1.318 0(4)	0.289(2)	1.317 9(4)	40.8(2)	0.159 4(8)
	0.25	1.264 5(4)	0.267(1)	1.264 1(4)	37.6(2)	0.146 8(7)
	0.5	1.205 7(4)	0.244(1)	1.207 0(4)	34.1(2)	0.133 1(7)
	0.75	1.141 6(4)	0.226(1)	1.139 5(4)	30.8(2)	0.120 3(6)
	1.0	1.066 9(5)	0.201(1)	1.064 9(5)	26.9(1)	0.105 1(6)
	1.25	0.978 3(7)	0.176(1)	0.979 8(7)	22.9(1)	0.089 3(4)
	1.5	0.871 3(2)	0.146 3(9)	0.869(2)	18.3(1)	0.071 5(4)
	1.75*	0.720(2)	0.109(1)	0.723(1)	12.8(1)	0.050 0(4)
6^4	0	1.309 2(8)	0.286(3)	1.307 5(8)	216(2)	0.167(2)
	0.25	1.257 4(8)	0.270(3)	1.257 3(8)	199(2)	0.154(2)
	0.5	1.198 2(9)	0.248(3)	1.199 6(8)	182(2)	0.141(2)
	0.75	1.134 2(7)	0.227(2)	1.134 1(7)	164(2)	0.126(1)
	1.0	1.063 9(7)	0.204(2)	1.060 4(7)	144(2)	0.111(1)
	1.25	0.977 4(6)	0.181(2)	0.977 6(6)	124(1)	0.096(1)
	1.5	0.871 4(6)	0.154(2)	0.871 6(6)	99(1)	0.076 6(9)
	1.75	0.727 9(7)	0.119(1)	0.728 6(7)	70.0(7)	0.054 0(5)
	2.0	0.502(2)	0.071 5(8)	0.505(2)	34.6(4)	0.026 7(3)
	2.1*	0.379(1)	0.046 5(9)	0.379(2)	19.8(4)	0.015 3(3)
TDL	0	1.291 6(4)	0.280(2)	1.286 5(4)	—	0.181 9(7)
	0.25	1.243 2(4)	0.275(1)	1.243 5(4)	—	0.168 1(7)
	0.5	1.183 2(4)	0.254(1)	1.184 8(3)	—	0.155 9(6)
	0.75	1.119 5(3)	0.229(1)	1.123 4(3)	—	0.138 7(6)
	1.0	1.057 9(4)	0.212(1)	1.051 3(4)	—	0.122 7(4)
	1.25	0.975 6(5)	0.191 2(9)	0.973 2(5)	—	0.108 8(4)
	1.5	0.871 7(5)	0.168 4(8)	0.876 9(5)	—	0.086 7(4)
	1.75*	0.743 9(6)	0.140 2(7)	0.739 9(6)	—	0.061 9(3)

and on $V = 4^4$ and 6^4 lattices. In addition, the values of the chiral condensate in the thermodynamic limit $V = \infty$, extrapolated from $V = 4^4$ and 6^4 , are listed for the coupling range $0 \leq \beta \leq 1.75$.³

The pseudo-scalar decay constant F in Eqs. (55) and (56) is essentially the proportionality constant between the perturbation strength μ_1 or e and the crossover parameter ρ . In Fig. 7 we exhibit the ratios $\bar{\rho}/\mu_1$ for the SU(2)+ICP model (left) and the $\bar{\rho}/e$ for SU(2) $\times$ U(1) model (right), both on $V = 6^4$ and at various β . For each model we observe excellent stability of the ratios as the perturbation μ_1 or e is varied. This justifies the use of combining $\bar{\rho}/\mu_1$ or $\bar{\rho}/e$ for all available μ_1 or e at each β , so that errors in the combined ratios are significantly reduced. Pseudo-scalar decay constants determined in this way are exhibited in the fourth column (SU(2)+ICP) and in the sixth column (SU(2) $\times$ U(1)) of Table 7. For the SU(2)+ICP model, Σ and F^2 are determined with 10^{-4} and 10^{-3} precision, respectively, on our small lattices of 4^4 and 6^4 . This observation, first advocated in Ref. [8] in the context of SU(3) gauge theory with isospin ICP, enables us to extrapolate their value to the thermodynamic limit $V \rightarrow \infty$, listed in the bottom rows of Table 7.

³ Although the fitting range of $P_k(s)$ at $\beta = 1.75$ on $V = 4^4$ is compromised as compared to $\beta \leq 1.5$, we dare to estimate the thermodynamic limit of the chiral condensate and pseudo-scalar decay constant using these data.

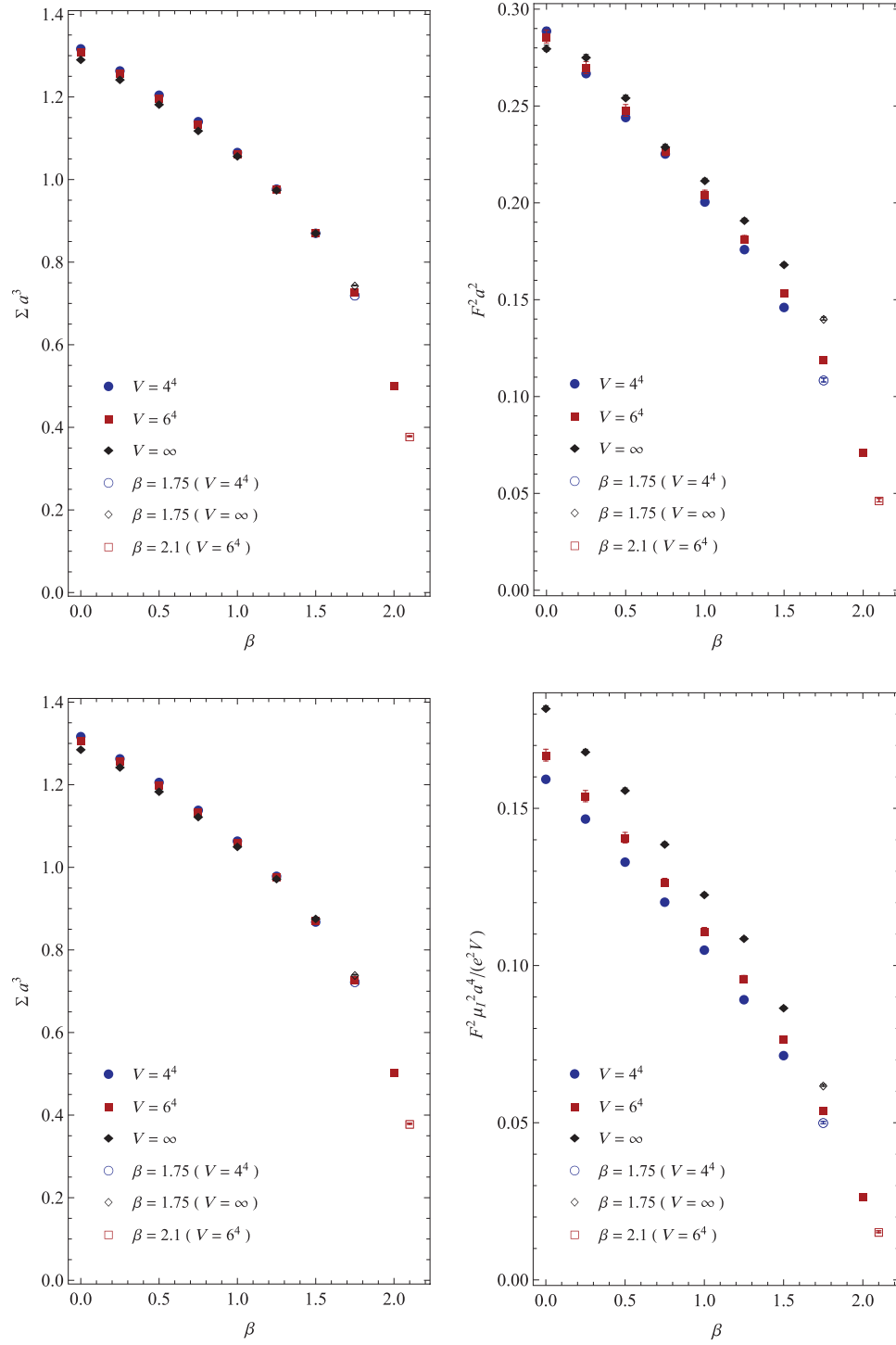


Fig. 8. Coupling dependence of low-energy constants Σ (left) and F^2 (right) derived from the SU(2)+ICP model (top) and from the SU(2)×U(1) model (bottom). Blue and red filled symbols represent the values on $V = 4^4$ and 6^4 , and black filled symbols their extrapolated values to $V \rightarrow \infty$. Empty symbols correspond to the cases marked * in Tables 1–7 (using limited ranges for fitting $P_k(s)$) and extrapolations thereof.

On the other hand, for the $SU(2) \times U(1)$ model the combination $F^2 \mu_1^2 / e^2$ is found to scale linearly with the lattice volume V . Thus we listed the extrapolated values of $F^2 \mu_1^2 / (e^2 V)$ in the thermodynamic limit.

Finally, in Fig. 8 we exhibit $SU(2)$ -coupling dependences of the low-energy constants on $V = 4^4, 6^4$, and ∞ in the $SU(2)$ -ICP model (top) and in the $SU(2) \times U(1)$ model. The error bars are so small that they are almost obscured by the symbols.

4. Conclusions

We have analytically evaluated the k th smallest eigenvalue distributions $p_k(s)$ for a random matrix ensemble interpolating chGSE and chGUE using a Nyström-type method applied to the Fredholm Pfaffian and resolvents of the quaternion kernel. These random matrix results are applied to fit the spectra of fundamental, staggered Dirac operators of $SU(2)$ gauge theory with imaginary chemical potential and of $SU(2) \times U(1)$ gauge theory on small lattices, from the strong-coupling to the near-scaling regions. Combined use of the first four nondegenerate Dirac eigenvalue distributions in place of the spectral density or the smallest eigenvalue distribution of unperturbed $SU(2)$ gauge theory enables us to determine the chiral condensate with $\mathcal{O}(10^{-4})$ precision. Excellent one-parameter fitting of $\chi^2/\text{d.o.f.} < 2$ between *non-cumulative* individual distributions and eigenvalue histograms is achieved for almost all cases of $U(1)$ perturbations. Combined use of the first four eigenvalue distributions also contributed to a reduction of the errors in the crossover parameter ρ . The acute sensitivity of $p_k(s)$ on ρ , and the observed linear dependence of ρ on the perturbation strength (AB flux φ or $U(1)$ coupling e) resulted in determination of the pseudo-scalar decay constant F (i.e., the coefficient of the pseudoreality-breaking term) with $\mathcal{O}(10^{-3})$ precision.

Our method of determining F in QCD-like theories, which has proved to be feasible on relatively small-sized lattices, is clearly advantageous over the conventional method of using axial current correlators, which inevitably requires a large temporal dimension. A possible application of our method would be towards technicolor candidate gauge theories with fermions in (pseudo)real representations, such as $SU(N)$ gauge theory with two adjoint flavors [35], which corresponds to the chGSE class if simulated with an overlap Dirac operator [7]. Hyper-precise determination of its “low-energy” constants (i.e., Higgs couplings) from lattice simulations using our ICP method would, upon comparison with knowledge from collider experiments, contribute to single out a credible scenario from such BSM candidates.

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