

Lectures on Random Matrices (Spring 2025)

Lecture 11: Some universal asymptotics of Dyson Brownian Motion

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1 Recap

1.1 Dyson Brownian Motion (DBM)

We introduced a time-dependent model of random matrices by letting an $N \times N$ Hermitian matrix $\mathcal{M}(t)$ evolve in time so that each off-diagonal entry follows independent Brownian increments (real or complex depending on the symmetry class). Setting

$$\mathcal{M}(t) = \frac{1}{\sqrt{2}}(X(t) + X^\dagger(t)),$$

where $X(t)$ is an $N \times N$ matrix of i.i.d. Brownian motions, produces a self-adjoint matrix with a stochastically evolving spectrum. This model is full-rank matrix Brownian motion, and works well for $\beta = 1, 2, 4$. For other β , we need an SDE to describe the evolution of the eigenvalues (particles).

1.2 Eigenvalue SDE

Denote by $\lambda_1(t) \geq \dots \geq \lambda_N(t)$ the ordered eigenvalues of $\mathcal{M}(t)$. Dyson showed that these eigenvalues form a continuous-time Markov process satisfying the SDE

$$d\lambda_i(t) = \frac{\beta}{2} \sum_{j \neq i} \frac{dt}{\lambda_i(t) - \lambda_j(t)} + dW_i(t), \quad i = 1, \dots, N,$$

where $\beta > 0$ and $W_i(t)$ are independent standard real Brownian motions. For classical random matrix ensembles ($\beta = 1, 2, 4$), this SDE describes how the eigenvalues evolve under real symmetric (GOE), Hermitian (GUE), or quaternionic (GSE) Brownian motion — in the last [Lecture 10](#) we discussed the cases $\beta = 1, 2$ in detail. A key feature is the *repulsion* term $\frac{1}{\lambda_i - \lambda_j}$, which prevents collisions (and ensures the ordering remains intact).

1.3 Preservation of $G\beta E$ density

A fundamental result is that starting from all eigenvalues at 0, the distribution of $\lambda(t)$ at time t has the joint density proportional to

$$\prod_{i < j} |\lambda_i - \lambda_j|^\beta \exp\left\{-\frac{1}{2t} \sum_i \lambda_i^2\right\},$$

matching the Gaussian β -Ensemble ($G\beta E$) law. Hence DBM provides a dynamical realization of $G\beta E$. Invariance can be checked by verifying that this density is annihilated by the generator of the SDE.

1.4 Harish–Chandra–Itzykson–Zuber (HCIZ) integral

The HCIZ integral is a key tool for computing matrix integrals involving traces. For two Hermitian matrices A and B with eigenvalues (a_1, \dots, a_N) and (b_1, \dots, b_N) , it states (in one common normalization):

$$\int_{U(N)} \exp(\text{Tr}(A U B U^\dagger)) dU = \prod_{k=1}^{N-1} k! \frac{\det[e^{a_i b_j}]_{i,j=1}^N}{\prod_{1 \leq i < j \leq N} (a_j - a_i) \prod_{1 \leq i < j \leq N} (b_j - b_i)}.$$

This formula is instrumental in deriving transition densities for $\beta = 2$ Dyson Brownian Motion.

2 Optional: proof of HCIZ integral via representation theory

In this section, we outline a standard argument (adapted from the theory of symmetric functions and representation theory of the unitary group) that leads to a proof of the Harish–Chandra–Itzykson–Zuber formula. It is often referred to as the “orbital integral” or “character expansion” approach.

Step 1. Setting up the integral and Schur expansions. Let A and B be two $N \times N$ diagonalizable matrices, with eigenvalues a_1, \dots, a_N and $\lambda_1, \dots, \lambda_N$ respectively. Denote by $D_a = \text{diag}(a_1, \dots, a_N)$ and $D_\lambda = \text{diag}(\lambda_1, \dots, \lambda_N)$. We want to evaluate the integral

$$I = \int_{U(N)} \exp(\text{Tr}(D_a U D_\lambda U^\dagger)) dU$$

over the Haar measure on $U(N)$.

Since $\text{Tr}(B) = p_1(B)$ in the language of power sums (where $p_1(x_1, x_2, \dots) = x_1 + x_2 + \dots$), we have

$$\exp(\text{Tr}(B)) = \exp(p_1(B)).$$

One can use a known expansion [Mac95]

$$e^{p_1(B)} = \sum_{m=0}^{\infty} \frac{p_1^m(B)}{m!} = \sum_{m=0}^{\infty} \frac{1}{m!} \sum_{\mu: |\mu|=m} \dim(\mu) s_\mu(B),$$

where the sum is over all partitions μ of size m , and $s_\mu(\cdot)$ is the Schur polynomial (or Schur function) indexed by μ . The coefficient $\dim(\mu)$ is the dimension of the corresponding representation of S_m .

We set $B = D_a U D_\lambda U^\dagger$ and write

$$I = \int_{U(N)} \exp(\text{Tr}(D_a U D_\lambda U^\dagger)) dU = \int_{U(N)} \sum_{m=0}^{\infty} \frac{1}{m!} \sum_{\mu: |\mu|=m} \dim(\mu) s_\mu(D_a U D_\lambda U^\dagger) dU.$$

One can exchange the integral and the sum (the series converges absolutely for all matrix arguments), giving

$$I = \sum_{m=0}^{\infty} \sum_{\mu: |\mu|=m} \frac{\dim(\mu)}{m!} \int_{U(N)} s_\mu(D_a U D_\lambda U^\dagger) dU. \quad (2.1)$$

Step 2. Orthogonality of characters and the Unitary group. The Schur functions $s_\mu(\cdot)$ can be seen as irreducible characters of the unitary group $U(N)$ (up to a normalization factor) when restricted to N -tuples of eigenvalues.¹

¹ s_μ for $\ell(\mu) \leq N$ can be viewed as the character of the corresponding polynomial representation of $GL(N, \mathbb{C})$, then restricted to $U(N)$. If $\ell(\mu) > N$, the function s_μ vanishes on $U(N)$. Thus, we need to impose the condition $|a_i| = |\lambda_i| = 1$ (so that $D_a, D_\lambda \in U(N)$) to ensure immediate applicability of representation theory of $U(N)$, then extend to general $\{a_i\}$ and $\{\lambda_i\}$ by analytic continuation.

Proposition 2.1 (Functional equation for characters of compact groups). *Let G be a compact group with normalized Haar measure dh , and let χ be an irreducible character of a finite-dimensional representation of G . Then for any elements $g_1, g_2 \in G$, the following relation holds:*

$$\int_G \chi(g_1 h g_2 h^{-1}) dh = \frac{\chi(g_1) \chi(g_2)}{\dim V}, \quad (2.2)$$

where $\dim V = \chi(e)$ is the dimension of the representation space.

Remark 2.2. A similar relation holds for characters of finite groups.

By Proposition 2.1, the integral over $U(N)$ in (2.1) can be evaluated as

$$\int_{U(N)} s_\mu(D_a U D_\lambda U^\dagger) dU = \frac{1}{\text{Dim}_N(\mu)} s_\mu(a) s_\mu(\lambda),$$

where $\text{Dim}_N(\mu)$ is the dimension of the corresponding irreducible representation of $U(N)$. Substituting back into (2.1) yields

$$I = \sum_{m=0}^{\infty} \sum_{\mu: |\mu|=m, \ell(\mu) \leq N} \frac{\dim(\mu)}{m!} \frac{1}{\text{Dim}_N(\mu)} s_\mu(a) s_\mu(\lambda),$$

where $\ell(\mu) \leq N$ is needed for $s_\mu(\cdot)$ not to vanish on $U(N)$.

Step 3. Hook-length formulas and the final determinant. Next, one applies the hook-length formula and the hook-content formula to dimensions:

$$\dim \mu = \frac{|\mu|!}{\prod_{\square \in \mu} h(\square)}, \quad \text{Dim}_N(\mu) = \frac{\prod_{\square \in \mu} (N + c(\square))}{\prod_{\square \in \mu} h(\square)},$$

We have

$$\prod_{\square \in \mu} (N + c(\square)) = \prod_{i=1}^N \frac{(\mu_i + N - i)!}{(N - i)!},$$

so identifying $m_i = \mu_i + N - i$ gives

$$I = 0!1! \cdots (N-1)! \sum_{m_1 > \dots > m_N \geq 0} \frac{s_\mu(a) s_\mu(\lambda)}{m_1! \cdots m_N!},$$

which yields the HCIZ formula by the Cauchy-Binet summation.

3 Determinantal structure for $\beta = 2$

3.1 Transition density

Theorem 3.1 ($\beta = 2$ Dyson Brownian Motion Transition Probabilities). *For $\beta = 2$, let $\lambda(t) = (\lambda_1(t) \geq \dots \geq \lambda_N(t))$ follow Dyson Brownian Motion starting at $\lambda(0) = \mathbf{a} = (a_1 \geq \dots \geq a_N)$. Then for each fixed time $t > 0$,*

$$P(\lambda(t) = \mathbf{x} \mid \lambda(0) = \mathbf{a}) = N! \left(\frac{1}{\sqrt{2\pi t}} \right)^N \prod_{1 \leq i < j \leq N} \frac{x_i - x_j}{a_i - a_j} \det \left[\exp \left(-\frac{(x_i - a_j)^2}{2t} \right) \right]_{i,j=1}^N,$$

where $x_1 \geq \dots \geq x_N$.

Proof. Consider an $N \times N$ Hermitian matrix process $X(t)$ whose entries perform independent complex Brownian motions (so that $X(t)$ is distributed as $A + \sqrt{t}$ GUE at each fixed time, with $A = \text{diag}(a_1, \dots, a_N)$). Its eigenvalues $\lambda_1(t) \geq \dots \geq \lambda_N(t)$ evolve exactly according to the $\beta = 2$ Dyson Brownian Motion.

The density of X at time t , viewed as a random matrix, is proportional to

$$\exp\left(-\frac{1}{2t} \text{Tr}(X - A)^2\right).$$

If we replace A by $U A U^\dagger$ for any fixed unitary U , the law of X remains the same (this follows from the unitary invariance of the GUE). Thus the distribution of the eigenvalues of X is unchanged by such conjugation.

One writes

$$\int_{U(N)} \exp\left(-\frac{1}{2t} \text{Tr}(X - U A U^\dagger)^2\right) dU = (\text{const.}) \times [\text{HCIZ integral in the variables } (X, A)],$$

which by the Harish–Chandra–Itzykson–Zuber formula leads to a product of determinants and a factor that is precisely

$$\exp\left(-\frac{1}{2t} \sum_{i=1}^N x_i^2 - \frac{1}{2t} \sum_{i=1}^N a_i^2\right) \frac{\det\left[\exp\left(\frac{x_i a_j}{t}\right)\right]}{\prod_{i < j} (x_i - x_j)(a_i - a_j)},$$

where x_1, \dots, x_N are the eigenvalues of X .

To convert this matrix distribution into a distribution on eigenvalues alone, we multiply by the usual Vandermonde Jacobian $\prod_{i < j} (x_i - x_j)^2$ (which comes from integrating out the unitary degrees of freedom). This produces exactly

$$N! \left(\frac{1}{\sqrt{2\pi t}}\right)^N \prod_{i < j} \frac{x_i - x_j}{a_i - a_j} \det\left[\exp\left(-\frac{(x_i - a_j)^2}{2t}\right)\right].$$

Hence we obtain the stated transition probability for the Dyson Brownian Motion at $\beta = 2$. \square

Remark 3.2. The factor $N! \left(\frac{1}{\sqrt{2\pi t}}\right)^N$ arises naturally from normalizing the Gaussian increments and accounts for the ordering $\lambda_1 \geq \dots \geq \lambda_N$. The determinant and product factors encode the eigenvalue “repulsion” characteristic of $\beta = 2$ random matrices.

3.2 Determinantal correlations

Theorem 3.3 (Determinantal structure for $\beta = 2$ DBM). *Let $\{x_1(t), \dots, x_n(t)\}$ be the eigenvalues at time $t > 0$ of the $\beta = 2$ Dyson Brownian Motion started at initial locations (a_1, \dots, a_n) at time 0. Equivalently, these $x_i(t)$ are the eigenvalues of*

$$A + \sqrt{t} G,$$

where $A = \text{diag}(a_1, \dots, a_n)$ and G is a random Hermitian matrix from the GUE. Then the (random) point configuration $\{x_i(t)\}$ forms a determinantal point process with correlation kernel

$$K_t(x, y) = \frac{1}{(2\pi)^2 t} \oint \oint \exp\left(\frac{w^2 - 2yw}{2t}\right) \Big/ \exp\left(\frac{z^2 - 2xz}{2t}\right) \prod_{i=1}^n \frac{w - a_i}{z - a_i} \frac{dw dz}{w - z}.$$

Here z goes around all the points a_1, \dots, a_n in the positive direction, and the w contour passes from $-i\infty$ to $i\infty$, to the right of the z contour.

- If $a_1 = \dots = a_n = 0$ and $t = 1$, this kernel reduces to the familiar correlation kernel of the GUE (see [Lecture 6](#)).
- One can use this formula to study the Baik–Ben Arous–Péché (BBP) [\[BBP05\]](#) phase transition for $\beta = 2$, which deals with finite rank perturbations of the GUE random matrix ensemble. Indeed, rank r perturbation corresponds to taking $a_1, \dots, a_r \neq 0$, and $a_{r+1} = \dots = a_n = 0$.

3.3 On the proof of determinantal structure

The idea of the proof of Theorem [3.3](#) is to represent the measure (the transition density) as a product of determinants. In general, if a measure is given as a product of determinants, there is a well-studied method (biorthogonal ensembles and, more generally, the Eynard–Mehta theorem) to compute the determinantal correlation kernel. We refer to [\[BR05\]](#), [\[Bor11\]](#) for a detailed exposition in the discrete case (which is arguably more transparent). The first step for the Dyson Brownian Motion is as follows.

Lemma 3.4 (Density representation). *Let $P_t(x \rightarrow y)$ be the transition probability kernel of standard Brownian motion,*

$$P_t(x \rightarrow y) = \frac{1}{\sqrt{2\pi t}} \exp\left(-\frac{(x-y)^2}{2t}\right).$$

Then the density of the eigenvalues (x_1, \dots, x_N) of DBM started at (a_1, \dots, a_N) at time 0 admits the representation

$$\lim_{s \rightarrow \infty} \left(\frac{1}{Z}\right) \det\left[P_t(a_i \rightarrow x_j)\right]_{i,j=1}^N \det\left[P_s(x_i \rightarrow k-1)\right]_{i,k=1}^N. \quad (3.1)$$

Remark 3.5. This representation [\(3.1\)](#) is related to an alternative description of the $\beta = 2$ Dyson Brownian Motion as an ensemble of noncolliding Brownian motions (that is, independent Brownian motions, conditioned to never collide).

Proof of Lemma 3.4. The first determinant (as $s \rightarrow \infty$) matches the determinant we have in Theorem [3.1](#). It remains to analyze the second determinant

$$\det\left[P_s(x_j \rightarrow k-1)\right]_{j,k=1}^N = \det\left[\frac{1}{\sqrt{2\pi s}} \exp\left(-\frac{((k-1)-x_j)^2}{2s}\right)\right]_{j,k=1}^N.$$

We may ignore the factor $\frac{1}{\sqrt{2\pi s}}$ in each entry since it does not depend on x_j . Inside the exponential,

$$-\frac{((k-1)-x_j)^2}{2s} = -\frac{x_j^2}{2s} + \frac{x_j(k-1)}{s} - \frac{(k-1)^2}{2s}.$$

Thus, up to the factor $\exp(-\frac{(k-1)^2}{2s})$ (which depends only on k and hence is independent of each x_j), we can factor out $\exp(-\frac{x_j^2}{2s})$ from row j . Consequently, the nontrivial part of the determinant becomes

$$\det \left[e^{\frac{x_j (k-1)}{s}} \right]_{j,k=1}^N.$$

Recognize this as a Vandermonde-type determinant in the variables $e^{x_j/s}$. Indeed,

$$\det \left[e^{\frac{x_j (k-1)}{s}} \right]_{j,k=1}^N = \prod_{1 \leq i < j \leq N} \left(e^{\frac{x_i}{s}} - e^{\frac{x_j}{s}} \right).$$

As $s \rightarrow \infty$, we expand $e^{\frac{x_i}{s}} = 1 + \frac{x_i}{s} + O(\frac{1}{s^2})$, so each difference $(e^{\frac{x_i}{s}} - e^{\frac{x_j}{s}}) \sim \frac{x_i - x_j}{s}$. Hence,

$$\prod_{1 \leq i < j \leq N} \left(e^{\frac{x_i}{s}} - e^{\frac{x_j}{s}} \right) \sim \frac{1}{s^{\frac{N(N-1)}{2}}} \prod_{1 \leq i < j \leq N} (x_i - x_j).$$

Combining all these factors and matching with the first determinant (as $s \rightarrow \infty$) verifies the claimed product form, up to overall constants that do not depend on the variables x_j . This completes the proof. \square

Then, the product of determinants idea (biorthogonal ensembles) applies to the density (3.1) before the limit $s \rightarrow \infty$, and simplifies after taking the limit. We omit the details here, see Problem K.1.

4 Asymptotic analysis: signal plus noise

4.1 Setup

In this section, we provide a detailed derivation of how the rank-1 spike

$$A + \sqrt{G}, \quad A = \text{diag}(a, 0, \dots, 0) \quad \text{with } a \in \mathbb{R},$$

affects the large- n and large-time behavior of the Dyson Brownian Motion at $\beta = 2$. See the simulation at <https://lpetrov.cc/simulations/2025-01-28-bbp-transition/>.

We set $a_1 = a\sqrt{n}$ and $a_2 = a_3 = \dots = a_n = 0$, which simplifies the product:

$$\prod_{i=1}^n (w - a_i) = (w - a\sqrt{n}) w^{n-1}, \quad \prod_{i=1}^n (z - a_i) = (z - a\sqrt{n}) z^{n-1}.$$

Let us also take $t = 1$ for simplicity, so that the limit shape (at least in the case $a = 0$, but also in general) is supported by $[-2\sqrt{n}, 2\sqrt{n}]$. Let us also make the change of the integration variables $w \rightarrow w\sqrt{n}$, $z \rightarrow z\sqrt{n}$.

Hence, the correlation kernel becomes

$$K_t(x, y) = \frac{\sqrt{n}}{(2\pi)^2} \oint \oint \exp\left(\frac{nw^2 - 2yw\sqrt{n}}{2}\right) \Big/ \exp\left(\frac{nz^2 - 2xz\sqrt{n}}{2}\right) \frac{w - a}{z - a} \left(\frac{w}{z}\right)^{n-1} \frac{dw dz}{w - z}. \quad (4.1)$$

Here:

- The z -contour is a small positively oriented loop around $z = a$, and also around $z = 0$, so that it encircles these two singularities but excludes w .
- The w -contour is a vertical line (or an equivalent contour from $-i\infty$ to $i\infty$) passing to the right of all singularities (i.e. to the right of z).

Note that to capture the edge behavior, we need to set $x = y = 2$ plus lower order terms. Let us make this substitution $x = 2\sqrt{n} + x'$, $y = 2\sqrt{n} + y'$, and the scale of x', y' will be determined later (but for now we assume that they are $o(\sqrt{n})$).

4.2 Outline of the steepest descent approach

We aim to understand the behavior of (4.1) in the regime $n \rightarrow \infty$, especially near the largest eigenvalue $\lambda_1(t)$. Recall from standard GUE (i.e. $a = 0$) that the top of the spectrum is about $2\sqrt{n}$. The presence of the rank-1 spike a can drastically modify the top eigenvalue if a is large enough to produce an “outlier.” Our goal is to detect precisely how this occurs by analyzing the double contour integral via steepest descent.

For large n , the integral localizes around these double critical point. Any crossing from z - to w -contour may pick up residues, which account for separate contributions (leading, for instance, to the Airy kernel in the unperturbed GUE). We track how the spike a changes these deformations.

4.3 Exponent form and critical points

Set

$$S(w; y') = \frac{w^2}{2} - 2w - y'w/\sqrt{n} + \frac{n-1}{n} \ln(w).$$

Then the integrand in (4.1) is

$$\frac{\exp\{n[S(w; y') - S(z; x')]\}}{w - z} \frac{w - a}{z - a}.$$

To capture the Airy behavior, we can ignore y' , and find the double critical point of $S(w; 0)$. It is equal to $w_c = 1$, and we would like to bring the z and w contours to intersect at $w_c = 1$. Note however that the old z contour must encircle $z = a$ and $z = 0$, and $z = a$ is a pole of the integrand. The w contour must always be to the right of the z contour.

We see that there are three regimes:

- If $a < 1$, we can deform the z contour to encircle $z = 0$ and $z = a$, and the w contour to pass through $w = 1$. This will lead to the Airy kernel, and the derivation is the same as in [Lecture 7](#). We obtain

$$z = 1 + \frac{Z}{n^{1/3}}, \quad w = 1 + \frac{W}{n^{1/3}},$$

$i++i$

- If $a = 1$, the behavior is going to be critical — we still will be able to get the same scaling, but the limiting kernel will be different.
- Finally,

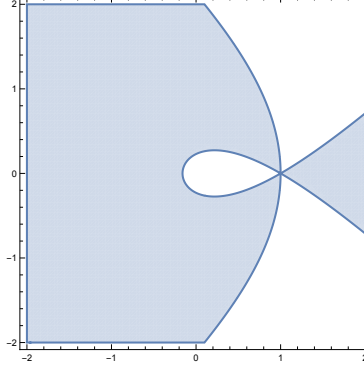


Figure 1: The plot of the region $\operatorname{Re} S(z) - \operatorname{Re} S(1) > 0$ at the edge, in the neighborhood of the double critical point $w_c = 1$. The new w contour should pass through the shaded region, and the new z must stay in the non-shaded region.

Remark 4.1. For $\beta = 2$ and $a = 0$, we typically expand about $z \approx 1$, $w \approx 1$ to capture the top edge at $2\sqrt{tn}$. Indeed, setting $z = 1 + u$, $n \rightarrow \infty$ leads to the cubic expansions in u . Once $a \neq 0$, we get an extra shift in z , w that might solve

$$\partial_z \Phi_{t,a} = 0, \quad \partial_w \Phi_{t,a} = 0.$$

We find solutions near $z = 1$, $w = 1$ if $|a| < 2$, but new solutions appear if $|a| > 2$.

Scaling near the top. We know from the unperturbed case ($a = 0$) that x, y near $2\sqrt{tn}$ is relevant. Let us adopt the usual *edge scaling*:

$$x = 2\sqrt{tn} + \frac{\xi}{n^{1/6}}, \quad y = 2\sqrt{tn} + \frac{\eta}{n^{1/6}}. \quad (4.2)$$

Then one can attempt expansions $z = 1 + u n^{-1/3}$, $w = 1 + v n^{-1/3}$ (plus additional shifts if $a \neq 0$). Indeed, if $|a| < 2$, the main location for $z_{\text{cr}}, w_{\text{cr}}$ remains close to 1. Conversely, if $|a| > 2$, we get a separate solution.

Let us isolate the product factor first. Suppose we rewrite

$$(w - a) w^{n-1} = (w - a) \exp((n-1) \ln w), \quad (z - a) z^{n-1} = (z - a) \exp((n-1) \ln z).$$

Then

$$\ln[(w - a) w^{n-1}] = (n-1) \ln w + \ln(w - a).$$

Hence these combined terms become

$$(n-1) [\ln w - \ln z] + \ln(w - a) - \ln(z - a).$$

Because $|a| < 2$ or $|a| > 2$ drastically changes the location(s) $z - a = 0$ or $w - a = 0$, the steepest descent approach tries to see where the real part of $\Phi_{t,a}(w, z)$ is stationary.

4.4 Controlling the integral: subcritical and supercritical analyses

4.4.1 Residue crossing and main contributions

As usual in the $\beta = 2$ story, we must deform the z -contour outward so that it crosses the w -contour. This can pick up a residue from the simple pole at $w = z$. In the pure GUE case ($a = 0$), that residue is known to yield the sine or Airy kernel asymptotics after expansions. Now, with the rank-1 spike, we get an additional factor from $(w - a)/(z - a)$. The upshot is that *if* the z contour crosses w in a certain region that does not include $w = a$, one obtains a residue that will lead to the same sine/Airy factor.

However, we must also check whether the z -contour encloses $z = a$ *and* the w -contour encloses $w = a$. The spike can produce an outlier contribution if $|a|$ is large enough for the new saddle point to dominate.

4.4.2 Case 1: $|a| < 2$ (subcritical)

For $|a| < 2$, one can show that the exponent $\Phi_{t,a}(w, z)$ remains effectively minimized near $z, w \approx 1$, and the presence of $(w - a)$ vs. $(z - a)$ does not shift that minimizer significantly. In more precise terms, one writes expansions:

$$z = 1 + \frac{u}{n^{1/3}}, \quad w = 1 + \frac{v}{n^{1/3}},$$

with x, y as in (4.2), and obtains a leading cubic term in u, v . The “ a ” correction modifies subleading expansions but does *not* cause a different order in n . Repeating the same steps as in the standard GUE edge analysis (cf. §??), one sees that the limit of $\frac{1}{n^{1/6}} K_t(x, y)$ is the Airy kernel

$$K_{\text{Ai}}(\xi, \eta) = \int_0^\infty \dots$$

(as in the usual expression, see ?? for one common form).

Consequently, in the subcritical regime $|a| < 2$, the top eigenvalue

$$x_{\max}(t) = x_1(t)$$

stays merged in the main bulk near $2\sqrt{tn}$ and has *Tracy–Widom* fluctuations on the $n^{-1/6}$ scale. The factor a only shifts lower-order terms in that expansion.

4.4.3 Case 2: $|a| > 2$ (supercritical)

When $|a| > 2$, the spike is large. One expects an *outlier* eigenvalue near $x \approx a\sqrt{t}$ (rather than near $2\sqrt{tn}$). Indeed, analyzing $\Phi_{t,a}$ in that scenario, one finds a separate solution to $\partial\Phi/\partial w = 0$ near $w \approx a \neq 1$ (and similarly for z). The portion of the z -contour around $z = a$ then yields a new local expansion.

Concretely, if $a > 2$ (the case $a < -2$ is similar), the factor $(w - a)$ can be near zero, so that the exponent can develop a more favorable real part at $w \approx a$. Meanwhile, the factor w^{n-1} is enormous if $|w| > 1$, but $a > 2$ means $|a| > 1$. As a result, the outlier solution with $w \approx a$ can produce a stable stationary phase. Indeed, one obtains a local *quadratic* expansion around

$w = a$ and $z = a$, thus leading to a *Gaussian* fluctuation scale of the resulting outlier eigenvalue, typically $\sigma n^{-1/2}$ for some $\sigma > 0$.

Hence, in the supercritical regime $|a| > 2$, the top eigenvalue $x_{\max}(t)$ *detaches* from $2\sqrt{tn}$. Its leading order becomes $a\sqrt{t}$ plus smaller corrections, and the local fluctuations are usually governed by a (shifted) Gaussian (or in some references, an incomplete gamma distribution, depending on subtle normalizations). The remaining $n-1$ eigenvalues fill out the usual semicircle from $-2\sqrt{tn}$ to $2\sqrt{tn}$, with an Airy edge near $2\sqrt{tn}$.

4.4.4 Case 3: $|a| = 2$ (critical)

At the boundary $|a| = 2$, the system is at a “double root” transition. One finds that $z = 1$ (the standard GUE edge) merges with $z = a$ in a certain sense, leading to a triple or quartic expansion if $a = 2$, for instance. Careful expansions show that the outlier is *just* beginning to separate from the main bulk. The correlation kernel in this regime is a special transition “higher-order” Airy-like kernel, often described as a *crossover kernel* in the BBP (Baik–Ben Arous–Péché) transition [BBP05]. The top eigenvalue distribution is neither purely Tracy–Widom nor purely Gaussian but an interpolating limit distribution.

4.5 Summary of the BBP transition from $K_t(x, y)$

Putting the above cases together, we see that analyzing the DBM correlation kernel (4.1) under the spiked initial condition ($a_1 = a \neq 0$, $(a_2 = \dots = a_n = 0)$) reveals:

- **If $|a| < 2$:** The spike is *subcritical*. The top eigenvalue remains at the GUE edge $\approx 2\sqrt{tn}$, with Tracy–Widom fluctuations on scale $n^{-1/6}$.
- **If $|a| > 2$:** The spike is *supercritical*. An outlier eigenvalue emerges near $a\sqrt{t}$, detaching from the main bulk. The fluctuation scale around that outlier is typically $n^{-1/2}$ and converges (after centering and scaling) to a simpler, often Gaussian-type limit. The remainder of the spectrum still exhibits the semicircle on $[-2\sqrt{tn}, 2\sqrt{tn}]$.
- **If $|a| = 2$:** *Critical regime*. The top eigenvalue experiences a BBP crossover distribution, with a kernel that can be viewed as a limit of the general (4.1) by carefully choosing the saddle point expansions.

This completes the steepest descent derivation of the rank-1 BBP phase transition from the Dyson Brownian Motion correlation kernel in Theorem 3.3. Detailed error estimates and rigorous Fredholm determinant limits (turning kernel statements into actual distributional statements about $x_{\max}(t)$) can be found in [BBP05], [BloemendalVirag], and other references cited therein.

Remark 4.2. For r -rank spikes (i.e. $a_1, \dots, a_r \neq 0$), one can repeat the same argument: the factor $\prod_{i=1}^n (w - a_i)/(z - a_i)$ in (4.1) splits into a product $(w - a_1) \dots (w - a_r) w^{n-r}$ vs. $(z - a_1) \dots (z - a_r) z^{n-r}$. Each $|a_i| > 2$ can produce a distinct outlier, while any $|a_i| < 2$ merges with the main bulk.

K Problems (due 2025-04-29)

K.1 Biorthogonal ensembles

Derive Theorem 3.3 from Lemma 3.4 using the orthogonalization process similar to Lecture 5, and then taking the limit as $s \rightarrow \infty$.

K.2 Scaling of the kernel

Let $a_i = 0$ in Theorem 3.3. Find α such that $t^\alpha K_t(x/\sqrt{t}, y/\sqrt{t})$ is independent of t . Can you explain this value of α ?

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