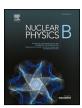
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Quantum Field Theory and Statistical Systems



Phase transitions in the decomposition of SU(N) representations

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ABSTRACT

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We study the multiplicity of irreducible representations in the decomposition of n fundamentals of SU(N) weighted by a power of their dimension in the large n and large N double scaling limit. A nontrivial scaling is obtained by keeping n/N^2 fixed, which plays the role of an order parameter. We find that the system generically undergoes a fourth order phase transition in this parameter, from a dense phase to a dilute phase. The transition is enhanced to third order for the unweighted multiplicity, and disappears altogether when weighting with the first power of the dimension. This corresponds to the infinite temperature partition function of non-Abelian ferromagnets, and the results should be relevant to the thermodynamic limit of such ferromagnets at high temperatures.

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1. Introduction

Nonabelian unitary groups SU(N) play a crucial role in particle physics [1,2], and, indirectly through matrix models [3], in string theory and gravity. Ungauged and gauged SU(2) and SU(3) groups are the most common, representing spin, flavor, or color degrees of freedom. Matrix models, on the other hand, involve systems invariant under a larger SU(N) symmetry, and eventually N is taken to infinity to achieve a particular scaling limit.

Independently, magnetic systems with SU(N) symmetry have been considered in the context of ultracold atoms [4–8], spin chains [9,10], or of interacting atoms on lattice sites [11–17] and in the presence of SU(N) magnetic fields [18–20]. Such systems consist of a large number of components ("atoms"), each carrying an irreducible representation of the symmetry group, and the full system transforms under the same symmetry in the direct sum of the representations of the components. The decomposition of the states of the full system into irreducible representations (irreps) of the symmetry group is, then, of physical relevance.

In previous work we derived the statistics of such decompositions [21] and investigated the properties of n coupled SU(N) atoms in the ferromagnetic regime, that is, in the regime where the mutual interaction of SU(N) components would tend to "align" their charges, also in the presence of an external nonabelian magnetic field coupled to the system [22]. We studied the system in its thermodynamic limit of a large number of atoms $n \gg 1$ and uncovered a rich phase structure, qualitatively and quantitatively different from that of usual SU(2) ferromagnets, involving various critical temperatures, hysteresis effects, coexistence of phases, and latent heat transfer during phase transitions.

The above raise the obvious question of the properties of such systems when both the large-n (thermodynamic) and large-N limits are taken. This is the topic of the present work. We consider the number of irreducible components in the decomposition of n fundamental irreps of SU(N) weighted by a (positive) power of their dimension, which can be viewed as the infinite-temperature partition function of an exotic ferromagnet. We derive the large-(N,n) properties of this quantity in the appropriate scaling in that limit, which requires $n/N^2 \sim 1$ as $N,n \gg 1$, and demonstrate that, in that limit, the representation content of the system undergoes a phase transition in this parameter. This transition is model-independent, as the Hamiltonian becomes irrelevant in the infinite-temperature limit, and thus represents a universal feature of such systems at high temperatures. Although the large-N limit may appear somewhat unnatural for physical applications, the weaker scaling of $N \sim \sqrt{n}$ required for this limit allows it to be achieved for reasonable values of N even for large systems.

In the following sections, we review the relevant mathematics of decomposing a large number of fundamental SU(N) representations into irreducible components, stressing the fermion momentum parametrization of irreps and related duality properties. Subsequently, we consider the multiplicity of each irrep weighted by various powers of the number of its states (dimension) in the large-(N,n) limit and solve for the saddle point irrep that maximizes it. We uncover a *fourth-order* phase transition in the order parameter $t = n/(4N^2)$ in the generic situation, which becomes a stronger, third-order transition between a duality preserving and a duality breaking phase for the unweighted multiplicity, and disappears altogether when weighting by the number of states. We will conclude with some remarks on possible applications and directions for further investigation.

2. Fermions and SU(N) group theory: a brief review

We review here the description of irreducible representations of SU(N) as N-fermion energy eigenstates on the circle [23] and the corresponding composition rules in the fermion picture focusing on the results relevant for our considerations.

The correspondence of irreps of SU(N) and N-fermion energy eigenstates is most readily established by considering the action of a particle moving freely on the group manifold U(N) [24] with Lagrangian

$$\mathcal{L} = -\frac{1}{2} \text{Tr} \left(U^{-1} \dot{U} \right)^2 , \qquad (2.1)$$

where U is an N-dimensional unitary matrix and overdot signifies time derivative. Classically, the particle performs geodesic motion on U(N). Quantum mechanically, energy eigestates are matrix elements of irreps of U(N), the energy corresponding to the quadratic Casimir of the irrep.

The Lagrangian (2.1) is invariant under unitary conjugations of U (as well as left- and right-multiplications by constant unitary matrices), with a conserved generator $P=i[U^{-1},\dot{U}]$, so we can impose the constraint that this conserved generator vanishes. Quantum mechanically, this amounts to choosing states invariant under unitary conjugations of U. Energy eigenstates in this subspace become the conjugation-invariant linear combinations $\sum_a r_{aa}(U) = \mathrm{Tr} r(U) = \chi_r(U)$, the latter being the character of the irrep, and depend only on the eigenvalues of the matrix U.

Classically, we can implement the conjugation invariance constraint at the Lagrangian level by setting

$$U(t) = V \operatorname{diag}\{z_j\} V^{-1}, \quad z_j := e^{ix_j}, \qquad x_j \equiv x_j + 2\pi.$$
 (2.2)

The vanishing of P implies that V can be chosen time independent, and (2.1) becomes the Lagrangian of N free particles on the unit circle with coordinates x_j . Since exchanging the eigenvalues is a special case of unitary conjugation, states $\phi(x_1,\ldots,x_N)$ are invariant under exchange of the x_j and are, in principle, bosonic. However, upon quantization, the change of variables from U to x_j introduces in the measure the absolute square of the Vandermonde determinant $|\Delta(\mathbf{z})|^2$, with

$$\Delta(\mathbf{z}) = (z_1 \cdots z_N)^{-\frac{N-1}{2}} \begin{vmatrix} z_1^{N-1} & z_1^{N-2} & \cdots & z_1 & 1 \\ z_1^{N-1} & z_1^{N-2} & \cdots & z_2 & 1 \\ \vdots & \vdots & \ddots & \vdots & \vdots \\ z_N^{N-1} & z_N^{N-2} & \cdots & z_N & 1 \end{vmatrix} . \tag{2.3}$$

(boldface \mathbf{z} stands for the full set of variable z_1, \dots, z_N , and similarly for other sets of N quantities). The prefactor $(z_1 \cdots z_N)^{-\frac{N-1}{2}}$ is a pure phase introduced to make $\Delta(\mathbf{z})$ invariant under uniform shifts of the x_i 's. Upon incorporating one factor $\Delta(\mathbf{z})$ into the wavefunction

$$\psi(\mathbf{z}) = \Delta(\mathbf{z}) \, \phi(\mathbf{z}) \,, \tag{2.4}$$

the integration measure in the x_j becomes flat and the Hamiltonian becomes the standard free N-particle Hamiltonian on the x_j . Because of the prefactor $\Delta(\mathbf{z})$ the states $\psi(\mathbf{z})$ are antisymmetric upon interchanging any two of the x_i . This establishes the correspondence between the conjugation-invariant sector of the model (2.1) and free fermions on the circle, and thus of irreps of SU(N) and free fermion energy eigenstates. In particular, the fermion energy eigenstate corresponding to an irrep r with energy $E = C_2(r)$ in terms of the corresponding quadratic Casimir becomes

$$\psi_r(\mathbf{z}) = \Delta(\mathbf{z}) \, \chi_r(\mathbf{z}) \,, \tag{2.5}$$

with $\chi_r(\mathbf{z})$ the character of irrep r in terms of the eigenvalues z_i of U.

The single-particle spectrum on the circle consists of discrete momentum eigenstates with eigenvalue $k=0,\pm 1,\pm 2,\ldots$ and energy $E_k=k^2/2$. An N-fermion energy eigenstate corresponds to filling n of the single-particle states with fermions. Call $k_1>k_2>\cdots>k_N$ the momenta of these states in decreasing order. The (unnormalized) wavefunction corresponding to this state is given by the Slater determinant

$$\psi_{\mathbf{k}}(\mathbf{z}) = \begin{vmatrix} z_1^{k_1} & z_1^{k_2} & \cdots & z_1^{k_{N-1}} & z_1^{k_N} \\ z_1^{k_1} & z_1^{k_2} & \cdots & z_1^{k_{N-1}} & z_1^{k_N} \\ z_2^{k_1} & z_2^{k_2} & \cdots & z_2^{k_{N-1}} & z_2^{k_N} \\ \vdots & \vdots & \ddots & \vdots & \vdots \\ z_N^{k_1} & z_N^{k_2} & \cdots & z_N^{k_{N-1}} & z_N^{k_N} \end{vmatrix} .$$

$$(2.6)$$

The total momentum of the fermions $K=k_1+\cdots+k_n$ corresponds to the states picking up a phase e^{icK} upon the shift $x_j\to x_j+c$, that is, upon $U\to e^{ic}U$. It thus represents the U(1) charge of the state. We may shift all momenta by a constant, changing k and the U(1) charge without affecting the SU(N) part of the states, which can then be labeled by the N-1 shift-invariant integers $k_1-k_N>k_2-k_N>\cdots>k_{N-1}-k_N>0$. Alternatively, we can neutralize the U(1) charge by introducing the prefactor $(z_1\dots z_N)^{-\sum_i k_i/N}$ in (2.6), similarly to the prefactor introduced in $\Delta(\mathbf{z})$. With this additional prefactor, (2.6) maps to (2.3) for the singlet representation for which $k_i=N-i$, $i=1,\dots,N$.

Finally, we note the correspondence with the standard Young tableau. This is done by expressing $k_j - k_N$ in terms of variables ℓ_j as

$$\ell_{j} = k_{j} - k_{N} + j - N , \qquad \ell_{1} \geqslant \ell_{2} \geqslant \cdots \geqslant \ell_{N-1} \geqslant 0 . \tag{2.7}$$

The non-negative, ordered integers ℓ_j represent the length of rows $j=1,2,\ldots,N-1$ of the Young tableau of the irrep corresponding to the fermionic state. The transition from k_j to ℓ_j is, in fact, bosonization (in the sense that any two consecutive ℓ , unlike the k, can be equal), the ℓ_j corresponding to the possible momenta of N bosons on the circle.

2.1. Composition of (fundamental) representations

Consider the direct product $r_1 \times r_2$ of two (possibly reducible) representations r_1 and r_2 . The basic relation

$$\chi_{r_1 \times r_2}(U) = \text{Tr}(r_1 \times r_2)(U) = \text{Tr}(r_1(U)) + \text{Tr}(r_2(U)) = \chi_{r_1}(U) \chi_{r_2}(U), \qquad (2.8)$$

implies, through ((2.4) and (2.5)), that their corresponding fermion states are related as

$$\psi_{r_1 \times r_2}(\mathbf{z}) = \Delta(\mathbf{z}) \chi_{r_1}(\mathbf{z}) \chi_{r_2}(\mathbf{z}) = \frac{\psi_{r_1}(\mathbf{z}) \psi_{r_2}(\mathbf{z})}{\Delta(\mathbf{z})} = \psi_{r_1}(\mathbf{z}) \chi_{r_2}(\mathbf{z}) = \psi_{r_2}(\mathbf{z}) \chi_{r_1}(\mathbf{z}). \tag{2.9}$$

Here we need the case of the fundamental f for which the character is simply

$$\chi_f(\mathbf{z}) = \sum_{i=1}^{N} z_i \ . \tag{2.10}$$

¹ In standard physics convention, we use the term tableau instead of the term diagram that is used in mathematics, where tableau stands for a diagram with properly ordered integer entries in its boxes.

The composition of two irreps $r \otimes f$, the second one being the fundamental, is then straightforward. The fermion wavefunction of the state corresponding to $r \otimes f$ is, using (2.9) and (2.10), simply

$$\psi_{r\otimes f}(\mathbf{z}) = \psi_r(\mathbf{z})\chi_f(\mathbf{z}) = \psi_r(\mathbf{z})\sum_{i=1}^N z_i. \tag{2.11}$$

This can be used to obtain the fermionic state corresponding to the composition of several (n in number) fundamental irreps f. The original, singlet state is simply $\Delta(\mathbf{z})$ and an iteration of the above formula yields

$$\psi_{N,n}(\mathbf{z}) = \Delta(\mathbf{z}) \left(\sum_{i=1}^{N} z_i \right)^n := \sum_{\mathbf{k}} d_{n;\mathbf{k}} \prod_{i=1}^{N} z_i^{k_i} . \tag{2.12}$$

Since $\psi_{N,n}$ is antisymmetric in the z_i , the coefficients $d_{n;\mathbf{k}}$ appearing above are fully antisymmetric in the k_i . When the k_i are in decreasing order, $d_{n;\mathbf{k}}$ gives the multiplicity of the irrep labeled by $k_1 > \cdots > k_N$.

To derive an explicit combinatorial expression for the multiplicity we first focus on the coefficients produced by the term $(z_1 + \cdots + z_N)^n$, denoted by $D_{n:\mathbf{k}}$. We have

$$\left(\sum_{i=1}^{N} z_{i}\right)^{n} := \sum_{k_{1},\dots,k_{N}} D_{n;\mathbf{k}} \prod_{i=1}^{N} z_{i}^{k_{i}}, \qquad D_{n;\mathbf{k}} = \delta_{k_{1}+\dots+k_{N},n} \frac{n!}{\prod_{i=1}^{N} k_{i}!}, \qquad k_{i} \geqslant 0.$$

$$(2.13)$$

Incorporating the Vandermonde factor in (2.12) we eventually obtain [21]

$$d_{n;\mathbf{k}} = n! \frac{\prod_{j>i=1}^{N} (k_i - k_j)}{\prod_{i=1}^{N} k_i!} = n! \frac{\Delta(\mathbf{k})}{\prod_{i=1}^{N} k_i!} \quad \text{with} \quad \sum_{i=1}^{N} k_i = n + \frac{N(N-1)}{2} \ . \tag{2.14}$$

The dimension of the irrep expressed in terms of the k_i 's becomes

$$\dim(\mathbf{k}) = \prod_{j>i=1}^{N} \frac{k_i - k_j}{j - i} = \frac{\Delta(\mathbf{k})}{\prod_{s=1}^{N-1} s!} . \tag{2.15}$$

2.2. Momentum density and a group duality

We conclude by giving a "second quantized" expression for the d_{n,k_1,\ldots,k_N} that is useful in the large-N,n limit. Thinking of the k_i as a distribution of fermions on the positive momentum lattice $s=0,1,\ldots$, we define the discrete momentum density of fermions ρ_s equal to one on points s of the momentum lattice where there is a fermion and zero elsewhere, that is,

$$\rho_{s} = \sum_{i=1}^{N} \delta_{s,k_{i}} . \tag{2.16}$$

Clearly ρ_s , and in accordance with (2.14), satisfies the relations

$$\sum_{s=0}^{M} \rho_s = N , \qquad \sum_{s=0}^{M} s \, \rho_s = K = n + \frac{N(N-1)}{2} , \qquad (2.17)$$

where M is a cutoff momentum that can be chosen arbitrarily as long as it is bigger than all the k_i 's. Then, it can be easily seen that (2.14) can be written as [21]

$$d_{n,\mathbf{k}} = n! \prod_{t>s=0}^{M} (t-s)^{(\rho_s-1)\rho_t} . \tag{2.18}$$

The integer M could in principle be taken to infinity. However, keeping it finite serves to demonstrate an interesting particle-hole duality of the formulae. Define

$$\tilde{\rho}_s = 1 - \rho_{M-s}$$
, $s = 0, 1, \dots, M$. (2.19)

Clearly $\tilde{\rho}_s$ is the density of holes on the lattice [0,M] with the momentum reversed. Moreover, using (2.17), $\tilde{\rho}_s$ satisfies

$$\sum_{s=0}^{M} \tilde{\rho}_s = M - N + 1 , \qquad \sum_{s=0}^{M} s \, \tilde{\rho}_s = n + \frac{(M - N + 1)(M - N)}{2} . \tag{2.20}$$

Therefore, $\tilde{\rho}_s$ represents an irrep of SU(M-N+1) with the same excitation n (total number of boxes) but with the rows in the Young tableau of the SU(N) irrep turned into columns for SU(M-N+1), which defines the dual irrep. One can check that

$$d_{n,\mathbf{k}} = n! \prod_{t>s=0}^{M} (t-s)^{(\tilde{\rho}_s-1)\tilde{\rho}_t} . \tag{2.21}$$

That is, in the decomposition of the tensor product of n fundamentals of SU(N), the multiplicity of any given irrep is the same as the one for its dual irrep in the product of n fundamentals of SU(M-N+1). Note that this relation holds for any M such that $M \ge k_1 > k_2 > \cdots > k_N$.

This duality between SU(N) and SU(M-N+1) can be turned into a self-duality if we choose M=2N-1, which is possible if $k_1 \le 2N-1$, that is, $\ell_1 \le N$. This will be guaranteed in the case $n \le N$ since then, indeed, $\ell_1 \le n \le N$.

3. Large-N, n limits

The two parameters at our disposal in the SU(N) case, namely N and n, can be taken to be large in various ways, leading to different large-N, n limit regimes. The dimensionality of the Hilbert space of the system is N^n , and both the $n \gg 1$ and the $N \gg 1$ limits are driving it to infinity, although in qualitatively different ways: the limit $n \gg 1$ can be viewed as a standard thermodynamic limit increasing the number of individual components (fundamental irreps) of the system, while the $N \gg 1$ limit swells the available phase space per degree of freedom. The relative scaling between n and N becomes important and, as we shall see, the distribution of irreps undergoes a phase transition at some critical line $n \sim N^2$ in the two-dimensional parameter space spanned by N and n. The exact critical line is fixed by the statistical quantity of interest in the problem, namely, number of irreps, number of states, or a more general combination.

In the following we will analyze the large-n, N limit of the distribution of irreps, derive the dominant distribution, establish the existence of a phase transition, and determine the order of the transition.

3.1. $n \gg N \sim 1$

The thermodynamic limit $n \gg N \sim 1$ is the most straightforward. It was derived in [21] and used in [22] to analyze the SU(N) ferromagnet and determine its intricate phase transition diagram. We can use the Stirling approximation in the combinatorial formula for $d_{n,k}$ given in (2.14). In that limit, $k_i \gg 1$ since from (2.14) their sum is of order n. The result is

$$d(N, n; \mathbf{k}) = \prod_{s=1}^{N-1} s! \frac{N^{n+N^2/2} e^{\frac{N^2(N^2-1)}{24n}}}{\sqrt{2\pi}^{N-1} n^{(N^2-1)/2}} \dim(\mathbf{k}) e^{-\frac{N}{n} c_2(\mathbf{k})},$$
(3.1)

with the dimension of the irrep and the corresponding quadratic Casimir given by (2.14) and

$$c_2(\mathbf{k}) = \frac{1}{2} \sum_{i=1}^{N} \left(k_i - \frac{K}{N} \right)^2 - \frac{N(N^2 - 1)}{24} , \qquad (3.2)$$

where factors subleading in 1/n are eliminated. Distribution (3.1) implies that in the limit $n \gg 1$ the deviations of k_i from their mean value $K/N \sim n/N$ scale as

$$k_i - \frac{K}{N} \sim \sqrt{n} \ . \tag{3.3}$$

3.2. $N, n \gg 1$

The limit where both N and n are large is more interesting. It that limit it does not make sense any more to define a continuous distribution $d(N, k_1, \ldots, k_N)$, as the dimensionality of the space of k_i grows to infinity for large N. Instead, it is possible to define a *density of points* $\rho(k)$ that is the continuous version of ρ_s defined in (2.16) smoothed over the position of lattice points s around momentum s, and express the number of irreps s a functional of this density s. Alternatively, s can be defined through the continuous momentum function

$$k(j) \simeq k_j \ (j = 1, ..., N), \qquad \rho(k) = -\frac{dj}{dk(j)} \text{ at } k(j) = k$$
 (3.4)

(the minus sign in the definition of $\rho(k)$ is needed to ensure a positive ρ since k_j is decreasing with j). Then, using the continuous version of (2.18), the logarithm of the number of irreps $d_{n,\mathbf{k}} := d_n[\rho(k)]$ becomes in the continuous approximation

$$\ln \frac{d_n[\rho(k)]}{n!} = \int_0^\infty dk \int_0^k dk' \, \rho(k) \rho(k') \ln(k - k') - \int_0^\infty dk \int_0^k dk' \, \rho(k) \ln(k - k'), \tag{3.5}$$

up to a constant of $\mathcal{O}(1)$ and where we have moved the ρ -independent term $\ln n!$ to the left hand side. Rewriting the double integral in a way symmetric in k and k' and performing the integral over k' in the second term, we obtain

$$\ln \frac{d_n[\rho(k)]}{n!} = \frac{1}{2} \int_0^\infty dk \int_0^\infty dk' \, \rho(k) \rho(k') \ln |k - k'| - \int_0^\infty dk \, \rho(k) \, k (\ln k - 1) \,. \tag{3.6}$$

Integrals involving singular kernels, such as $\ln |k - k'|$, are always defined via their principal value, since the discrete version omits the points $k_i = k_j$ while including points $k_i = k_j \pm 1$, leading to a symmetric regularization. Similarly, the logarithm of the dimension of the irrep (2.15) becomes in the continuum limit

$$\ln \dim[\rho(k)] = -\ln \prod_{s=1}^{N-1} s! + \int_{0}^{\infty} dk \int_{0}^{k} dk' \, \rho(k) \rho(k') \ln(k-k') \,. \tag{3.7}$$

The distribution $\rho(k)$ satisfies the constraints

$$0 \le \rho(k) \le 1$$
, $\int_{0}^{\infty} dk \, \rho(k) = N$, $\int_{0}^{\infty} dk \, k \, \rho(k) = \frac{N^2}{2} + n$. (3.8)

The first constraint arises from the fermionic property $k_i \ge k_{i+1} + 1$, while the other two are the continuous, large-*N* version of (2.17).

In the large-n, N limit the statistics of irreps will be dominated by one particular distribution $\rho(k)$ corresponding to one irrep, the contribution from other irreps falling off exponentially as the density deviates from that distribution. The determination of the dominant irrep depends on the quantity of interest. In a pure mathematical context, the number of irreps $d_{n;k}$ could be the quantity of interest, and we would need to maximize it with respect to $\rho(k)$. That is, we would maximize the expression in (3.5) under the constraints (3.8). In statistical physics applications, on the other hand, the relevant quantity is the total number of states at given energy and other thermodynamic state variables. Assuming that the irrep determines all such variables (energy etc.), the relevant quantity is the total number of a given irrep times its dimensionality (number of states), that is,

$$m_{n;\mathbf{k}} = \dim(\mathbf{k}) d_{n;\mathbf{k}} = \frac{n! \, \Delta(\mathbf{k})^2}{\prod_{s=1}^{N-1} s! \prod_{i=1}^{N} k_i!} , \quad \text{with} \quad \sum_{i=1}^{N} k_i = n + \frac{N(N-1)}{2} ,$$
(3.9)

where we used ((2.14) and (2.15)). In terms of the (discrete) fermion density ρ_s we have that

$$m_{n;\rho} = \frac{n!}{\prod_{s=1}^{N-1} s!} \prod_{t>s=0}^{M} (t-s)^{(2\rho_s-1)\rho_t} . \tag{3.10}$$

The above obeys the formal duality invariance $\rho_s \to \tilde{\rho}_s = \frac{1}{2} - \rho_{M-s}$. However, this is not a true duality since, given that $\rho_s = 0$ or 1, $\tilde{\rho}_s$ takes the values $\pm \frac{1}{2}$. In the large-n, N limit in which $\rho_s \to \rho(k)$ we obtain from ((3.5) and (3.7)) that

$$\ln m_n[\rho(k)] = 2 \int_0^\infty dk \int_0^k dk' \, \rho(k) \rho(k') \ln(k - k') - \int_0^\infty dk \, \rho(k) \, k(\ln k - 1) \,, \tag{3.11}$$

up to a ρ -independent constant. This is identical in form to $\ln(d_n[\rho(k)]/n!)$ in (3.5), the only difference being the factor of 2 in front of the double integral.

We can thus consider the general form

$$S_{w,n}[\rho(k)] = \frac{w}{2} \int_{0}^{M} dk \int_{0}^{M} dk' \, \rho(k) \rho(k') \ln|k - k'| - \int_{0}^{M} dk \, \rho(k) \, k(\ln k - 1) \,, \tag{3.12}$$

where we ignored any overall ρ -independent constant and introduced an upper cutoff M for the k integrals, with the understanding that $\rho(k)=0$ for k>M. This reproduces the cases $\ln d_n[\rho]$ for w=1 and $\ln m_n[\rho(k)]$ for w=2, but can describe a more general situation. Note that the case of general w would correspond to starting, instead of (3.9), with

$$m_{w,n;\mathbf{k}} = \left(\dim(\mathbf{k})\right)^{w-1} d_{n;\mathbf{k}} = \frac{n! \left(\Delta(\mathbf{k})\right)^w}{\left(\prod_{s=1}^{N-1} s!\right)^{w-1} \prod_{i=1}^{N} k_i!} , \qquad \sum_{i=1}^{N} k_i = n + \frac{N(N-1)}{2} . \tag{3.13}$$

There is a clear distinction between the cases w > 1 and w < 1. The latter one is rather exotic, and perhaps unphysical, as it would correspond to a statistical model with entropy decreasing as the dimensionality of the irrep increases.

The quantity $S_{w,n}[\rho(k)]$ is invariant under the formal duality transformation

$$\rho(k) \to \tilde{\rho}(k) = w^{-1} - \rho(M - k)$$
 (3.14)

analogous to the one for the w=2 case (3.10). For $\tilde{\rho}$ to obey the fermionic constraint $0<\tilde{\rho}<1$, ρ must satisfy

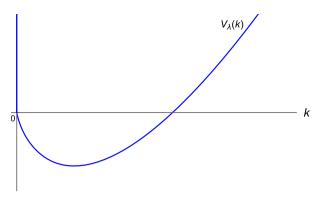


Fig. 1. The potential $V_{\lambda}(k)$ for a generic value of λ . It has a "rigid wall" at k=0 and forms a well for all values of λ , allowing for a distribution $\rho(k)$ localized inside the well, either touching k=0 or with support entirely at k>0.

$$w^{-1} - 1 < \rho < w^{-1}$$
, (3.15)

which implies that $\tilde{\rho}$ will also satisfy it. Further, $\tilde{\rho}$ satisfies the integral constraints (3.8) with modified parameters \tilde{N} and \tilde{n}

$$\tilde{N} = \frac{M}{w} - N$$
, $\tilde{n} = n + \frac{(w - 1)M(M - 2wN)}{2w^2}$. (3.16)

For the special value M = 2wN we see that $\tilde{N} = N$ and $\tilde{n} = n$ and therefore the transformation (3.14) becomes a self-duality. Note, however, that this duality holds only for densities satisfying (3.15) so it remains a restricted invariance. Its domain, in particular, does not include the singlet.

To calculate the dominant irrep, that is, the distribution $\rho(k)$ maximizing $S_{w,n}$, we maximize $S_{w,n}[\rho(k)]$ while enforcing the constrains (3.8) via two Lagrange multipliers. That is, we extremize

$$S_{w,n}[\rho(k)] - \mu \left(\int_{0}^{\infty} dk \, \rho(k) - N \right) - \lambda \left(\int_{0}^{\infty} dk \, k \, \rho(k) - n - \frac{N^2}{2} \right). \tag{3.17}$$

where from now on we take $M \to \infty$ keeping in mind that, in general, $\rho(k)$ will vanish outside a finite range. Using (3.5) and setting the functional derivative with respect to $\rho(k)$ to zero yields

$$w \int_{0}^{\infty} dk' \, \rho(k') \ln|k - k'| = k(\ln k - 1) + \mu + \lambda k \,. \tag{3.18}$$

Further differentiating (3.18) with respect to k we obtain

$$w \int_{0}^{\infty} dk' \frac{\rho(k')}{k - k'} = \ln k + \lambda . \tag{3.19}$$

The above equation must hold for k such that $\rho(k) \neq 0$ and $\rho(k) \neq 1$, since in empty regions with $\rho(k) = 0$ there are no k_i to vary, and in fully filled ones with $\rho(k) = 1$ the k_i cannot vary.

Hence, from (3.18) we see that the problem amounts to finding the equilibrium distribution of a large number of particles repelling each other with a logarithmic potential of strength w, inside an external potential given by the right hand side of (3.18), that is

$$V_{\lambda}(k) = k(\ln k - 1) + \lambda k$$
, (3.20)

where we have omitted the constant term μ . This potential is depicted in Fig. 1. It has a rigid "wall" at k=0 and goes to infinity as $k\to\infty$. Therefore, it may always support a finite lump of particles, spread around its minimum value $V_{\lambda,\min}=-e^{-\lambda}$ arising at $k=e^{-\lambda}$. The issue is whether this lump extends all the way to the boundary at k=0, and if it respects the condition $0\leqslant\rho(k)\leqslant1$. As we shall see, this depends on the values of N,n and w.

In the following sections, we solve the minimization problem and obtain the dominant $\rho(k)$. We will first treat the case w = 1 (maximizing $d_n[\rho(k)]$), since the solution simplifies and has some special properties, and then extend it to general w.

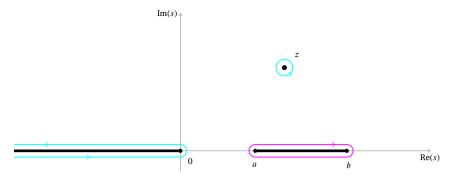


Fig. 2. Contour of integration in the *s*-plane. The original (magenta) contour around the square root cut on (a, b) is pulled back to the two (cyan) contours around the pole at z and the logarithm cut on $(-\infty, 0)$.

3.3. w = 1, maximal d_n

In this case the equation satisfied by $\rho(k)$ is simply (3.19) with w = 1

$$\int_{0}^{\infty} dk' \frac{\rho(k')}{k - k'} = \ln k + \lambda , \qquad (3.21)$$

together with the constraints (3.8). Moreover, for w=1 the duality relation (3.14) becomes exact, as it preserves the range of ρ and maps the singlet to itself. Depending on whether the inequality constraint $\rho \le 1$ is saturated in a finite domain, we distinguish two cases, corresponding to broken or unbroken duality symmetry.

3.3.1. Duality breaking phase $n > N^2/4$

We start by assuming that the distribution $\rho(k)$ does not reach the "wall" on the left at k = 0, i.e., it is nonzero inside an interval 0 < a < k < b and vanishes outside. Then solving (3.19) becomes a standard single-cut Cauchy problem. We define the resolvent

$$u(z) = \int dk \, \frac{\rho(k)}{z - k} \,, \tag{3.22}$$

with z on the upper complex plane. Its real and imaginary part on the real axis reproduce $\rho(k)$ and its Hilbert transform

$$u(k+i\epsilon) = \int dk' \frac{\rho(k')}{k-k'} - i\pi\rho(k) . \tag{3.23}$$

Therefore, a function that is analytic on the upper half plane and its real part on the real axis equals $\ln k + \lambda$ will equal u(z) up to an additive constant, and its imaginary part will fix $\rho(k)$. In standard fashion, we write

$$u(z) = \frac{1}{2\pi i} \sqrt{(z-a)(z-b)} \oint ds \, \frac{\ln s + \lambda}{(s-z)\sqrt{(s-a)(s-b)}} \,, \tag{3.24}$$

where the contour winds in the clockwise direction around the cut of the square root but does not include the singularity at z and the cut of the logarithm (see Fig. 2).

Pulling back the contour we pick up the pole at s = z and the integral around the cut of the logarithm

$$u(z) = \ln z + \lambda - \sqrt{(z-a)(z-b)} \int_{0}^{\infty} \frac{ds}{(s+z)\sqrt{(s+a)(s+b)}}$$

$$= \ln z + \lambda - i\cos^{-1}\frac{2z - a - b}{b - a} + i\cos^{-1}\frac{(a+b)z - 2ab}{(b-a)z}.$$
(3.25)

For z = k real and between a and b (the region in which $\rho(k)$ does not vanish) the last two terms are purely imaginary (the square root factor multiplying the integral provides a factor +i, since we assume that z approaches k from the upper-half complex plane). Then, according to (3.23), we determine $\rho(k)$ as

$$\rho(k) = \frac{1}{\pi} \cos^{-1} \frac{2k - a - b}{b - a} - \frac{1}{\pi} \cos^{-1} \frac{(a + b)k - 2ab}{(b - a)k}$$

$$= \frac{2}{\pi} \cos^{-1} \frac{\sqrt{k} + \sqrt{ab/k}}{\sqrt{a} + \sqrt{b}}, \quad a \le k \le b.$$
(3.26)

The last expression above makes clear that $\rho(k)$ indeed vanishes at k=a and k=b. It also makes clear that $\rho(k)$ is positive and never becomes larger than 1 (since the argument of \cos^{-1} never becomes negative and thus the angle does not exceed $\pi/2$), so $\rho(k)$ satisfies the constraint $0 \le \rho(k) \le 1$.

The parameters a, b and λ can be determined by matching the asymptotics of u(z)

$$u(z) = z^{-1} \int_{0}^{\infty} dk \, \rho(k) + z^{-2} \int_{0}^{\infty} dk \, k \, \rho(k) + \mathcal{O}(z^{-3})$$

$$= z^{-1} \, N + z^{-2} \left(n + \frac{N^2}{2} \right) + \mathcal{O}(z^{-3}) ,$$
(3.27)

where in the second line we used (3.8), with those implied from (3.25). We obtain

$$N = \frac{(\sqrt{b} - \sqrt{a})^2}{2} ,$$

$$n + \frac{N^2}{2} = \frac{2(b - a)^2 + (\sqrt{b} - \sqrt{a})^4}{16} ,$$

$$\lambda = -2 \ln \frac{\sqrt{a} + \sqrt{b}}{2} ,$$
(3.28)

from which the parameters a, b and λ are determined in terms of N and n as

$$\sqrt{a} = \sqrt{\frac{N}{2}} \left(\sqrt{\frac{1+t}{2}} - 1 \right),$$

$$\sqrt{b} = \sqrt{\frac{N}{2}} \left(\sqrt{\frac{1+t}{2}} + 1 \right),$$

$$\lambda = -\ln \frac{N}{4} - \ln \left(1 + t \right),$$
(3.29)

where we defined

$$t = 4\frac{n}{N^2} \,. \tag{3.30}$$

The above expressions make clear that, in order for both parameters n, N to remain relevant, their scaling must be $n \sim N^2$, identifying t in (3.30) as the relevant order parameter. Under that scaling, k, a, b, and $e^{-\lambda}$ all scale like $N \sim \sqrt{n}$. The form of $\rho(k)$ in (3.26) becomes, upon substituting (3.29),

$$\rho(k) = \frac{2}{\pi} \cos^{-1} \frac{\sqrt{\frac{k}{N}} + \frac{1}{4} \sqrt{\frac{N}{k}} (t - 1)}{\sqrt{t + 1}} , \quad a \le k \le b .$$
 (3.31)

It can be checked that the above density indeed obeys (3.8).

Since $\sqrt{a} > 0$, (3.29) shows that this solution will exist for $n > N^2/4$, that is, t > 1. For $n < N^2/4$ the above solution is not valid, and the point $n = N^2/4$ marks a transition. Clearly $\rho(k)$ in (3.31) does not satisfy the self-duality condition $\rho(k) = 1 - \rho(2N - k)$, so the phase $n > N^2/4$ is a duality breaking one.

3.3.2. Duality preserving phase $n < N^2/4$

For $n = N^2/4$ (t = 1) the parameter a is driven to zero and the solution (3.26) of the previous section becomes

$$\rho(k) = \frac{1}{\pi} \cos^{-1} \frac{2k - b}{b} = \frac{2}{\pi} \cos^{-1} \sqrt{\frac{k}{b}}, \qquad 0 < k < b.$$
(3.32)

The parameters b and λ in this case can be obtained from the $a \to 0$ limit of the corresponding expressions (3.29) as

$$b = 4\sqrt{n} , \qquad \lambda = -\frac{1}{2} \ln n . \tag{3.33}$$

We note that, now, $\rho(0) = 1$. This marks a transition to a phase where the density $\rho(k)$ saturates to 1 over a finite interval when $n < N^2/4$.

In fact, in this phase the expression for the maximal $d_n[\rho(k)]$ develops a flat region for a range $0 \le k \le a$ with the rest of it becoming N-independent, for all $n < N^2/4$. This is already clear in the case of finite N > n: the expression (2.14) for $d_{n,k}$ also holds if N is replaced by N' > N. Specifically, define the new "extended" momenta k'_i with i = 1, 2, ..., N'

$$k'_{i} = \begin{cases} k_{i} + N' - N, & i = 1, \dots, N \\ N' - i, & i = N + 1, \dots, N' \end{cases}$$
(3.34)

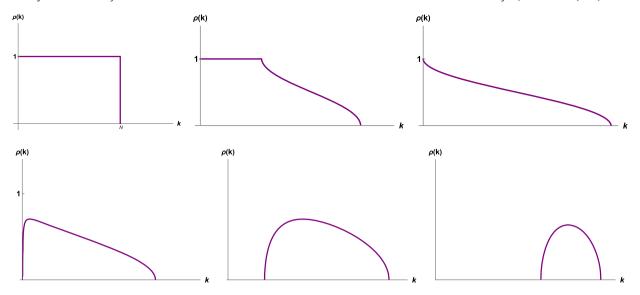


Fig. 3. The distribution ρ(k) for various values of n/N^2 . For n = 0 (first panel) the distribution is a step function corresponding to the singlet. For $0 < n < N^2/4$ (second panel) the edge of the distribution deforms into an inverse cosine. For $n = N^2/4$ (third panel) the deformation reaches k = 0, signaling a phase transition. As soon as n exceeds $N^2/4$ (fourth panel) the left edge of the distribution drops to ρ(0) = 0, and as n increases (fifth panel) ρ(x) has support on a positive interval. For $n \gg N^2/4$ (sixth panel) it approaches a Wigner semicircle distribution.

identical to the old ones but shifted to the right by N' - N with the "Fermi sea" filled to their left (the second line above). Then we have

$$d_{n,\mathbf{k}} = d_{n,\mathbf{k}'}. ag{3.35}$$

This can be shown, e.g., inductively. For N' = N + 1,

$$\frac{\prod_{j>i=1}^{N'}(k_i'-k_j')}{\prod_{i=1}^{N'}k_i'!} = \frac{\prod_{j>i=1}^{N}(k_i-k_j)\prod_{i=1}^{N}(k_i+1)}{\prod_{i=1}^{N}(k_i+1)!} = \frac{\prod_{j>i=1}^{N}(k_i-k_j)}{\prod_{i=1}^{N}k_i!}$$
(3.36)

and by induction we can reach any value N' > N. It is clear that

$$\sum_{i=1}^{N'} k'_i = n + \frac{N'(N'-1)}{2} \,, \tag{3.37}$$

where we used (2.14) for k. Therefore, $\{k_i'\}$ represents the same irrep as $\{k_i\}$ (same Young tableau) arising from the direct product of n fundamentals of SU(N'). This translates to a corresponding relation in the momentum density description. It can be checked that the new "extended" density

$$\rho'(k) = \begin{cases} 1, & 0 < k < N' - N, \\ \rho(k + N - N'), & k > N' - N, \end{cases}$$
(3.38)

produces the same $d_n[\rho(k)]$ in (3.5).

This provides an extension of the solution (3.32), which was ostensibly valid only for $n = N^2/4$, to any $n > N^2/4$. Specifically, applying (3.38) with $N \to 2\sqrt{a}$ and $N' \to N$ for the configuration (3.32), and taking also into account that $b = 4\sqrt{n}$, we obtain

$$\rho(k) = \begin{cases} 1, & 0 < k < N - 2\sqrt{n}, \\ \frac{1}{\pi} \cos^{-1} \frac{k - N}{2\sqrt{n}}, & N - 2\sqrt{n} < k < N + 2\sqrt{n} \end{cases}$$
(3.39)

and zero elsewhere. It can be checked that the above density indeed satisfies (3.8). Moreover, it is self-dual, satisfying

$$\rho(k) = 1 - \rho(2N - k) , \qquad 0 < k < 2N . \tag{3.40}$$

Therefore, $n < N^2/4$ represents a duality preserving phase. An evolution of $\rho(k)$ for various values of n/N^2 is presented in Fig. 3.

We conclude by pointing out that the distribution (3.39) reproduces the VKLS limiting Young tableau shape for a large number of boxes n weighted by the Plancherel measure [25,26]. The VKLS parametrization of Young tableau shapes is achieved by reflecting the Young tableau about its top row, rotating it by $\pi/4$ in the positive direction to produce a V-based shape, and rescaling it by a factor of $1/\sqrt{n}$. The coordinates (x, y) of the last box of row ℓ_i , for large n, will be

$$\ell_i = \sqrt{\frac{n}{2}}(x+y), \qquad i = \sqrt{\frac{n}{2}}(y-x)$$
 (3.41)

Going over to our variables k_i , we recall relation (2.7) between ℓ_i and k_i . In fact, we will slightly modify this relation to $k_i = \ell_i + N - i$, so that ℓ_N is not necessarily 0. This has the advantage that the total number of boxes in the tableau is n, at the price of potentially introducing columns of length N in the tableau, although in the duality preserving phase this will never happen since $k_N = 0$. Then, in the continuum limit $N, n \gg 1$, (3.41) implies

$$k = N + \sqrt{2n} x \tag{3.42}$$

The corresponding density of k_i will be

$$\rho(k) = -\frac{di}{dk} = \frac{1}{2} \left(1 - \frac{dy}{dx} \right) . \tag{3.43}$$

Upon substituting (3.39) and integrating, we obtain

$$y = \int \left(1 - 2\rho(N + \sqrt{2nx})\right) dx = \begin{cases} \frac{2}{\pi} \left(\sqrt{2 - x^2} + x \sin^{-1} \frac{x}{\sqrt{2}}\right), & |x| \le \sqrt{2}, \\ |x|, & -\frac{N}{\sqrt{2n}} \le x \le -\sqrt{2} & \text{or} & x \ge \sqrt{2}, \end{cases}$$
(3.44)

where the integration constant was fixed by the condition that k = 0 for i = N. This is the VKLS distribution, with the left branch of |x| truncated at $x = -N/\sqrt{2n}$, and is valid for all $n < N^2/4$. The relation of our work with the Plancherel process and related distributions will be analyzed in an upcoming publication.

3.3.3. The Wigner semicircle limit $1 \ll N \ll \sqrt{n}$

In the limit $n \gg N^2/4 \gg 1$ ($t \gg 1$), deep in the duality breaking region, the particle momenta k_i lump near the minimum of the effective potential in (3.20), where it can be approximated as a harmonic oscillator. The minimum arises at

$$k_0 = e^{-\lambda} = \frac{n}{N} + \frac{N}{4} \simeq \frac{n}{N} = \frac{N}{4}t$$
(3.45)

and the potential can be approximated by

$$V(k) \simeq -\frac{n}{N} + \frac{N}{2n} \left(k - \frac{n}{N} \right)^2 . \tag{3.46}$$

As a consequence, the distribution $\rho(k)$ will become a Wigner semicircle around k_0 with radius $\sqrt{2n} = N\sqrt{t/2}$

$$\rho(k) = \frac{N}{\pi n} \sqrt{2n - \left(k - \frac{n}{N}\right)^2} = \frac{2}{\pi} \sqrt{\frac{2}{t}} \sqrt{1 - 2\frac{(k/N - t/4)^2}{t}} \ . \tag{3.47}$$

The same result is obtained by taking a limit of $\rho(k)$ in (3.26): we set $k/N = t/4 + \sqrt{t} y$, where the form of the last term is dictated by the fact that k has range $b - a \simeq N\sqrt{2t}$ for $t \gg 1$. Expanding $\rho(k)$ for large t we obtain to leading order

$$\rho(k) \simeq \frac{2}{\pi} \cos^{-1} \left(1 - \frac{1 - 2y^2}{t} \right) \simeq \frac{2}{\pi} \sqrt{\frac{2}{t}} \sqrt{1 - 2y^2} \ . \tag{3.48}$$

Reinstating the variable k/N we recover (3.47).

Interestingly, this result also derives from the $n\gg N\sim 1$ distribution (3.1). The exponent of the gaussian exponential corresponds to k_i having a potential $(N/2n)k_i^2$. which is the same potential as (3.46). The Vandermonde factor in front endows the k_i with a logarithmic mutual repulsive potential, while the delta-function puts the average momentum $(1/N)\sum_i k_i$ to the value n/N, leading to the Wigner semicircle distribution (3.47). So the $N\sim 1$ result reliably reproduces the $1\ll N\ll 2\sqrt{n}$ situation, but not the one for $N\sim 2\sqrt{n}$, which leads to distorted Wigner semicircle distributions and, eventually, to the phase transition to the self-dual phase.

3.3.4. Phase transition

The point $n = N^2/4$, or t = 1, clearly marks a phase transition. To identify the order of the phase transition and the properties of the two phases we calculate the maximal d_n for the dominant $\rho(k)$ in each phase. We will compute its first few derivatives with respect the order parameter t and will discover a discontinuity in its third derivative at t = 1, identifying it as a third-order transition.

We will use the standard result that the derivative with respect to n of any functional $F[\rho(k)]$ of $\rho(k)$ that does not explicitly depend on n at its maximum in $\rho(k)$ subject to the constraints (3.8) is given by the Lagrange multiplier λ for n. Indeed,

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$$\frac{\partial}{\partial n} F[\rho(k)] = \int_{0}^{\infty} dk \, \frac{\delta F[\rho(k)]}{\delta \rho(k)} \, \frac{\partial \rho}{\partial n} = \int_{0}^{\infty} dk \, (\mu + \lambda k) \frac{\partial \rho}{\partial n}$$

$$= \mu \frac{\partial}{\partial n} \int_{0}^{\infty} dk \, \rho + \lambda \frac{\partial}{\partial n} \int_{0}^{\infty} dk \, k \rho = \mu \frac{\partial N}{\partial n} + \lambda \frac{\partial}{\partial n} \left(n + \frac{N^{2}}{2} \right) = \lambda,$$
(3.49)

where we used the saddle point condition and the constraints. The quantity $\ln(d_n[\rho(k)]/n!)$ that was maximized with respect to $\rho(k)$ indeed does not involve n explicitly, and thus

$$\frac{\partial}{\partial n} \ln \frac{d_n[\rho(k)]_{\text{max}}}{n!} = \lambda , \qquad (3.50)$$

or, in terms of the order parameter $t = 4n/N^2$.

$$\frac{\partial}{\partial t} \ln \frac{d_n[\rho(k)]_{\text{max}}}{n!} = \frac{N^2}{4} \lambda . \tag{3.51}$$

where we have explicitly indicated that it is the density at the maximal configuration.

Before proceeding to compute further derivatives, we may integrate (3.50) above to find $d_n[\rho(k)]_{\text{max}}$ itself. In the duality preserving phase $n < N^2/4$ (t < 1), for which $\lambda = -\frac{1}{2} \ln n$ from (3.33), we obtain

$$\ln \frac{d_n[\rho(k)]_{\text{max}}}{n!} = -\frac{N^2}{8}t\left(\ln t - 1 + 2\ln\frac{N}{2}\right) = -\frac{n}{2}(\ln n - 1) \simeq -\frac{1}{2}\ln n!,$$
(3.52)

where the integration constant has been fixed such that in the limit $n \to 0$, in which we are left with the fundamental irrep, $d_n \to 1$. Therefore,

duality preserving phase:
$$d_{n \text{ max}} = d_n[\rho(k)]_{\text{max}} \simeq \sqrt{n!}$$
. (3.53)

Similarly, for the duality breaking phase, λ is given by (3.29), so we find

$$\ln \frac{d_n[\rho(k)]_{\text{max}}}{n!} = n - \left(n + \frac{N^2}{4}\right) \ln \left(\frac{n}{N} + \frac{N}{4}\right) - \frac{N^2}{4} \left(\frac{1}{2} - \ln \frac{N}{2}\right)$$
(3.54)

where we fixed the integration constant by matching the result at $n = N^2/4$ with the one in the duality preserving phase, since $\rho(k)$ has no discontinuity at the transition point. Thus

duality breaking phase:
$$d_{n,max} = d_n[\rho_n(k)] = N^n \frac{n! \left(\frac{N^2}{4}\right)!}{\left(n + \frac{N^2}{4}\right)!} \left(2e^{-1/2}\right)^{N^2/4}$$
 (3.55)

The above was written in a suggestive form, one of several equivalent forms at the $N, n \gg 1$ limit, to display the leading behavior: the multiplicity of the dominant irrep becomes a fraction of the total number of states N^n . As n increases, this fraction becomes

$$\frac{d_{n,max}}{N^n} \simeq \left(\frac{N^2}{4}\right)! \left(\frac{2}{ne^{1/2}}\right)^{N^2/4}, \quad n \gg N^2/4.$$
 (3.56)

It is clear that $\ln d_n/n!$, playing the role of free energy, has no discontinuity at the transition point, and neither does its first derivative, since λ is continuous across the transition. It turns out that the second derivative in t for fixed N evaluated at t=1 is also continuous. Specifically,

$$\frac{\partial}{\partial t} \ln \frac{d_n[\rho(k)]_{\text{max}}}{n!} \Big|_{t \to 1^-} = -\frac{N^2}{4} \ln \frac{N}{2} = \frac{\partial}{\partial t} \ln \frac{d_n[\rho(k)]_{\text{max}}}{n!} \Big|_{t \to 1^+},
\frac{\partial^2}{\partial t^2} \ln \frac{d_n[\rho(k)]_{\text{max}}}{n!} \Big|_{t \to 1^-} = -\frac{N^2}{8} = \frac{\partial^2}{\partial t^2} \ln \frac{d_n[\rho(k)]_{\text{max}}}{n!} \Big|_{t \to 1^+},$$
(3.57)

However, the third derivative is discontinuous:

$$\frac{\partial^{3}}{\partial t^{3}} \ln \frac{d_{n} [\rho(k)]_{\text{max}}}{n!} \Big|_{t \to 1^{-}} = \frac{N^{2}}{8} ,
\frac{\partial^{3}}{\partial t^{3}} \ln \frac{d_{n} [\rho(k)]_{\text{max}}}{n!} \Big|_{t \to 1^{+}} = \frac{N^{2}}{16} .$$
(3.58)

Therefore, this is a 3rd-order phase transition.

The picture that emerges in terms of fermion momenta is that, for large N, the singlet irrep corresponds to a filled Fermi sea with Fermi level k = N. Multiplying with fundamental irreps excites the state by one unit of momentum per irrep and results in excitations around the Fermi level. As long as $n < N^2/4$ the excitations remain localized around the Fermi level and are N-independent. For

 $n = N^2/4$ the excitations reach the bottom of the sea (k = 0), marking a phase transition, and for $n > N^2/4$ the entire Fermi sea is excited and lifted above k = 0. We remark that, in the case $1 \ll 2\sqrt{n} < N < n$, there are in principle irreps with all the k_i excited above their ground state (singlet) values, but such irreps have subleading multiplicities and are irrelevant in the large-N limit.

3.4. General repulsion $w \neq 1$, maximal $S_{w,n}$

For general w, the equation for $\rho(k)$ that maximizes $S_{w,n}[\rho(k)]$ is (3.19) with the constraints (3.8). The solution proceeds along similar lines as the w=1 case. Rather than "duality preserving" and "duality violating" phases, we will talk about "condensed" and "dilute" cases, the former being one where the density reaches its saturation value $\rho(k)=1$ for a range of values of k, the latter one with $\rho(k)$ always less than 1. As we shall see, for w>1 the condensed phase always involves a saturation region [0,a] for k, while for w<1 it can saturate in a region $[a_1,a_2]$ with $0< a_1< a_2$. We shall focus on the physically more relevant case $w\geqslant 1$ from now on, for reasons explained below (3.13).

3.4.1. Dilute phase $n > n_w$

In this case $\rho(k) < 1$ for all k and the solution can be obtained from the w = 1 solution with a simple rescaling. Specifically,

$$\bar{\rho}(k) = w\rho(k) , \quad \overline{N} = wN , \quad \bar{n} = wn + \frac{w(1-w)N^2}{2} ,$$
 (3.59)

satisfy the same equations and constraints for w = 1. We can use the solution in that case to find

$$\rho(k) = \frac{2}{u\pi} \cos^{-1} \frac{\sqrt{k} + \sqrt{ab/k}}{\sqrt{a} + \sqrt{b}}, \qquad a \leqslant k \leqslant b,$$
(3.60)

with

$$\sqrt{a} = \sqrt{\frac{(2-w)N}{4} + \frac{n}{N}} - \sqrt{\frac{wN}{2}},$$

$$\sqrt{b} = \sqrt{\frac{(2-w)N}{4} + \frac{n}{N}} + \sqrt{\frac{wN}{2}},$$

$$\lambda = -\ln\left(\frac{(2-w)N}{4} + \frac{n}{N}\right).$$
(3.61)

Clearly $\rho(k) < w^{-1}$ and thus $\rho(k) \le 1$ for $w \ge 1$. The above solution will exist as long as $\sqrt{a} \ge 0$, and thus for n above a critical value n_w

$$n > n_w = \frac{(3w - 2)N^2}{4} \implies t > 3w - 2,$$
 (3.62)

for the order parameter t defined in (3.30). For w = 1 we recoved the transition at $n = N^2/4$, while for the number of states case w = 2 the transition happens at $n = N^2$.

3.4.2. Condensed phase $n < n_w$

For $n < n_w$ (t < 3w - 2) the solution (3.60) ceases to exist and we enter the condensation phase. Similar to the solution (3.39), we set

$$\rho(k) = \begin{cases} 1, & 0 < k < a, \\ \rho_0(k-a), & a < k < a+b, \end{cases}$$
 (3.63)

and zero elsewhere, with a,b two positive constants. The function $\rho_0(k)$ satisfies the constraints

$$0 \le \rho_0(k) \le 1 \;, \quad \int_0^b dk \, \rho_0(k) = N - a \;, \quad \int_0^b dk \, k \, \rho_0(k) = \frac{(N - a)^2}{2} + n \;. \tag{3.64}$$

Substitution of $\rho(k)$ in (3.12) yields, upon changing variable $k \to k + a$,

$$\begin{split} S_{w,n}[\rho_0(k)] &= \frac{w}{2} \int\limits_0^\infty dk \int\limits_0^\infty dk' \, \rho_0(k) \rho_0(k') \ln |k-k'| \\ &- w \int\limits_0^\infty dk \, \rho_0(k) \, k(\ln k - 1) + (w - 1) \int\limits_0^\infty dk \, \rho_0(k) (k + a) \Big(\ln (k + a) - 1 \Big) \;, \end{split} \tag{3.65}$$

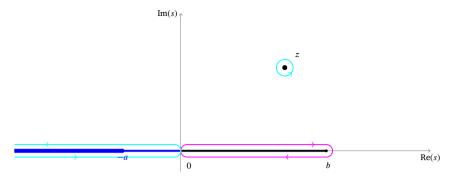


Fig. 4. Contour of integration in the *s* plane. The original (magenta) contour around the (black) square root cut on (0,b) is pulled back to the two (cyan) contours around the pole at *z* and the two (blue) logarithm cuts on $(-\infty,0)$ and $(-\infty,-a)$.

where we omitted terms set to be constant by the constraints and harmlessly extended the integration range to infinity since $\rho_0(k)$ has finite range. Adding to this the constraints (3.64) with appropriate Lagrange multipliers and following the procure that led to (3.18) we obtain the equilibrium equation for ρ_0

$$w \int_{0}^{\infty} dk' \, \rho_0(k') \ln|k - k'| = wk(\ln k - 1) - (w - 1)(k + a) \left(\ln(k + a) - 1 \right) + \mu + \lambda k \,. \tag{3.66}$$

Taking the k-derivative we obtain the analog of (3.19), i.e.

$$\int_{0}^{\infty} dk' \frac{\rho_0(k')}{k - k'} = \ln k - \gamma \ln(k + a) + \lambda/w , \quad \text{when } \rho_0(k) > 0$$
 (3.67)

and where we defined for convenience the parameter²

$$\gamma = 1 - \frac{1}{w}, \quad \gamma \in [0, 1].$$
 (3.68)

We see that the equation for $\rho_0(k)$ now has a two-logarithm potential, given by the right hand side of (3.66) as

$$V_{\lambda w a}(k) = wk(\ln k - 1) - (w - 1)(k + a)(\ln(k + a) - 1) + \lambda k.$$
(3.69)

(up to a constant term μ), while for w = 1 the second logarithm drops out.

To solve for $\rho_0(k)$, we define as before the resolvent

$$u_0(z) = \int dk \frac{\rho_0(k)}{z - k},$$
 (3.70)

reproducing $\rho_0(k)$ and its Hilbert transform as

$$u_0(k+i\epsilon) = \int dk' \frac{\rho_0(k')}{k-k'} - i\pi \rho_0(k) . \tag{3.71}$$

In analogy to (3.24), we set

$$u_0(z) = \frac{1}{2\pi i} \sqrt{z(z-b)} \oint ds \, \frac{\ln s - \gamma \ln(s+a) + \lambda/w}{(s-z)\sqrt{s(s-b)}} \,, \tag{3.72}$$

where the contour winds in the clockwise direction around the cut of the square root [0,b] but does not include the singularity at z nor the cuts of the logarithms (so it "threads" the real line at k=0) (see Fig. 4). Pulling back the contour we pick up the pole at s=z and the integral around the cuts of the logarithms and obtain

$$u_0(z) = \ln z - \gamma \ln(z+a) + \frac{\lambda}{w} - \sqrt{z(z-b)} \left[\int\limits_0^\infty -\gamma \int\limits_a^\infty \right] \frac{ds}{(s+z)\sqrt{s(s+b)}} \ . \tag{3.73}$$

For z = k real and between 0 and b (the region in which $\rho_0(k)$ does not vanish) the last term proportional to $\sqrt{z(z-b)}$ is purely imaginary and, according to (3.71), determines $\rho_0(k)$ as

² A similar equation to (3.67) was found in [27] (eq. (5.2) of that work) in a matrix model approach to black hole microstates.

$$\rho_0(k) = \sqrt{k(b-k)} \left[(1-\gamma) \int_0^\infty + \gamma \int_0^a \frac{ds}{(s+k)\sqrt{s(s+b)}} \right]. \tag{3.74}$$

Upon performing the integrals,

$$\rho_0(k) = \frac{2}{w\pi} \cos^{-1} \sqrt{\frac{k}{b}} + \frac{2(w-1)}{w\pi} \cos^{-1} \sqrt{\frac{(a+b)k}{(a+k)b}} \,. \tag{3.75}$$

The density $\rho_0(k)$ is a decreasing function of k, with $\rho_0(0) = 1$ and $\rho_0(b) = 0$. For w = 1 this recovers the result (3.32), while for the number of states case w = 2 it reduces to

$$\rho_0(k) = \frac{1}{\pi} \cos^{-1} \sqrt{\frac{k}{b}} + \frac{1}{\pi} \cos^{-1} \sqrt{\frac{(a+b)k}{(a+k)b}} = \frac{1}{\pi} \cos^{-1} \frac{(1+\sqrt{1+b/a})k - b}{b\sqrt{1+k/a}} \,. \tag{3.76}$$

The parameters a, b and λ can be related to N and n by matching the asymptotic expansion of $u_0(z)$

$$\begin{split} u_0(z) &= z^{-1} \int\limits_0^\infty dk \, \rho_0(k) + z^{-2} \int\limits_0^\infty dk \, k \, \rho_0(k) + \mathcal{O}\!\left(z^{-3}\right) \\ &= z^{-1} \left(N - a\right) + z^{-2} \left(n + \frac{(N - a)^2}{2}\right) + \mathcal{O}\!\left(z^{-3}\right) \,, \end{split} \tag{3.77}$$

from (3.70) using (3.64) to those from (3.72). We obtain

$$\lambda = 2(w-1)\ln\frac{\sqrt{a} + \sqrt{a+b}}{2} - w\ln\frac{b}{4},$$

$$N - a = \frac{b}{2} - \frac{\gamma}{2}\left(\sqrt{a+b} - \sqrt{a}\right)^{2},$$

$$\frac{(N-a)^{2}}{2} + n = \frac{3b^{2}}{16} + \frac{\gamma}{4}\left(-\frac{3b^{2}}{4} + 2a^{2} + (b-2a)\sqrt{a(a+b)}\right).$$
(3.78)

Although the system (3.78) is quite complicated, it can be explicitly solved for a, b. We set

$$a = N\left(\sqrt{\left(\frac{w}{2} - 1\right)x + 1} - \sqrt{\frac{wx}{2}}\right)^{2},$$

$$b = 2N\sqrt{wx\left((w - 2)x + 2\right)}.$$

$$(3.79)$$

These satisfy the last two of (3.78) provided that

$$x = \frac{4n/N^2}{w + \sqrt{w^2 + 4(w - 2)n/N^2}}.$$
(3.80)

The above give for λ

$$\lambda = -\ln\frac{N}{2} - \frac{w}{2}\ln\left(\frac{4wn}{N^2}\right) + \left(\frac{w}{2} - 1\right)\ln\left(2 - w + \sqrt{w^2 + (w - 2)4n/N^2}\right) + \frac{w}{2}\ln\left(w + \sqrt{w^2 + (w - 2)4n/N^2}\right). \tag{3.81}$$

Note that the argument of the square roots are strictly positive for $n < n_w$.

As a consistency check, for w = 1 the above reproduce the results (3.28) (for a = 0), while for $n = n_w = (3w - 2)/4N^2$, x = 1 and thus a = 0, b = 2wN and the results match the results of the dense case at the critical point. Also, n = 0 implies x = 0, and thus a = N, b = 0, reproducing the singlet distribution.

For w = 2 the results simplify considerably and we obtain

$$w = 2: \qquad a = \left(\sqrt{N} - \sqrt{\frac{n}{N}}\right)^2, \qquad b = 4\sqrt{n}, \qquad \lambda = -\ln\frac{n}{N}. \tag{3.82}$$

Remarkably, these are just the dilute case results (3.61) for w = 2 (noting that b in that case maps to a + b in the present case), analytically continued for negative values of $n - N^2$.

The behavior of $\rho(k)$ for w>1 is qualitatively similar as for w=1, with the notable exception that the maximal value of $\rho(k)$ in the dilute phase is $\rho=1/w$, achieved for $k\simeq 0$ as $n\to n_w+\epsilon$. However, $\rho(0)$ jumps from $\rho(0)=0$ for $n=n_w+\epsilon$ to $\rho(0)=1$ for $n=n_w-\epsilon$. This behavior is displayed in Fig. 5.

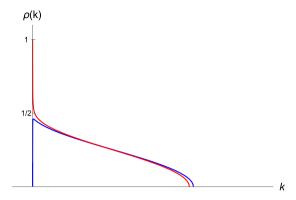


Fig. 5. The distribution $\rho(k)$ for w=2 around the critical point $n=N^2$. Both the dilute (blue) density and the dense (red) density have a value close to $w^{-1}=1/2$ near k=0, but the dilute one sharply dips to 0 and the dense one sharply rises to 1, for an increasingly sharp transition as n crosses the critical value N^2 .

Table 1 Phase transitions for various values of $w \ge 1$.

transition	3rd order	4th order	no transition
$w = 1$ $w > 1 (w \neq 2)$	✓ (crossover)	√	
w = 2			✓

3.4.3. Phase transition

The general w case exhibits an interesting pattern of phase transitions, depending on the value of w. As in the w=1 case, the "free energy" functional $S_w[\rho(k)]$ appearing in (3.12) does not depend explicitly on n, so its derivative with respect to n at the dominant configuration is still given by the value of the Lagrange multiplier λ . We have explicit expressions for λ in both phases, dilute (3.61) and condensed (3.81), so we may directly calculate its derivatives on either side of the critical point $n=n_w$. Equivalently, we give its expansion around the critical point. For the condensed phase $n < n_w$ the expansion is

$$n < n_w: \qquad \lambda = -\ln \frac{wN}{2} - \frac{2}{wN^2} (n - n_w) + \frac{2}{w^2 N^4} (n - n_w)^2 - \frac{8 - 14w + 7w^2}{3w^3 (1 - w)^2 N^6} (n - n_w)^3 + \dots$$
(3.83)

while for the dilute phase $n > n_{uv}$ we obtain

$$n > n_w: \qquad \lambda = -\ln \frac{wN}{2} - \frac{2}{wN^2} (n - n_w) + \frac{2}{w^2 N^4} (n - n_w)^2 - \frac{8}{3w^3 N^6} (n - n_w)^3 + \dots$$
(3.84)

We see that for all values of w > 1, except w = 2, λ and its first two derivatives in n (equivalently in t) are continuous, while its third derivative is discontinuous at $n = n_w$. This means that the first three derivatives of S_w are continuous but the *fourth* one is discontinuous, signaling a *fourth-order* phase transition at $n = n_w$.

The values w = 2 and w = 1 are special. For w = 2, all derivatives of λ are continuous across $n = n_2 = N^2$, since λ in this case is given by a unique analytic function of n (3.82), so there is *no* phase transition.

The case w=1 is trickier. From (3.83) and (3.84) we would deduce that the second derivative of λ is continuous, in contradiction to the results of section 3.3.4. What in fact happens is that for $w=1+\epsilon$, the second derivative $d^2\lambda/dn^2$ in the condensed phase displays an increasingly sharp transition: it approaches $8/N^4$ as n approaches $n_1=N^2/4$, which is the value for w=1, but at $n \sim N^2/4 - \epsilon^2$ it starts sharply dropping, and as n exceeds $N^2/4$ it reaches $4/N^4$, the value consistent with (3.83) (see Fig. 6). Physically, for w close to 1 the system goes through a crossover near $n=n_w$, which is increasingly sharp as w approaches 1 and becomes a full phase transition at w=1. Mathematically, the limits $w\to 1$ and $n\to n_w$ in the condensed phase do not commute. By contrast, in the dilute phase the $w\to 1$ limit is smooth. The various phase transitions are summarized in Table 1.

As in the w=1 case, we can integrate λ with respect to n to find the quantity $m_{w,n}$ for general w. Note that the relation of $m_{w,n}$ of (3.13) with $S_{w,n}$ of (3.12) is

 $^{^{3}}$ In [27] a third order phase transition was mentioned with an order parameter related to w. Its relationship to our present work is not obvious.

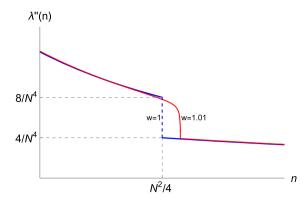


Fig. 6. Plots of $\lambda''(n) = d^2 \lambda/dn^2$ for w = 1 (blue curve) and w = 1.01 (red curve) around the w = 1 critical point $n = N^2/4$. For w = 1.01 there is a sharp transition from $\lambda'' \simeq 4/N^4$ to $\lambda'' \simeq 8/N^4$ but no discontinuity, while for w = 1 the transition evolves into a discontinuity. The w = 1.01 curve has a cusp at its critical point, signifying a discontinuous 3^{rd} derivative and a fourth-order phase transition.

$$\ln m_{w,n} = S_{w,n} + \ln n! - (w - 1) \sum_{s=1}^{N-1} \ln s!$$
(3.85)

and is defined such that for the singlet irrep it becomes 1. This fixes the integration constant in the dense case, and continuity at $n = n_w$ fixes the integration constant in the dilute case. We obtain for the dense phase $n < n_w$

$$\ln m_{w,n;\max} = \frac{2n(w-1)}{w + \sqrt{\frac{4n(w-2)}{N^2} + w^2}} - \frac{n}{2} \left((w-2) \ln n + w \ln w + 1 \right)$$

$$+ \frac{1}{2} \left(w \left(n + N^2 \right) - 2n - N^2 \right) \ln \left(\sqrt{\frac{4n(w-2)}{N^2} + w^2} - w + 2 \right)$$

$$+ \frac{nw}{2} \ln \left(\sqrt{\frac{4n(w-2)}{N^2} + w^2} + w \right) + (w-1) \left(n \ln \frac{N}{2} - \frac{N^2}{2} \ln 2 \right),$$
(3.86)

which indeed vanishes for n = 0. For the dilute phase $n > n_w$ we have that

$$\ln m_{w,n;\max} = n \ln n - N \left(\frac{(2-w)N}{4} + \frac{n}{N} \right) \ln \left(\frac{(2-w)N}{4} + \frac{n}{N} \right) + \frac{N^2}{8} \left(4 \ln N + w - 2 + 2w \ln \frac{w}{2N} \right).$$
(3.87)

In the special case w = 2 the expressions (3.86) and (3.87) simplify and unify, giving the result

$$m_{2\,n\text{max}} = N^n \,, \tag{3.88}$$

as expected, since the saddle point in the large-N, n limit must give the full number of states N^n up to subleading terms (determinants).

We conclude with a couple of remarks. In the case 0 < w < 1, $\rho(k)$ in (3.60) can reach or exceed the value 1 for n below a critical value *higher* than n_w , or equivalently a critical value for t

$$t < t_c = \frac{2w}{\sin^2 \frac{\pi w}{2}} + w - 2. \tag{3.89}$$

For $t < t_c$, or $n < t_c N^2/4$, the solution (3.60) is not valid any more, as it exceeds 1 at some interval. The true solution is one with

$$\rho(k) = \begin{cases} 0, & 0 < k < a_1 \text{ or } a_2 < k, \\ \rho_1(k), & a_1 < k < b_1, \\ 1, & b_1 < k < b_2, \\ \rho_2(k), & b_2 < k < a_2. \end{cases}$$

$$(3.90)$$

leading to a genuine two-cut Cauchy problem. We will not explore this solution, as w < 1 would correspond to a statistical model with entropy decreasing as the dimensionality of the irrep increases (see (3.13)). This is rather unphysical, although it could conceivably find application in some exotic situation.

Finally, the case w < 0 is even more unphysical, representing a drastic reduction of entropy. Its large-(N, n) limit would correspond to N particles *attracting* each other with two-body logarithmic potentials. In the concave external potential $V_{\lambda}(k)$ of (3.20) the only stable configuration is one with all particles coalescing to an interval of length N, for a density of 1, corresponding to the singlet irrep.

4. Conclusions

We considered the multiplicity of irreps arising in the decomposition of n fundamental representations of SU(N), weighted by a power of their dimension. We showed that a nontrivial double scaling limit exists in which both n and N become large keeping the ratio n/N^2 fixed, and uncovered novel phase transitions in which this ratio plays the role of the order parameter. The system generically undergoes a fourth order phase transition, from a dense to a dilute phase, enhanced to a third order one for the unweighted multiplicity, and ceasing to exist altogether when weighting with the first power of the dimension, which corresponds to the infinite temperature partition function of nonabelian ferromagnets.

Our results are model independent, not involving a Hamiltonian, and should thus be relevant to the thermodynamics of nonabelian ferromagnets at high temperatures. In this respect, it is interesting to reconsider the phase structure of the ferromagnetic model in which n atoms mutually interact via SU(N) components, which we recently investigated in [22] in the thermodynamic limit $n \gg 1$ but for fixed finite N. We expect that the present double scaling limit will qualitatively modify the phase structure, provided $N \sim \sqrt{n}$. The generalization of our results for a product of irreps other than the fundamental would also be interesting, the adjoint being the most natural alternative choice. The related question of the relation of our results to Markov processes in the space of tableaux, such as the Plancherel process, is also of mathematical interest.

Finally, the relevance of our results to matrix models and large-N Yang-Mills theories should be explored. Of particular interest is the understanding of microstates in the two-dimensional black hole of [28,29] with matrix models, along the lines of [30,31] and more recently of [27,32], and of the deconfinement/Hagedorn transition in large-N gauge theories [33–35], especially in the setting of [36,37]. These and other related questions are the subject of ongoing investigation.

Declaration of competing interest

The authors declare that they have no known competing financial interests or personal relationships that could have appeared to influence the work reported in this paper.

Data availability

No data was used for the research described in the article.

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