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On the principal components of sample covariance matrices

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We introduce a class of $M \times M$ sample covariance matrices $\mathcal{Q}$ which subsumes and generalizes several previous models. The associated population covariance matrix $\Sigma = \mathbb{E}\mathcal{Q}$ is assumed to differ from the identity by a matrix of bounded rank. All quantities except the rank of $\Sigma - I_M$ may depend on M in an arbitrary fashion. We investigate the principal components, i.e. the top eigenvalues and eigenvectors, of $\mathcal{Q}$. We derive precise large deviation estimates on the generalized components $\langle \mathbf{w}, \boldsymbol{\xi}_i \rangle$ of the outlier and non-outlier eigenvectors $\boldsymbol{\xi}_i$. Our results also hold near the so-called BBP transition, where outliers are created or annihilated, and for degenerate or near-degenerate outliers. We believe the obtained rates of convergence to be optimal. In addition, we derive the asymptotic distribution of the generalized components of the non-outlier eigenvectors. A novel observation arising from our results is that, unlike the eigenvalues, the eigenvectors of the principal components contain information about the *subcritical* spikes of Σ .

The proofs use several results on the eigenvalues and eigenvectors of the uncorrelated matrix $\mathcal{Q}$, satisfying $\mathbb{E}\mathcal{Q} = I_M$, as input: the isotropic local Marchenko-Pastur law established in [9], level repulsion, and quantum unique ergodicity of the eigenvectors. The latter is a special case of a new universality result for the joint eigenvalue-eigenvector distribution.

1. Introduction

In this paper we investigate $M \times M$ sample covariance matrices of the form

$$\mathcal{Q} = \frac{1}{N} A A^* = \left(\frac{1}{N} \sum_{\mu=1}^N A_{i\mu} A_{j\mu} \right)_{i,j=1}^M, \quad (1.1)$$

where the *sample matrix* $A = (A_{i\mu})$ is a real $M \times N$ random matrix. The main motivation to study such models stems from multivariate statistics. Suppose we are interested in the statistics of M mean-zero variables $\mathbf{a} = (a_1, \dots, a_M)^*$ which are thought to possess a certain degree of interdependence. Such problems of multivariate statistics commonly arise in population genetics, economics, wireless communication, the physics of mixtures, and statistical learning [3, 24, 30]. The goal is to unravel the interdependencies among the variables $\mathbf{a}$ by finding the *population covariance matrix*

$$\Sigma = \mathbb{E} \mathbf{a} \mathbf{a}^* = (\mathbb{E} a_i a_j)_{i,j=1}^M. \quad (1.2)$$

To this end, one performs a large number, N , of repeated, independent measurements, called “samples”, of the variables $\mathbf{a}$. Let $A_{i\mu}$ denote the value of a_i in the μ -th sample. Then the sample covariance matrix (1.1) is the empirical mean approximating the population covariance matrix Σ .

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In general, the mean of the variables $\mathbf{a}$ is nonzero and unknown. In that case, the population covariance matrix (1.2) has to be replaced with the general form

$$\Sigma = \mathbb{E}[(\mathbf{a} - \mathbb{E}\mathbf{a})(\mathbf{a} - \mathbb{E}\mathbf{a})^*].$$

Correspondingly, one has to subtract from $A_{i\mu}$ the empirical mean of the i -th row of A , which we denote by $[A]_i := \frac{1}{N} \sum_{\mu=1}^N A_{i\mu}$. Hence, we replace (1.1) with

$$\dot{Q} := \frac{1}{N-1} A(I_N - \mathbf{e}\mathbf{e}^*)A = \left(\frac{1}{N-1} \sum_{\mu=1}^N (A_{i\mu} - [A]_i)(A_{j\mu} - [A]_j) \right)_{i,j=1}^M, \quad (1.3)$$

where we introduced the vector

$$\mathbf{e} := N^{-1/2}(1, 1, \dots, 1)^* \in \mathbb{R}^N. \quad (1.4)$$

Since $\dot{Q}$ is invariant under the shift $A_{i\mu} \mapsto A_{i\mu} + m_i$ for any deterministic vector $(m_i)_{i=1}^M$, we may assume without loss of generality that $\mathbb{E}A_{i\mu} = 0$. We shall always make this assumption from now on.

It is easy to check that $\mathbb{E}Q = \mathbb{E}\dot{Q} = \Sigma$. Moreover, we shall see that the principal components of Q and $\dot{Q}$ have identical asymptotic behaviour. For simplicity of presentation, in the following we focus mainly on Q , bearing in mind that every statement we make on Q also holds verbatim for $\dot{Q}$ (see Theorem 2.22 below).

By the law of large numbers, if M is fixed and N taken to infinity, the sample covariance matrix Q converges almost surely to the population covariance matrix Σ . In many modern applications, however, the population size M is very large and obtaining samples is costly. Thus, one is typically interested in the regime where M is of the same order as N , or even larger. In this case, as it turns out, the behaviour of Q changes dramatically and the problem becomes much more challenging. In *principal component analysis*, one seeks to understand the correlations by considering the *principal components*, i.e. the top eigenvalues and associated eigenvectors, of Q . These provide an effective low-dimensional projection of the high-dimensional data set A , in which the significant trends and correlations are revealed by discarding superfluous data.

The fundamental question, then, is how the principal components of $\Sigma = \mathbb{E}Q$ are related to those of Q .

1.1. The uncorrelated case. In the “null” case, the variables $\mathbf{a}$ are uncorrelated and $\Sigma = I_M$ is the identity matrix. The global distribution of the eigenvalues is governed by the *Marchenko-Pastur law* [28]. More precisely, defining the dimensional ratio

$$\phi \equiv \phi_N := \frac{M}{N}, \quad (1.5)$$

the empirical eigenvalue density of the rescaled matrix $Q = \phi^{-1/2} \dot{Q}$ has the same asymptotics for large M and N as

$$\frac{\sqrt{[(x - \gamma_-)(\gamma_+ - x)]_+}}{2\pi\sqrt{\phi}x} dx + (1 - \phi^{-1})_+ \delta(dx), \quad (1.6)$$

where we defined

$$\gamma_{\pm} := \phi^{1/2} + \phi^{-1/2} \pm 2 \quad (1.7)$$

to be the edges of the limiting spectrum. Hence, the unique nontrivial eigenvalue 1 of Σ spreads out into a bulk spectrum of Q with diameter $4\phi^{1/2}$. Moreover, the local spectral statistics are universal; for instance, the top eigenvalue of Q is distributed according to the Tracy-Widom-1 distribution [22, 23, 39, 40]. Finally, Q satisfies *quantum unique ergodicity*, i.e. its eigenvectors are uniformly distributed on the unit sphere of $\mathbb{R}^M$. We refer to Theorem 8.3 and Remark 8.4 below for precise statements.

1.2. Examples and outline of the model. The problem becomes much more interesting if the variables $\mathbf{a}$ are correlated. Several models for correlated data have been proposed in the literature, starting with the Gaussian spiked model from the seminal paper of Johnstone [23]. Here we propose a general model which includes many previous models as special cases. We motivate it using two examples.

- (1) Let $\mathbf{a} = T\mathbf{b}$, where the entries of $\mathbf{b}$ are independent with zero mean and unit variance, and T is a deterministic $M \times M$ matrix. This may be interpreted as an observer studying a complicated system whose randomness is governed by many independent internal variables $\mathbf{b}$. The observer only has access to the external variables $\mathbf{a}$, which may depend on the internal variables $\mathbf{b}$ in some complicated and unknown fashion. Assuming that this dependence is linear, we obtain $\mathbf{a} = T\mathbf{b}$. The sample matrix for this model is therefore $A = TB$, where B is an $M \times N$ matrix with independent entries of unit variance. The population covariance matrix is $\Sigma = TT^*$.

- (2) Let $r \in \mathbb{N}$ and set

$$\mathbf{a} = \mathbf{z} + \sum_{l=1}^r y_l \mathbf{u}_l.$$

Here $\mathbf{z} \in \mathbb{R}^M$ is a vector of “noise”, whose entries are independent with zero mean and unit variance. The “signal” is given by the contribution of r terms of the form $y_l \mathbf{u}_l$, whereby $y_1, \dots, y_r$ are independent, with zero mean and unit variance, and $\mathbf{u}_1, \dots, \mathbf{u}_r \in \mathbb{R}^M$ are arbitrary deterministic vectors. The sample matrix is

$$A = Z + \sum_{l=1}^r \mathbf{u}_l \mathbf{y}_l^*,$$

where, writing $Y := [\mathbf{y}_1, \dots, \mathbf{y}_r] \in \mathbb{R}^{N \times r}$, the $(M+r) \times N$ matrix $B := \begin{pmatrix} Z \\ Y \end{pmatrix}$ has independent entries with zero mean and unit variance. Writing $U := [\mathbf{u}_1, \dots, \mathbf{u}_r] \in \mathbb{R}^{M \times r}$, we therefore have

$$A = TB, \quad T := (I_M, U).$$

The population covariance matrix is $\Sigma = TT^* = I_M + UU^*$.

Below we shall refer to these examples as Examples (1) and (2) respectively. Motivated by them, we now outline our model. Let B be an $(M+r) \times N$ matrix whose entries are independent with zero mean and unit variance. We choose a deterministic $M \times (M+r)$ matrix T , and set

$$\mathcal{Q} = \frac{1}{N} T B B^* T^*. \tag{1.8}$$

We stress that we do not assume that the underlying randomness is Gaussian; we shall allow the entries of B to be arbitrary (not necessarily identically distributed) random variables possessing enough moments. Moreover, the assumption that the variance of the entries of B be equal to one is trivial: it can always be satisfied by a simple rescaling the rows of B , which may be absorbed into the deterministic matrix T . It is easy to check that the population covariance matrix associated with the model (1.8) is given by $\Sigma := TT^*$. As in our second example above, we always assume that r and the rank of $\Sigma - I_M$ are bounded.

As explained around (1.3), in addition to $\mathcal{Q}$ we also consider the matrix

$$\dot{\mathcal{Q}} = \frac{1}{N-1} T B (I_N - \mathbf{e}\mathbf{e}^*) B^* T^*, \tag{1.9}$$

whose principal components turn out to have the same asymptotic behaviour as those of $\mathcal{Q}$.

1.3. Sketch of behaviour of the principal components of $\mathcal{Q}$. To guide the reader, we now give a heuristic description of the behaviour of principal components of $\mathcal{Q}$. For convenience, we shall always work with the rescaled sample covariance matrix

$$Q := \phi^{-1/2} \mathcal{Q}. \tag{1.10}$$

The motivation behind this rescaling is that, as observed in (1.6), it ensures that the bulk spectrum of Q has asymptotically a fixed diameter, 4, for arbitrary N and M . Since TB is an $M \times N$ matrix, we find that Q has

$$K \equiv K_N := \min\{M, N\} \quad (1.11)$$

nontrivial (i.e. nonzero) eigenvalues. We use the notation

$$\Sigma := TT^* = \sum_{i=1}^M \sigma_i \mathbf{v}_i \mathbf{v}_i^* = I_M + \phi^{1/2} \sum_{i=1}^M d_i \mathbf{v}_i \mathbf{v}_i^* \quad (1.12)$$

for the spectral decomposition of the nonnegative matrix Σ , where $\{\mathbf{v}_i\}_{i=1}^M$ is a real orthonormal basis of $\mathbb{R}^M$ and $\{\sigma_i\}_{i=1}^M$ are the eigenvalues of Σ . Here we introduce the representation

$$\sigma_i = 1 + \phi^{1/2} d_i$$

for the eigenvalues σ_i . We always order the values d_i such that $d_1 \geq d_2 \geq \dots \geq d_M$. By assumption on Σ , only a bounded number of d_i 's are nonzero. They, together with the associated eigenvectors $\mathbf{v}_i$, are commonly referred to as the *spikes* of Σ .

We are interested in the eigenvalues of Q , denoted by

$$\mu_1 \geq \mu_2 \geq \dots \geq \mu_M,$$

and the associated unit eigenvectors of Q , denoted by

$$\xi_1, \xi_2, \dots, \xi_M \in \mathbb{R}^M.$$

The spectrum of Q consists of a *bulk spectrum* and of *outliers*—eigenvalues separated from the bulk. The bulk contains an order K eigenvalues, which are distributed on large scales according to the Marchenko-Pastur law (1.6). In addition, if $\phi > 1$ there are $M - K$ trivial eigenvalues at zero. Each d_i satisfying $|d_i| > 1$ gives rise to an outlier located near its *classical location*

$$\theta(d) := \phi^{1/2} + \phi^{-1/2} + d + d^{-1}. \quad (1.13)$$

Any d_i satisfying $|d_i| < 1$ does not result in an outlier. We summarize this picture in Figure 1.1. The

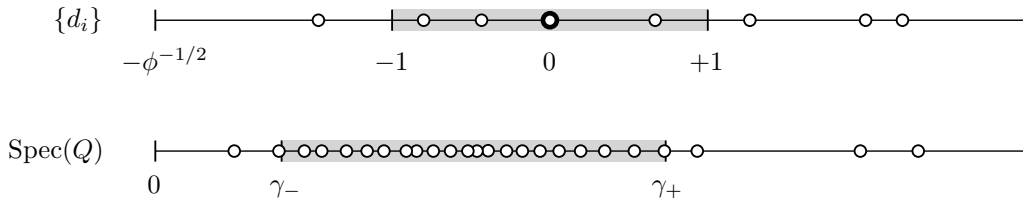


FIGURE 1.1. A typical configuration of $\{d_i\}$ (above) and the resulting spectrum of Q (below). An order M of the d_i 's are zero, which is symbolized by the thicker dot at 0. Any d_i inside the grey interval $[-1, 1]$ does not give rise to an outlier, while any d_i outside the grey interval gives rise to an outlier located near its classical location $\theta(d_i)$ and separated from the bulk $[\gamma_-, \gamma_+]$.

creation or annihilation of an outlier as a d_i crosses ± 1 is known as the *BBP phase transition* [3]. It takes place on the scale $||d_i| - 1| \sim K^{-1/3}$. This scale has a simple heuristic explanation (we focus on the right edge of the spectrum). Suppose that $d_1 \in (0, 1)$ and all other d_i 's are zero. Then the top eigenvalue μ_1 exhibits universality, and fluctuates on the scale $K^{-2/3}$ around γ_+ (see Theorem 8.3 and Remark 8.7 below). Increasing d_1 beyond the critical value 1, we therefore expect μ_1 to become an outlier when its

classical location $\theta(d_1)$ is located at a distance greater than $K^{-2/3}$ from γ_+ . By a simple Taylor expansion of θ , the condition $\theta(d_1) - \gamma_+ \gg K^{-2/3}$ becomes $d_1 - 1 \gg K^{-1/3}$.

We conclude this subsection by outlining the distribution of the outlier eigenvectors. Let μ_i be an outlier with associated eigenvector ξ_i . Then ξ_i is concentrated on a cone [7, 30] with axis parallel to $\mathbf{v}_i$, the corresponding eigenvector of the population covariance matrix Σ . More precisely, assuming that the eigenvalue $1 + \phi^{1/2}d_i$ of Σ is simple, we have with high probability

$$\langle \mathbf{v}_i, \xi_i \rangle^2 \approx u(d_i), \quad (1.14)$$

where we defined

$$u(d_i) \equiv u_\phi(d_i) := \frac{\sigma_i}{\phi^{1/2}\theta(d_i)}(1 - d_i^{-2}) \quad (1.15)$$

for $d_i > 1$. The function u determines the aperture $2 \arccos \sqrt{u(d_i)}$ of the cone. Note that $u(d_i) \in (0, 1)$ and $u(d_i)$ converges to 1 as $d_i \rightarrow \infty$. See Figure 1.2.

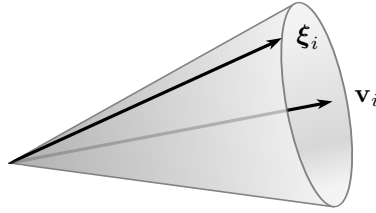


FIGURE 1.2. The eigenvector ξ_i associated with an outlier μ_i is concentrated on a cone with axis parallel to $\mathbf{v}_i$. The aperture of the cone is determined by $u(d_i)$ defined in (1.15).

1.4. Summary of previous related results. There is an extensive literature on spiked covariance matrices. So far most of the results have focused on the outlier eigenvalues of Example (1), with the nonzero d_i independent of N and ϕ fixed. Eigenvectors and the non-outlier eigenvalues have seen far less attention.

For the uncorrelated case $\Sigma = I_M$ and Gaussian B in (1.8) with fixed ϕ , it was proved in [22] for the complex case and in [23] for the real case that the top eigenvalue, rescaled as $K^{2/3}(\mu_1 - \gamma_+)$, is asymptotically distributed according to the Tracy-Widom law of the appropriate symmetry class [39, 40]. Subsequently, these results were shown to be universal, i.e. independent of the distribution of the entries of B , in [33, 37]. The assumption that ϕ be fixed was relaxed in [15, 32].

The study of covariance matrices with nontrivial population covariance matrix $\Sigma \neq I_M$ goes back to the seminal paper of Johnstone [23], where the Gaussian spiked model was introduced. The BBP phase transition was established by Baik, Ben Arous, and P       [3] for the case of Example (1) with complex Gaussian B , fixed rank of $\Sigma - I_M$, and fixed ϕ . Subsequently, the results of [3] were extended to the other Gaussian symmetry classes, such as real covariance matrices, in [10, 11]. The proofs of [3, 31] use an asymptotic analysis of Fredholm determinants, while those of [10, 11] use an explicit tridiagonal representation of BB^* ; both of these approaches rely heavily on the Gaussian nature of B .

For the model from Example (1) with fixed nonzero $\{d_i\}$ and ϕ , the almost sure convergence of the outliers was established in [4]. It was also shown in [4] that if $|d_i| < 1$ for all i , the top eigenvalue μ_1 converges to γ_+ . For this model, a central limit theorem of the outliers was proved in [2]. In [1], the almost sure convergence of the outliers was proved for a generalized spiked model whose population covariance matrix is of the block diagonal form $\Sigma = \text{diag}(A, T)$, where A is a fixed $r \times r$ matrix and T is chosen so that the associated sample covariance matrix has no outliers.

In [7], the almost sure convergence of the projection of the outlier eigenvectors onto the finite-dimensional spike subspace was established, under the assumption that ϕ and the nonzero d_i are fixed, and that B and T are both random and one of them is orthogonally invariant. In particular, the cone concentration from

(1.14) was established in [7]. In [30], under the assumption that B is Gaussian and ϕ and the nonzero d_i are fixed, a central limit theorem for a certain observable, the so-called sample vector, of the outlier eigenvectors was established. The result of [30] was extended to non-Gaussian entries for a special class of Σ in [36].

Moreover, in [8, 29] result analogous to those of [7] were obtained for the model from Example (2). Finally, a related class of models, so-called deformed Wigner matrices, have been the subject of much attention in recent years; we refer to [25, 26, 34, 35] for more details; in particular, the joint distribution of all outliers was derived in [26].

1.5. Outline of results. We consider the general model from (1.8), and make the following assumptions: (i) r is bounded; (ii) $\Sigma - I_M$ has bounded rank; (iii) $\log N$ is comparable to $\log M$; (iv) the entries of B are independent, with zero mean and unit variance, and have a sufficient number of bounded moments. Note that, as observed after (1.4), the assumption that the entries of B have zero mean is made for convenience and does not entail a loss in generality. We emphasize that everything apart from r and the rank of $\Sigma - I_M$ is allowed to depend on N in an arbitrary fashion. For definiteness, we explain our results for $Q = \phi^{-1/2} \dot{Q}$, but they also hold verbatim for $\dot{Q} = \phi^{-1/2} Q$.

We establish results on the eigenvalues μ_i and the eigenvectors ξ_i of Q . Our results consist of large deviation bounds and asymptotic laws. We believe that all of our large deviation bounds from Theorems 2.3, 2.6, 2.9, 2.14, and 2.15 are optimal (up to the technical conditions in the definition of $\prec$ given in Definition 2.1). We do not prove this. However, we expect that, combining our method with the techniques of [26], one may also derive the asymptotic laws of all quantities on which we establish large deviation bounds, in particular proving the optimality of our large deviation bounds.

Our results on the eigenvalues of Q consist of two parts. First, we derive large deviation bounds on the locations of the outliers (Theorem 2.3). Second, we prove *eigenvalue sticking* for the non-outliers (Theorem 2.6), whereby each non-outlier “sticks” with high probability and very accurately to the eigenvalues of a related covariance matrix satisfying $\Sigma = I_M$ and whose top eigenvalues exhibit universality. As a corollary (Remark 8.7), we prove that the top non-outlier eigenvalue of Q has asymptotically the Tracy-Widom-1 distribution. This sticking is very accurate if all d_i ’s are separated from the critical point 1, and becomes less accurate if a d_i is in the vicinity of 1. Eventually, it breaks down precisely on the BBP transition scale $|d_i - 1| \sim K^{-1/3}$, at which the Tracy-Widom-1 distribution is known not to hold for the top non-outlier eigenvalue. These results generalize those from [25, Theorem 2.7].

Next, we outline our results for the eigenvectors ξ_i of Q . We consider the *generalized components* $\langle \mathbf{w}, \xi_i \rangle$ of ξ_i , where $\mathbf{w} \in \mathbb{R}^M$ is an arbitrary deterministic vector. In our first result on the eigenvectors (Theorems 2.9 and 2.14), we establish large deviation bounds on the generalized components of outlier eigenvectors (and, more generally, of the outlier spectral projections defined in (2.16) below). This result gives a quantitative version of the cone concentration from (1.14), which in particular allows us to track the strength of the concentration in the vicinity of the BBP transition and for overlapping outliers. Our results also establish the complete delocalization of an outlier eigenvector ξ_i in any direction orthogonal to the spike direction $\mathbf{v}_i$, provided the outlier μ_i is well separated from the bulk spectrum and other outliers. We say that the vector ξ_i is *completely delocalized, or unbiased, in the direction $\mathbf{w}$* if $\langle \mathbf{w}, \xi_i \rangle^2 \prec M^{-1}$, where “ $\prec$ ” denotes a high probability bound up to powers of M^ε (see Definition 2.1).

If the outlier μ_i approaches the bulk spectrum or another outlier, the cone concentration becomes less accurate. For the case of two nearby outlier eigenvalues, for instance, the cone concentration (1.14) of the eigenvectors breaks down when the distributions of the outlier eigenvalues have a nontrivial overlap. In order to understand this behaviour in more detail, we introduce the deterministic projection

$$\Pi_A := \sum_{i \in A} \mathbf{v}_i \mathbf{v}_i^*, \quad (1.16)$$

where $A \subset \{1, \dots, M\}$. Then the cone concentration from (1.14) may be written as $|\Pi_{\{i\}} \xi_i|^2 \approx u(d_i) |\xi_i|^2$. In contrast, in the degenerate case $d_1 = d_2 > 1$ and all other d_i ’s being zero, (1.14) is replaced with

$$\langle \xi_i, \Pi_{\{1,2\}} \xi_j \rangle \approx \delta_{ij} u(d_1) |\xi_i| |\xi_j|, \quad (1.17)$$

where $i, j \in \{1, 2\}$. We deduce that each ξ_i lies on the cone

$$|\Pi_{\{1,2\}}\xi_i|^2 \approx u(d_1)|\xi_i|^2, \quad (1.18)$$

and that $\Pi_{\{1,2\}}\xi_1 \perp \Pi_{\{1,2\}}\xi_2$. Moreover, we prove that ξ_i is completely delocalized in any direction orthogonal to $\mathbf{v}_1$ and $\mathbf{v}_2$. The interpretation is that ξ_1 and ξ_2 both lie on the cone (1.18), that they are orthogonal on both the range and null space of $\Pi_{\{1,2\}}$, and that beyond these constraints their distribution is unbiased (i.e. isotropic). Finally, we note that the preceding discussion remains unchanged if one interchanges ξ_i and $\mathbf{v}_i$. We refer to Example 2.13 below for more details.

In our second result on the eigenvectors (Theorem 2.15), we establish delocalization bounds for the generalized components of non-outlier eigenvectors ξ_i . In particular, we prove complete delocalization of non-outlier eigenvectors in directions orthogonal to any spike $\mathbf{v}_j$ whose value d_j is near the critical point 1. In addition, we prove that non-outlier eigenvectors away from the edge are completely delocalized in all directions. The complete delocalization in the direction $\mathbf{v}_j$ breaks down if $|d_j - 1| \ll 1$. The interpretation of this result is that any spike d_j near the BBP transition point 1 causes all non-outlier eigenvectors ξ_i near the upper edge of the bulk spectrum to have a bias in the direction $\mathbf{v}_j$, in contrast to the completely delocalized case where ξ_i is uniformly distributed on the unit sphere.

In our final result on the eigenvectors (Theorem 2.18), we give the asymptotic law of the generalized component $\langle \mathbf{w}, \xi_i \rangle$ of a non-outlier eigenvector ξ_i . In particular, we prove that this generalized component is asymptotically Gaussian and has a variance predicted by the delocalization bounds from Theorem 2.15. For instance, we prove that if $|d_j - 1| \gg K^{-1/3}$ then

$$\langle \mathbf{v}_j, \xi_i \rangle^2 = \frac{\sigma_i}{M(d_j - 1)^2} \Theta, \quad (1.19)$$

for all non-outlier indices i that are not too large (see Theorem 2.18 for a precise statement). Here Θ is a random variable that converges in distribution to a chi-squared variable. If ξ_i were completely delocalized in the direction $\mathbf{v}_j$, the right-hand side would be of order M^{-1} . Suppose for simplicity that ϕ is of order one. The bias of ξ_i in the direction $\mathbf{v}_j$ emerges as soon as $|d_j - 1| \ll 1$, and reaches a magnitude of order $M^{-1/3}$ for d_j near the BBP transition. This is much larger than the unbiased M^{-1} . Note that this phenomenon applies simultaneously to all non-outlier eigenvectors near the right edge: the right-hand side of (1.19) does not depend on i . Note also that the right-hand side of (1.19) is insensitive to the sign of $d_j - 1$. In particular, the bias is also present for *subcritical* spikes. We conclude that even subcritical spikes are observable in the principal components. In contrast, if one only considers the eigenvalues of the principal components, the subcritical spikes cannot be detected; this follows from the eigenvalue sticking result in Theorem 2.6.

Finally, the proofs of universality of the non-outlier eigenvalues and eigenvectors require the universality of Q for the uncorrelated case $\Sigma = I_M$ as input. This universality result is given in Theorem 8.3, which is also of some independent interest. It establishes the joint, fixed-index, universality of the eigenvalues and eigenvectors of Q (and hence, as a special case, the so-called *quantum unique ergodicity* of the eigenvectors of Q). It works for all eigenvalue indices i satisfying $i \leq K^{1-\tau}$ for any fixed $\tau > 0$.

We conclude this subsection by outlining the key novelties of our work.

- (i) We introduce the general models given in (1.8) and (1.9), which subsume and generalize several models considered previously in the literature¹. We allow the entries of B to be arbitrary random variables (up to a technical assumption on their tails). All quantities except r and the rank of $\Sigma - I_M$ may depend on N . We make no assumption on T beyond the bounded-rank condition of $TT^* - I_M$. The dimensions M and N may be wildly different, and are only subject to the technical condition (2.3).
- (ii) We study the behaviour of the principal components of Q near the BBP transition and when outliers collide. Our results hold for generalized components $\langle \mathbf{w}, \xi_i \rangle$ of the eigenvectors in arbitrary directions $\mathbf{w}$.

¹In particular, the current paper is the first to study the principal components of a realistic sample covariance matrix (1.3) instead of the zero mean case (1.1).

- (iii) We obtain quantitative bounds (i.e. rates of convergence) on the outlier eigenvalues and the generalized components of the eigenvectors. We believe these bounds to be optimal.
- (iv) We obtain precise information about the non-outlier principal components. A novel observation is that, provided there is a d_i satisfying $|d_i - 1| \ll 1$ (i.e. Q is near the BBP transition), all non-outlier eigenvectors near the edge will be biased in the direction of $\mathbf{v}_i$. In particular, non-outlier eigenvectors, unlike non-outlier eigenvalues, retain some information about the subcritical spikes of Σ .
- (v) We establish the joint, fixed-index, universality of the eigenvalues and eigenvectors for the case $\Sigma = I_M$. This result holds for any eigenvalue indices i satisfying $i \leq K^{1-\tau}$ for an arbitrary $\tau > 0$. Note that previous works [27, 38] (established in the context of Wigner matrices) required either the much stronger condition $i \leq (\log K)^{C \log \log K}$ or a four-moment matching condition.

We remark that the large deviation bounds derived in this paper also allow one to derive the joint distribution of the generalized components of the outlier eigenvectors; this will be the subject of future work.

1.6. A few remarks on applications to statistics. We conclude this introduction with a few remarks on what our results imply for applications to statistics. We assume throughout that the population covariance matrix satisfies

$$\max_k |\Sigma_{kk}| \leq C \quad (1.20)$$

for some constant C . We consider the following simple model problem. Suppose there is some (unknown) set $S \subset \{1, \dots, M\}$ whose associated variables $(a_k)_{k \in S}$ are strongly correlated. For simplicity, let us assume that the correlations are given by a single spike in Σ , i.e.

$$\Sigma = 1 + (\sigma - 1)\mathbf{v}\mathbf{v}^*, \quad \sigma - 1 = \phi^{1/2}d,$$

where the spike direction $\mathbf{v} = (v(k))_{k=1}^M$ is given by

$$v(k) := \begin{cases} |S|^{-1/2} & \text{if } k \in S \\ 0 & \text{if } k \notin S. \end{cases}$$

(We choose this precise form for $\mathbf{v}$ so as to simplify the presentation as much as possible. The following discussion also holds if $\mathbf{v}$ is essentially supported on S , but the magnitude of its entries is not necessarily constant.) Moreover, for simplicity we assume that $T = \Sigma^{1/2}$. In components, we have

$$\Sigma_{kl} = \delta_{kl} + (\sigma - 1)v(k)v(l). \quad (1.21)$$

The goal is to recover the set S from an observed realization of the sample covariance matrix $\mathcal{Q}$. (In the more general case where $\mathbf{v}$ is not constant on S , one may easily recover its entries from the submatrix $(\mathcal{Q}_{kl})_{k,l \in S}$ once S has been determined.)

The most naive way to proceed is to compare the entries of $\mathcal{Q}$ with those of Σ . Using (1.20) it is not hard to conclude that

$$\mathcal{Q}_{kl} = \Sigma_{kl} + O_{\prec}(N^{-1/2}),$$

where $O_{\prec}(N^{-1/2})$ denotes a random error term that is bounded with high probability by $N^{\varepsilon-1/2}$ for all $\varepsilon > 0$; see Definition 2.1 below for a precise definition. We look at the off-diagonal terms of $\mathcal{Q}$ and infer that k belongs to S if there exists an index l such that $\mathcal{Q}_{kl}$ is much larger than $N^{-1/2}$. For this approach to work, we require that $|\Sigma_{kl}| \gg N^{-1/2}$, which reads $(\sigma - 1)v(k)v(l) \gg N^{-1/2}$. We conclude that this naive entrywise approach works provided that

$$\sigma - 1 \gg \frac{|S|}{\sqrt{N}}. \quad (1.22)$$

In contrast, according to Theorems 2.3 and 2.9, looking at the principal components of $\mathcal{Q}$ allows us to determine S from the top eigenvector ξ_1 provided the spike σ is supercritical, i.e. gives rise to an outlier. This gives the condition

$$\sigma - 1 \geq \phi^{1/2}. \quad (1.23)$$

Comparing (1.22) and (1.23), we conclude that the principal component analysis works and when the naive componentwise approach does not if we are in the regime

$$|S| \gg \sqrt{M}, \quad \sigma - 1 \geq \phi^{1/2}.$$

Hence the principal component analysis for this example is very effective when the family of correlated variables is quite large, $|S| \gg \sqrt{M}$.

More generally, the principal component approach may work in the regime

$$|S| \gg \phi^{1/2} \quad (1.24)$$

and cannot work for smaller $|S|$. Indeed, the assumption (1.20) is satisfied for $\sigma \leq |S|$, so that in the case of the strongest possible correlations, $\sigma \sim |S|$, the condition (1.23) reduces to (1.24). On the other hand, if $|S| \ll \phi^{1/2}$, the assumption (1.20) implies that $\sigma - 1 \ll \phi^{1/2}$, in contradiction to (1.23).

Clearly, by (1.23), for the purposes of statistical inference it is desirable to make ϕ as small as possible. The interpretation is simply of having more samples per variable. It is therefore natural to attempt to make M smaller so as to reduce ϕ . Obviously, if we know a priori that S is contained in some subset R of $\{1, \dots, M\}$ of size $M/2$, then we simply discard all variables indexed by R^c and consider the correlations restricted to $(a_i)_{i \in R}$; we have halved ϕ in the process.

However, if we have no such a priori knowledge about S , discarding half of the variables a_i is a bad idea. In this case, the best one can do is to choose R at random. Thus, suppose that S is uniformly distributed among the subsets of $\{1, \dots, M\}$ of size $|S|$. We cut the sample space $\{1, \dots, M\}$ in half by keeping only the $M/2$ first elements. We therefore obtain a new family of variables with dimensional parameters

$$\widetilde{M} := M/2, \quad \widetilde{\phi} := \phi/2.$$

Let $\widetilde{\Sigma}$ be the $M/2 \times M/2$ matrix obtained from $\Sigma = 1 + \phi^{1/2} d \mathbf{v} \mathbf{v}^*$ by restricting it to the first $M/2$ elements, i.e. $\widetilde{\Sigma} := (\Sigma_{kl})_{k,l=1}^{M/2}$. Note that $\widetilde{\Sigma}$ is again a rank-one perturbation of the identity. We write it in the form

$$\widetilde{\Sigma} = 1 + \widetilde{\phi}^{1/2} d \widetilde{\mathbf{v}} \widetilde{\mathbf{v}}^*,$$

where $\widetilde{\mathbf{v}}$ is a unit vector. Since

$$|\{k : \widetilde{v}(k) \neq 0\}| \approx \frac{|S|}{2}$$

with high probability, we find that $\widetilde{v}(k) \approx \sqrt{2}v(k)$ for $k \in S$. Picking an entry $\Sigma_{kl} = \widetilde{\Sigma}_{kl}$ for $k, l \in S \setminus \{1, \dots, M/2\}$, we obtain

$$\phi^{1/2} d v(k) v(l) = \widetilde{\phi}^{1/2} d \widetilde{v}(k) \widetilde{v}(l),$$

we therefore find

$$\widetilde{d} \approx d/\sqrt{2}.$$

Hence, detecting spikes in the new problem is *more difficult* than in the original problem, and the halving of sample space is therefore counterproductive unless one has some good a priori information about Σ .

We conclude this discussion by referring to Remark 2.21 below for some further comments on using the *non-outlier* principal components for statistical inference.

Conventions. We use C to denote a generic large positive constant, which may depend on some fixed parameters and whose value may change from one expression to the next. Similarly, we use c to denote a generic small positive constant. For two positive quantities A_N and B_N depending on N we use the notation $A_N \asymp B_N$ to mean $C^{-1}A_N \leq B_N \leq CA_N$ for some positive constant C . For $a < b$ we set $\llbracket a, b \rrbracket := [a, b] \cap \mathbb{Z}$. We use the notation $\mathbf{v} = (v(i))_{i=1}^M$ for vectors in $\mathbb{R}^M$, and denote by $|\cdot| = \|\cdot\|_2$ the Euclidean norm of vectors and by $\|\cdot\|$ the corresponding operator norm of matrices. We use I_M to denote the $M \times M$ identity matrix, which we also sometimes write simply as 1 when there is no risk of confusion.

2. Model and results

2.1. Model. In this section we give the precise definition of our model and introduce some basic notations. Fix a constant $r = 0, 1, 2, 3, \dots$. We consider the $M \times M$ rescaled sample covariance matrix from (1.10) and (1.8),

$$Q := TXX^*T^*, \quad (2.1)$$

as well as the rescaled version of (1.3), defined by

$$\dot{Q} := \frac{N}{N-1}TX(I_N - \mathbf{e}\mathbf{e}^*)X^*T^*; \quad (2.2)$$

here the vector $\mathbf{e}$ was defined in (1.4). (Note that we replace the random matrix B from (1.8) with its rescaled version $X = (MN)^{-1/4}B$.) In these definitions, X is an $(M+r) \times N$ random matrix and T an $M \times (M+r)$ deterministic matrix. For definiteness, and bearing the motivation of sample covariance matrices from Section 1 in mind, we assume that the entries of X and T are real. However, our method also trivially applies to complex-valued X and T , with merely cosmetic changes to the proofs.

We always regard N as the fundamental large parameter, and write $M \equiv M_N$. Here, and throughout the following, in order to unclutter notation we omit the argument N in quantities, such as M , that depend on it. In other words, every symbol that is not explicitly a constant is in fact a sequence indexed by N . We assume that M and N satisfy the bounds

$$N^{1/C} \leq M \leq N^C \quad (2.3)$$

for some positive constant C . We recall the dimensional ratio $\phi := M/N$ from (1.5), and emphasize that it may depend on N and need not converge in $(0, \infty)$; in fact, as follows from (2.3), the only restriction on ϕ is $N^{1/C'} \leq \phi \leq N^{C'}$ for some constant $C' > 0$.

We assume that the entries $X_{i\mu}$ of the $M \times N$ matrix X are independent (but not necessarily identically distributed) random variables satisfying

$$\mathbb{E}X_{i\mu} = 0, \quad \mathbb{E}X_{i\mu}^2 = \frac{1}{\sqrt{NM}}. \quad (2.4)$$

In addition, we assume that, for all $p \in \mathbb{N}$, the random variables $(NM)^{1/4}X_{i\mu}$ have a uniformly bounded p -th moment. In other words, we assume that there is a constant C_p such that

$$\mathbb{E}|(NM)^{1/4}X_{i\mu}|^p \leq C_p. \quad (2.5)$$

The assumption that (2.5) hold for all $p \in \mathbb{N}$ may be easily relaxed. For instance, it is easy to check that our results and their proofs remain valid, after minor adjustments, if we only require that (2.5) holds for all $p \leq C$ for some large enough constant C . We do not pursue such generalizations further.

Recall the definition (1.12) of the the population covariance matrix $\Sigma := TT^*$ associated with T . We suppose that Σ is positive definite, so that each d_i lies in the interval

$$\mathcal{D} := (-\phi^{-1/2}, \infty). \quad (2.6)$$

Our only nontrivial assumption on T is that $\Sigma - I_M$ have bounded rank. Thus, we assume that all but a finite, fixed, number of d_i 's are nonzero. We introduce the index set $R := \{i : d_i \neq 0\}$ of the nontrivial values of d_i . We therefore assume that $|R| = O(1)$.

The following notion of a high-probability bound was introduced in [16], and has been subsequently used in a number of works on random matrix theory. It provides a simple way of systematizing and making precise statements of the form “ A is bounded with high probability by B up to small powers of N ”.

DEFINITION 2.1 (STOCHASTIC DOMINATION). *Let*

$$A = (A^{(N)}(u) : N \in \mathbb{N}, u \in U^{(N)}), \quad B = (B^{(N)}(u) : N \in \mathbb{N}, u \in U^{(N)})$$

be two families of nonnegative random variables, where $U^{(N)}$ is a possibly N -dependent parameter set. We say that A is stochastically dominated by B , uniformly in u , if for all (small) $\varepsilon > 0$ and (large) $D > 0$ we have

$$\sup_{u \in U^{(N)}} \mathbb{P} \left[A^{(N)}(u) > N^\varepsilon B^{(N)}(u) \right] \leq N^{-D}$$

for large enough $N \geq N_0(\varepsilon, D)$. Throughout this paper the stochastic domination will always be uniform in all parameters (such as matrix indices) that are not explicitly fixed. Note that $N_0(\varepsilon, D)$ may depend on the constants from (2.3) and (2.5) as well as any constants fixed in the assumptions of our main results. If A is stochastically dominated by B , uniformly in u , we use the notation $A \prec B$. Moreover, if for some complex family A we have $|A| \prec B$ we also write $A = O_{\prec}(B)$.

REMARK 2.2. Because of (2.3), all (or some) factors of N in Definition (2.1) could be replaced with M without changing the definition of stochastic domination.

The remainder of this section is devoted to main results of this paper. As explained in Section 1, we first state our results for the matrix Q from (2.1), bearing in mind that they also hold for $\dot{Q}$ from (2.2). The precise statement for $\dot{Q}$ is given in Theorem 2.22 below.

2.2. Eigenvalue locations. We begin with results on the locations of the eigenvalues of Q . These results will also serve as a fundamental input for the proofs of the results on eigenvectors presented in Sections 2.3 and 2.4.

Recall that Q has $M - K$ zero eigenvalues. We shall therefore focus on the K nontrivial eigenvalues $\mu_1 \geq \dots \geq \mu_K$ of Q . On the global scale, the eigenvalues of Q are distributed according to the Marchenko-Pastur law (1.6). This may be easily inferred from the fact that (1.6) gives the global density of the eigenvalues for the uncorrelated case $\Sigma = I_M$, combined with eigenvalue interlacing (see Lemma 4.1). In this section we focus on *local* eigenvalue information.

We introduce the set of *outlier indices*

$$\mathcal{O} := \{i \in R : |d_i| \geq 1 + K^{-1/3}\}. \quad (2.7)$$

As explained in Section 1.3, each $i \in \mathcal{O}$ gives rise to an outlier of Q near the classical location $\theta(d_i)$ defined in (1.13). In the definition (2.7), the lower bound $1 + K^{-1/3}$ is chosen for definiteness; it could be replaced with $1 + aK^{-1/3}$ for any fixed $a > 0$. We denote by

$$s_{\pm} := |\{i \in \mathcal{O} : \pm d_i > 0\}| \quad (2.8)$$

the number of outliers to the left (s_-) and right (s_+) of the bulk spectrum.

For $d \in \mathcal{D} \setminus [-1, 1]$ we define

$$\Delta(d) := \begin{cases} \frac{\phi^{1/2}\theta(d)}{1+(|d|-1)^{-1/2}} & \text{if } -\phi^{-1/2} < d < -1 \\ (d-1)^{1/2} & \text{if } 1 < d \leq 2 \\ 1 + \frac{d}{1+\phi^{-1/2}} & \text{if } d \geq 2. \end{cases}$$

The function $\Delta(d)$ will be used to give an upper bound on the magnitude of the fluctuations of an outlier associated with d . We give such a precise expression for Δ in order to obtain sharp large deviation bounds for all $d \in \mathcal{D} \setminus [-1, 1]$. (Note that the discontinuity of Δ at $d = 2$ is immaterial since Δ is used as an upper bound with respect to $\prec$. The ratio of the right- and left-sided limits at 2 of Δ lies in $[1, 3]$.)

Our result on the outlier eigenvalues is the following.

THEOREM 2.3 (OUTLIER LOCATIONS). *Fix $\tau > 0$. Then for $i \in \mathcal{O}$ we have the estimate*

$$|\mu_i - \theta(d_i)| \prec \Delta(d_i) K^{-1/2} \quad (2.9)$$

provided that $d_i > 0$ or $|\phi - 1| \geq \tau$.

Furthermore, the extremal non-outliers μ_{s_++1} and μ_{K-s_-} satisfy

$$|\mu_{s_++1} - \gamma_+| \prec K^{-2/3}, \quad (2.10)$$

and, assuming in addition that $|\phi - 1| \geq \tau$,

$$|\mu_{K-s_-} - \gamma_-| \prec K^{-2/3}. \quad (2.11)$$

REMARK 2.4. Theorem 2.3 gives large deviation bounds for the locations of the outliers to the right of the bulk. Since $\tau > 0$ may be arbitrarily small, Theorem 2.3 also gives the full information about the outliers to the left of the bulk except in the case $1 > \phi = 1 + o(1)$. Although our methods may be extended to this case as well, we exclude it here to avoid extraneous complications.

REMARK 2.5. By definition of s_- and $\mathcal{D}$, if $\phi > 1$ then $s_- = 0$. Hence, by (2.11), if $\phi > 1$ there are no outliers on the left of the bulk spectrum.

The locations of the non-outlier eigenvalues μ_i , $i \notin \mathcal{O}$, are governed by *eigenvalue sticking*, whereby the eigenvalues of Q “stick” with high probability to eigenvalues of a reference matrix which has a trivial population covariance matrix. The reference matrix is Q from (2.1) with uncorrelated entries. More precisely, we set

$$H := YY^*, \quad Y := (I_M, 0)OX, \quad (2.12)$$

$O \equiv O(T) \in O(M+r)$ is a deterministic orthogonal matrix. It is easy to check that $\mathbb{E}H = (NM)^{-1/2}I_M$, so that H corresponds to an uncorrelated population. The matrix $O(T)$ is explicitly given in (8.1) below. In fact, in Theorem 8.3 below we prove the universality of the joint distribution of non-bulk eigenvalues and eigenvectors of H . Here, by definition, we say that an index $i \in \llbracket 1, K \rrbracket$ is *non-bulk* if $i \notin \llbracket K^{1-\tau}, K - K^{1-\tau} \rrbracket$ for some fixed $\tau > 0$. In particular, the asymptotic distribution of the non-bulk eigenvalues and eigenvectors of H does not depend on the choice of O . Note that for the special case $r = 0$ the eigenvalues of H coincide with those of XX^* . We denote by

$$\lambda_1 \geq \lambda_2 \geq \dots \geq \lambda_M$$

the eigenvalues of H .

THEOREM 2.6 (EIGENVALUE STICKING). *Define*

$$\alpha_{\pm} := \min_{1 \leq i \leq M} |d_i \mp 1|. \quad (2.13)$$

Fix $\tau > 0$. Then we have for all $i \in \llbracket 1, (1-\tau)K \rrbracket$

$$|\mu_{i+s_+} - \lambda_i| \prec \frac{1}{K\alpha_+}. \quad (2.14)$$

Similarly, if $|\phi - 1| \geq \tau$ then we have for all $i \in \llbracket \tau K, K \rrbracket$

$$|\mu_{i-s_-} - \lambda_i| \prec \frac{1}{K\alpha_-}. \quad (2.15)$$

REMARK 2.7. As outlined above, in Theorem 8.3 below we prove that the asymptotic joint distribution of the non-bulk eigenvalues of H is universal, i.e. it coincides with that of the Wishart matrix $H_{\text{Wish}} = XX^*$ with $r = 0$ and X Gaussian. As an immediate corollary of Theorems 2.6 and 8.3, we obtain the universality of the non-outlier eigenvalues of Q with index $i \leq K^{1-\tau}\alpha_+^3$. This condition states simply that the right-hand side of (2.14) is much smaller than the scale on which the eigenvalue λ_i fluctuates, which is $K^{-2/3}i^{-1/3}$. See Remark 8.7 below for a precise statement.

REMARK 2.8. Theorem 2.6 is analogous to Theorem 2.7 of [25], where sticking was first established for Wigner matrices. Previously, eigenvalue sticking was established for a certain class of random perturbations of Wigner matrices in [5, 6]. We refer to [25, Remark 2.8] for a more detailed discussion.

Aside from holding for general covariance matrices of the form (2.1), Theorem 2.6 is stronger than its counterpart from [25] because it holds much further into the bulk: in [25, Theorem 2.7], sticking was established under the assumption that $i \leq (\log K)^{C \log \log K}$.

2.3. Outlier eigenvectors. We now state our main results for the outlier eigenvectors. Statements of results on eigenvectors requires some care, since there is some arbitrariness in the definition of the eigenvector ξ_i . In order to get rid of the arbitrariness in the sign (or, in the complex case, the phase) of ξ_i we consider products of generalized components,

$$\langle \mathbf{v}, \xi_i \rangle \langle \xi_i, \mathbf{w} \rangle.$$

It is easy to check that these products characterize the eigenvector ξ_i completely, up to the ambiguity of a global sign (or phase). More generally, one may consider the generalized components $\langle \mathbf{v}, (\cdot) \mathbf{w} \rangle$ of the (random) *spectral projection*

$$P_A := \sum_{i \in A} \xi_i \xi_i^*, \quad (2.16)$$

where $A \subset \mathcal{O}$.

In the simplest case $A = \{i\}$, and the generalized components of P_A characterize the generalized components of ξ_i . The need to consider higher-dimensional projections arises if one considers degenerate or almost degenerate outliers. Suppose for example that $d_1 \approx d_2$ and all other d_i 's are zero. Then the cone concentration (1.14) fails, to be replaced with (1.17). The failure of the cone concentration is also visible in our results as a blowup of the error bounds. This behaviour is not surprising, since for degenerate outliers $d_1 = d_2$ it makes no sense to distinguish the associated spike eigenvectors $\mathbf{v}_1$ and $\mathbf{v}_2$; only the eigenspace matters. Correspondingly, we have to consider the orthogonal projection onto the eigenspace of the outliers in A . See Example 2.13 below for a more detailed discussion.

For $i \in \llbracket 1, M \rrbracket$ we define $\nu_i \geq 0$ through

$$\nu_i \equiv \nu_i(A) := \begin{cases} \min_{j \notin A} |d_i - d_j| & \text{if } i \in A \\ \min_{j \in A} |d_i - d_j| & \text{if } i \notin A. \end{cases}$$

In other words, $\nu_i(A)$ is the distance from d_i to either $\{d_i\}_{i \in A}$ or $\{d_i\}_{i \notin A}$, whichever it does not belong to. For a vector $\mathbf{w} \in \mathbb{R}^M$ we also introduce the shorthand

$$w_i := \langle \mathbf{v}_i, \mathbf{w} \rangle$$

to denote the components of $\mathbf{w}$ in the eigenbasis of Σ .

For definiteness, we only state our results for the outliers on the right-hand side of the bulk spectrum. Analogous results hold for the outliers on the left-hand side. Since the behaviour of the fluctuating error term is different in the regimes $\mu_i - \gamma_+ \ll 1$ (near the bulk) and $\mu_i - \gamma_+ \gg 1$ (far from the bulk), we split these two cases into separate theorems.

THEOREM 2.9 (OUTLIER EIGENVECTORS NEAR BULK). *Fix $\tau > 0$. Suppose that $A \subset \mathcal{O}$ satisfies $1 + K^{-1/3} \leq d_i \leq \tau^{-1}$ for all $i \in A$. Define the deterministic positive quadratic form*

$$\langle \mathbf{w}, Z_A \mathbf{w} \rangle := \sum_{i \in A} u(d_i) w_i^2,$$

where we recall the definition (1.15) of $u(d_i)$. Then for any deterministic $\mathbf{w} \in \mathbb{R}^M$ we have

$$\langle \mathbf{w}, P_A \mathbf{w} \rangle = \langle \mathbf{w}, Z_A \mathbf{w} \rangle + O_{\prec} \left[\sum_{i \in A} \frac{w_i^2}{M^{1/2}(d_i - 1)^{1/2}} + \sum_{i=1}^M \frac{\sigma_i w_i^2}{M \nu_i(A)^2} + \langle \mathbf{w}, Z_A \mathbf{w} \rangle^{1/2} \left(\sum_{i \notin A} \frac{\sigma_i w_i^2}{M \nu_i(A)^2} \right)^{1/2} \right]. \quad (2.17)$$

Note that the last error term is zero if $\mathbf{w}$ is in the subspace $\text{Span}\{\mathbf{v}_i\}_{i \in A}$ or orthogonal to it.

REMARK 2.10. Theorem 2.9 may easily also be stated for more general quantities of the form $\langle \mathbf{v}, P_A \mathbf{w} \rangle$. We omit the precise statement; it is a trivial corollary of (5.2) below, which holds under the assumptions of Theorem 2.9.

We emphasize that the set A in Theorem 2.9 may be chosen at will. If all outliers are well-separated, then the choice $A = \{i\}$ gives the most precise information. However, as explained at the beginning of this subsection, the indices of outliers that are close to each other should be included in the same set A . Thus, the freedom to chose $|A| \geq 2$ is meant for degenerate or almost degenerate outliers. (In fact, as explained after (2.19) below, the correct notion of closeness of outliers is that of *overlapping*.)

We consider a few examples.

EXAMPLE 2.11. Let $A = \{i\}$ and $\mathbf{w} = \mathbf{v}_i$. Then we get from (2.17)

$$\langle \mathbf{v}_i, \boldsymbol{\xi}_i \rangle^2 = u(d_i) + O_{\prec} \left[\frac{1}{M^{1/2}(d_i - 1)^{1/2}} + \frac{\sigma_i}{M\nu_i^2} \right]. \quad (2.18)$$

This gives a precise version of the cone concentration from (1.14). Note that the cone concentration holds provided the error is much smaller than the main term $u(d_i)$, which leads to the conditions

$$d_i - 1 \gg K^{-1/3} \quad \text{and} \quad \nu_i \gg (d_i - 1)^{-1/2} K^{-1/2}; \quad (2.19)$$

here we used that $d_i \asymp 1$ and $M \asymp (1 + \phi)K$.

We claim that both conditions in (2.19) are natural and necessary. The first condition of (2.19) simply means that μ_i is an outlier. The second condition of (2.19) is a *non-overlapping* condition. To understand it, recall from (2.9) that μ_i fluctuates on the scale $(d_i - 1)^{1/2} K^{-1/2}$. Then μ_i is a non-overlapping outlier if all other outliers are located with high probability at a distance greater than this scale from μ_i . Recalling the definition of the classical location $\theta(d_i)$ of μ_i , the non-overlapping condition becomes

$$\min_{j \in \mathcal{O} \setminus \{i\}} |\theta(d_j) - \theta(d_i)| \gg (d_i - 1)^{1/2} K^{-1/2}. \quad (2.20)$$

After a simple estimate using the definition of θ , we find that this is precisely the second condition of (2.19). The degeneracy or almost degeneracy of outliers discussed at the beginning of this subsection is hence to be interpreted more precisely in terms of overlapping of outliers.

Provided μ_i is well-separated from both the bulk spectrum and the other outliers, we find that the error in (2.18) is of order $M^{-1/2}$.

EXAMPLE 2.12. Take $A = \{i\}$ and $\mathbf{w} = \mathbf{w}_j$ with $j \neq i$. Then we get from (2.17)

$$\langle \mathbf{v}_j, \boldsymbol{\xi}_i \rangle^2 \prec \frac{\sigma_j}{M(d_i - d_j)^2}. \quad (2.21)$$

Suppose for simplicity that $\phi \asymp 1$. Then, under the condition that $|d_i - d_j| \asymp 1$, we find that $\boldsymbol{\xi}_i$ is completely delocalized in the direction $\mathbf{v}_j$. In particular, if $\nu_i \asymp 1$ then $\boldsymbol{\xi}_i$ is completely delocalized in any direction orthogonal to $\mathbf{v}_i$.

As d_j approaches d_i the delocalization bound from (2.21) deteriorates, and eventually when μ_i and μ_j start overlapping, i.e. the second condition of (2.19) is violated, the right-hand side of (2.21) has the same size as the leading term of (2.18). This is again a manifestation of the fact that the individual eigenspaces of overlapping outliers cannot be distinguished.

EXAMPLE 2.13. Suppose that we have an $|A|$ -fold degenerate outlier, i.e. $d_i = d_j$ for all $i, j \in A$. Then from Theorem 2.9 and Remark 2.10 (see the estimate (5.2)) we get, for all $i, j \in A$,

$$\langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle = \delta_{ij} u(d_i) + O_{\prec} \left[\frac{1}{M^{1/2}(d_i - 1)^{1/2}} + \frac{\sigma_i}{M\nu_i(A)^2} \right].$$

Defining the $|A| \times |A|$ random matrix $M = (M_{ij})_{i,j \in A}$ through $M_{ij} := \langle \mathbf{v}_i, \boldsymbol{\xi}_j \rangle$, we may write the left-hand side as $(MM^*)_{ij}$. We conclude that $u(d_i)^{-1/2} M$ is approximately orthogonal, from which we deduce that

$u(d_i)^{-1/2}M^*$ is also approximately orthogonal. In other words, we may interchange the families $\{\mathbf{v}_i\}_{i \in A}$ and $\{\boldsymbol{\xi}_i\}_{i \in A}$. More precisely, we get

$$\langle \boldsymbol{\xi}_i, \Pi_A \boldsymbol{\xi}_j \rangle = (M^* M)_{ij} = \delta_{ij} u(d_i) + O_{\prec} \left[\frac{1}{M^{1/2}(d_i - 1)^{1/2}} + \frac{\sigma_i}{M \nu_i(A)^2} \right].$$

This is the correct generalization of (2.18) from Example 2.11 to the degenerate case. The error term is the same as in (2.18), and its size and relation to the main term is exactly the same as in Example 2.11. Hence the discussion following (2.18) may be take over verbatim to this case.

In addition, analogously to Example 2.12, for $i \in A$ and $j \notin A$ we find that (2.21) remains true. This establishes the delocalization of $\boldsymbol{\xi}_i$ in any direction within the null space of Π_A .

These estimates establish the general cone concentration, with optimal rate of convergence, for degenerate outliers outlined around (1.18). The eigenvectors $\{\boldsymbol{\xi}_i\}_{i \in A}$ are all concentrated on the cone defined by $|\Pi_A \boldsymbol{\xi}|^2 = u(d_i)|\boldsymbol{\xi}|^2$ (for some immaterial $i \in A$). Moreover, the eigenvectors $\{\boldsymbol{\xi}_i\}_{i \in A}$ are orthogonal on both the range and null space of Π_A . Provided that the group $\{d_i\}_{i \in A}$ is well-separated from 1 and all other d_i 's, the eigenvectors $\{\boldsymbol{\xi}_i\}_{i \in A}$ are completely delocalized on the null space of Π_A .

We conclude this example by remarking that a similar discussion also holds for a group of outliers that is not degenerate, but nearly degenerate, i.e. $|d_i - d_j| \ll |d_i - d_k|$ for all $i, j \in A$ and $k \notin A$. We omit the details.

The next result is the analogue of Theorem 2.9 for outliers far from the bulk.

THEOREM 2.14 (OUTLIER EIGENVECTORS FAR FROM BULK). *Fix $\tau > 0$. Suppose that $A \subset \mathcal{O}$ satisfies $d_i \geq 1 + \tau$ for all $i \in A$, and that there exists a positive d_A such that $\tau d_A \leq d_i \leq \tau^{-1} d_A$ for all $i \in A$. Then for any deterministic $\mathbf{w} \in \mathbb{R}^M$ we have*

$$\begin{aligned} \langle \mathbf{w}, P_A \mathbf{w} \rangle = \langle \mathbf{w}, Z_A \mathbf{w} \rangle + O_{\prec} & \left[\frac{1}{M^{1/2}(\phi^{1/2} + d_A)} \sum_{i \in A} \sigma_i w_i^2 + \left(1 + \frac{\phi^{1/2} d_A^2}{\phi^{1/2} + d_A} \right) \sum_{i=1}^M \frac{\sigma_i w_i^2}{M \nu_i(A)^2} \right. \\ & \left. + \frac{d_A}{\phi^{1/2} + d_A} \left(\sum_{i \in A} \sigma_i w_i^2 \right)^{1/2} \left(\sum_{i \notin A} \frac{\sigma_i w_i^2}{M \nu_i(A)^2} \right)^{1/2} \right]. \end{aligned} \quad (2.22)$$

We leave the discussion on the interpretation of the error in (2.22) to the reader; it is similar to that of Examples 2.11, 2.12, and 2.13.

2.4. Non-outlier eigenvectors. In this subsection we state our results on the *non-outlier* eigenvectors, i.e. on $\boldsymbol{\xi}_a$ for $a \notin \mathcal{O}$. Our first result is a delocalization bound. In order to state it, we define for $a \in \llbracket 1, K \rrbracket$ the *typical distance from μ_a to the spectral edges $\gamma_{\pm}$* through

$$\kappa_a := K^{-2/3} (a \wedge (K + 1 - a))^{2/3}. \quad (2.23)$$

This quantity should be interpreted as a deterministic version of $\min\{|\mu_a - \gamma_-|, |\mu_a - \gamma_+|\}$ for $a \notin \mathcal{O}$; see Theorem 3.5 below.

THEOREM 2.15 (DELOCALIZATION BOUND FOR NON-OUTLIERS). *Fix $\tau > 0$. For $a \in \llbracket 1, (1 - \tau)K \rrbracket \setminus \mathcal{O}$ and deterministic $\mathbf{w} \in \mathbb{R}^M$ we have*

$$\langle \mathbf{w}, \boldsymbol{\xi}_a \rangle^2 \prec \frac{|\mathbf{w}|^2}{M} + \sum_{i=1}^M \frac{\sigma_i w_i^2}{M((d_i - 1)^2 + \kappa_a)}. \quad (2.24)$$

Similarly, if $|\phi - 1| \geq \tau$ then for $a \in \llbracket \tau K, K \rrbracket \setminus \mathcal{O}$ and deterministic $\mathbf{w} \in \mathbb{R}^M$ we have

$$\langle \mathbf{w}, \boldsymbol{\xi}_a \rangle^2 \prec \frac{|\mathbf{w}|^2}{M} + \sum_{i=1}^M \frac{\sigma_i w_i^2}{M((d_i + 1)^2 + \kappa_a)}. \quad (2.25)$$

For the following examples, we take $\mathbf{w} = \mathbf{v}_i$ and $a \in \llbracket 1, (1 - \tau)K \rrbracket \setminus \mathcal{O}$. Under these assumptions (2.24) yields

$$\langle \mathbf{w}, \boldsymbol{\xi}_a \rangle^2 \prec \frac{1}{M} + \frac{\sigma_i}{M((d_i - 1)^2 + \kappa_a)} \quad (2.26)$$

EXAMPLE 2.16. Fix $\tau > 0$. If $|d_i - 1| \geq \tau$ (d_i is separated from the transition point) or $a \geq \tau K$ (μ_a is in the bulk), then the right-hand side of (2.26) reads $(1 + \sigma_i)/M$. In particular, if the eigenvalue σ_i of Σ is bounded, $\boldsymbol{\xi}_a$ is completely delocalized in the direction $\mathbf{v}_i$.

EXAMPLE 2.17. Suppose that $a \leq C$ (μ_a is near the edge), which implies that $\kappa_a \asymp K^{-2/3}$. Suppose moreover that d_i is near the transition point 1. Then we get

$$\langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2 \prec \frac{\sigma_i}{M((d_i - 1)^2 + K^{-2/3})}.$$

Therefore the delocalization bound for $\boldsymbol{\xi}_a$ in the direction of $\mathbf{v}_i$ becomes worse as d_i approaches the critical point (from either side), from $(1 + \phi)^{1/2} M^{-1}$ for d_i separated from 1, to $(1 + \phi)^{-1/6} M^{-1/3}$ for d_i at the transition point 1.

Next, we derive the law of the generalized component $\langle \mathbf{w}, \boldsymbol{\xi}_a \rangle$ for non-outlier a . In particular, this provides a lower bound complementing the upper bound from Theorem 2.15. Recall the definition (2.13) of α_+ .

THEOREM 2.18 (LAW OF NON-OUTLIERS). *Fix $\tau > 0$. Then, for any deterministic $a \in \llbracket 1, K^{1-\tau} \alpha_+^3 \rrbracket \setminus \mathcal{O}$ and $\mathbf{w} \in \mathbb{R}^M$, there exists a random variable $\Theta(a, \mathbf{w}) \equiv \Theta_N(a, \mathbf{w})$ satisfying*

$$\langle \mathbf{w}, \boldsymbol{\xi}_a \rangle^2 = \sum_i \frac{\sigma_i w_i^2}{M(d_i - 1)^2} \Theta(a, \mathbf{w}).$$

and

$$\Theta(a, \mathbf{w}) \longrightarrow \chi_1^2$$

in distribution, uniformly in a and $\mathbf{w}$. Here χ_1^2 is a chi-squared random variable (i.e. the square of a standard normal).

An analogous statement holds near the left spectral edge provided $|\phi - 1| \geq \tau$; we omit the details.

REMARK 2.19. More generally, our method also yields the asymptotic joint distribution of the family

$$\left(\mu_{a_1}, \dots, \mu_{a_k}, \langle \mathbf{u}_1, \boldsymbol{\xi}_{b_1} \rangle \langle \boldsymbol{\xi}_{b_1}, \mathbf{w}_1 \rangle, \dots, \langle \mathbf{u}_k, \boldsymbol{\xi}_{b_k} \rangle \langle \boldsymbol{\xi}_{b_k}, \mathbf{w}_k \rangle \right) \quad (2.27)$$

(after a suitable affine rescaling of the variables, as in Theorem (8.3) below), where $a_1, \dots, a_k, b_1, \dots, b_k \in \llbracket 1, K^{1-\tau} \alpha_+^3 \rrbracket \setminus \mathcal{O}$. We omit the precise statement, which is a universality result: it says essentially that the asymptotic distribution of (2.27) coincides with that under the standard Wishart ensemble (i.e. an uncorrelated Gaussian sample covariance matrix). The proof is a simple corollary of Theorem 2.6, Proposition 6.3, Proposition 6.4, and Theorem 8.3.

REMARK 2.20. The restriction $a \leq K^{1-\tau} \alpha_+^3$ is the same as in Remarks 2.7 and 8.7. There, it is required for the eigenvalue sticking to be effective in the sense that the right-hand side of (2.14) is much smaller than the scale on which the eigenvalue λ_a fluctuates. Here, it ensures the distribution of the eigenvector $\boldsymbol{\xi}_a$ is determined by the distribution of a single eigenvector of H (see Proposition 6.3).

REMARK 2.21. As outlined in the introduction, Theorem 2.18 implies that the non-outlier eigenvectors near the edge are all biased in the direction of $\mathbf{v}_i$ provided that d_i is near the BBP transition point 1. Suppose for simplicity that ϕ is bounded. Then Theorem 2.18 implies that for d_i near 1 we have $\langle \mathbf{w}, \boldsymbol{\xi}_a \rangle^2 \asymp (d_i - 1)^{-2} M^{-1}$ with high probability. We can therefore detect the spike d_i through the resulting bias in the direction $\mathbf{v}_i$ as long as $|d_i - 1| \ll 1$. (Recall that the unbiased, or completely delocalized, case corresponds to $\langle \mathbf{w}, \boldsymbol{\xi}_a \rangle^2 \asymp M^{-1}$.) Hence, the non-outlier eigenvectors $\boldsymbol{\xi}_a$ retain information even about

a *subcritical* spikes. This is in stark contrast to the eigenvalues μ_a , which, by Theorem 2.6, retain no information about the subcritical spikes.

If $d_i \leq 1 - \tau$ for some fixed τ , then the principal components of Q contain no information about the spike d_i . We illustrate this using the simple statistical problem from Section 1.6, whose notations we take over without further comment. Suppose we try to determine the set S by choosing the components of ξ_a that are much larger than the unbiased background value $M^{-1/2}$. This works provided that for $k \in S$ we have

$$\xi_a(k)^2 \asymp \frac{(1 + \phi)^{1/2} v(k)^2}{M} \gg \frac{1}{M}.$$

Using $v(k)^2 = |S|^{-1}$ for $k \in S$, we therefore get the condition $(1 + \phi)^{1/2} \gg |S|$. This however cannot hold, since we have by assumption (1.20) on Σ that

$$C \geq \Sigma_{kk} \asymp 1 + \phi^{1/2} |S|^{-1}.$$

Finally, we note that all of our results also hold for $\dot{Q}$ instead of Q .

THEOREM 2.22. *Theorems 2.3, 2.6, 2.9, 2.14, 2.15, and 2.18 hold with μ_i and ξ_i denoting the eigenvalues and eigenvectors of $\dot{Q}$ instead of Q . For Theorem 2.6, λ_i denotes the eigenvalues of $\frac{N}{N-1} Y(I_N - \mathbf{e}\mathbf{e}^*) Y^*$ instead of YY^* from (2.12).*

3. Preliminaries

The rest of this paper is devoted to the proofs of the results from Sections 2.2–2.4. To clarify the presentation of the main ideas of the proofs, we shall first assume that

$$r = 0 \quad \text{and} \quad T = \Sigma^{1/2}. \quad (3.1)$$

We make the assumption (3.1) throughout Sections 3–7. The additional arguments required to relax the assumption (3.1) are presented in Section 8. Under the assumption (3.1) we have

$$Q = \Sigma^{1/2} H \Sigma^{1/2}, \quad H = X X^*. \quad (3.2)$$

Moreover, the extension of our results from Q to $\dot{Q}$, and hence the proof of Theorem 2.22, is given in Section 9.

For an $M \times M$ matrix A and $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$ we abbreviate

$$A_{\mathbf{v}\mathbf{w}} := \langle \mathbf{v}, A\mathbf{w} \rangle.$$

We also write

$$A_{\mathbf{v}\mathbf{e}_i} \equiv A_{\mathbf{v}i}, \quad A_{\mathbf{e}_i\mathbf{v}} \equiv A_{i\mathbf{v}}, \quad A_{\mathbf{e}_i\mathbf{e}_j} \equiv A_{ij},$$

where $\mathbf{e}_i \in \mathbb{R}^M$ denotes the i -th standard basis vector.

3.1. The isotropic local Marchenko-Pastur law. In this section we collect the key tool of our analysis: the isotropic Marchenko-Pastur law from [9].

It is well known that the empirical distribution of the eigenvalues of the $N \times N$ matrix X^*X has the same asymptotics as the *Marchenko-Pastur* law

$$\varrho_\phi(dx) := \frac{\sqrt{\phi}}{2\pi} \frac{\sqrt{[(x - \gamma_-)(\gamma_+ - x)]_+}}{x} dx + (1 - \phi)_+ \delta(dx), \quad (3.3)$$

where we recall the edges $\gamma_\pm$ of the limiting spectrum defined in (1.7). Similarly, as noted in (1.6), the empirical distribution of the eigenvalues of the $M \times M$ matrix XX^* has the same asymptotics as $\varrho_{\phi^{-1}}$.

Note that (3.3) is normalized so that its integral is equal to one. The Stieltjes transform of the Marchenko-Pastur law (3.3) is

$$m_\phi(z) := \int \frac{\varrho_\phi(dx)}{x-z} = \frac{\phi^{1/2} - \phi^{-1/2} - z + i\sqrt{(z-\gamma_-)(\gamma_+ - z)}}{2\phi^{-1/2}z}, \quad (3.4)$$

where the square root is chosen so that m_ϕ is holomorphic in the upper half-plane and satisfies $m_\phi(z) \rightarrow 0$ as $z \rightarrow \infty$. The function $m_\phi = m_\phi(z)$ is also characterized as the unique solution of the equation

$$m + \frac{1}{z + z\phi^{-1/2}m - (\phi^{1/2} - \phi^{-1/2})} = 0 \quad (3.5)$$

satisfying $\text{Im } m(z) > 0$ for $\text{Im } z > 0$. The formulas (3.3)–(3.5) were originally derived for the case when $\phi = M/N$ is independent of N (or, more precisely, when ϕ has a limit in $(0, \infty)$ as $N \rightarrow \infty$). Our results allow ϕ to depend on N under the constraint (2.3), so that m_ϕ and ϱ_ϕ may also depend on N through ϕ .

Throughout the following we use a spectral parameter

$$z = E + i\eta,$$

with $\eta > 0$, as the argument of Stieltjes transforms and resolvents. Define the resolvent

$$G(z) := (XX^* - z)^{-1}. \quad (3.6)$$

For $z \in \mathbb{C}$, define $\kappa(z)$ to be the distance of $E = \text{Re } z$ to the spectral edges $\gamma_\pm$, i.e.

$$\kappa \equiv \kappa(z) := \min\{|\gamma_+ - E|, |\gamma_- - E|\}. \quad (3.7)$$

Throughout the following we regard the quantities $E(z)$, $\eta(z)$, and $\kappa(z)$ as functions of z and usually omit the argument unless it is needed to avoid confusion.

Sometimes we shall need the following notion of high probability.

DEFINITION 3.1. *An N -dependent event $\Xi \equiv \Xi_N$ holds with high probability if $1 - \mathbf{1}(\Xi) \prec 0$.*

Fix a (small) $\omega \in (0, 1)$ and define the domain

$$\mathbf{S} \equiv \mathbf{S}(\omega, K) := \{z \in \mathbb{C} : \kappa \leq \omega^{-1}, K^{-1+\omega} \leq \eta \leq \omega^{-1}, |z| \geq \omega\}. \quad (3.8)$$

Beyond the support of the limiting spectrum, one has stronger control all the way down to the real axis. For fixed (small) $\omega > 0$ define the region

$$\tilde{\mathbf{S}} \equiv \tilde{\mathbf{S}}(\omega, K) := \{z \in \mathbb{C} : E \notin [\gamma_-, \gamma_+], K^{-2/3+\omega} \leq \kappa \leq \omega^{-1}, |z| \geq \omega, 0 < \eta \leq \omega^{-1}\} \quad (3.9)$$

of spectral parameters separated from the asymptotic spectrum by $K^{-2/3+\omega}$, which may have an arbitrarily small positive imaginary part η . Throughout the following we regard ω as fixed once and for all, and do not track the dependence of constants on ω .

THEOREM 3.2 (ISOTROPIC LOCAL MARCHENKO-PASTUR LAW [9]). *Suppose that (2.4), (2.3), and (2.5) hold. Then*

$$|\langle \mathbf{v}, G(z)\mathbf{w} \rangle - m_{\phi^{-1}}(z)\langle \mathbf{v}, \mathbf{w} \rangle| \prec \sqrt{\frac{\text{Im } m_{\phi^{-1}}(z)}{M\eta}} + \frac{1}{M\eta} \quad (3.10)$$

uniformly in $z \in \mathbf{S}$ and any deterministic unit vectors $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$. Moreover,

$$|\langle \mathbf{v}, G(z)\mathbf{w} \rangle - m_{\phi^{-1}}(z)\langle \mathbf{v}, \mathbf{w} \rangle| \prec \sqrt{\frac{\text{Im } m_{\phi^{-1}}(z)}{M\eta}} \asymp \frac{1}{1+\phi}(\kappa + \eta)^{-1/4}K^{-1/2} \quad (3.11)$$

uniformly in $z \in \tilde{\mathbf{S}}$ and any deterministic unit vectors $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$.

REMARK 3.3. The probabilistic estimates (3.10) and (3.11) of Theorem 3.2 may be strengthened to hold simultaneously for all $z \in \mathbf{S}$ and for all $z \in \tilde{\mathbf{S}}$, respectively. For instance, (3.11) may be strengthened to

$$\mathbb{P} \left[\bigcap_{z \in \tilde{\mathbf{S}}} \left\{ |\langle \mathbf{v}, G(z) \mathbf{w} \rangle - m_{\phi^{-1}}(z) \langle \mathbf{v}, \mathbf{w} \rangle| \leq N^\varepsilon \frac{1}{1+\phi} (\kappa + \eta)^{-1/4} K^{-1/2} \right\} \right] \geq 1 - N^{-D},$$

for all $\varepsilon > 0$, $D > 0$, and $N \geq N_0(\varepsilon, D)$. See [9, Remark 2.6].

The next results are on the nontrivial (i.e. nonzero) eigenvalues of XX^* as well as the corresponding eigenvectors. The matrix XX^* has K nontrivial eigenvalues, which we order according to $\lambda_1 \geq \lambda_2 \geq \dots \geq \lambda_K$. (The remaining $M - K$ eigenvalues of XX^* are zero.) Moreover, we denote by $\zeta_1, \dots, \zeta_K \in \mathbb{R}^M$ the unit eigenvectors of XX^* associated with the nontrivial eigenvalues $\lambda_1 \geq \dots \geq \lambda_K$.

THEOREM 3.4 (ISOTROPIC DELOCALIZATION [9]). *Fix $\tau > 0$, and suppose that (2.4), (2.3), and (2.5) hold. Then for $i \in \llbracket 1, K \rrbracket$ we have*

$$\langle \zeta_i, \mathbf{v} \rangle^2 \prec M^{-1} \quad (3.12)$$

if either $i \leq (1 - \tau)K$ or $|\phi - 1| \geq \tau$.

The following result is on the rigidity of the nontrivial eigenvalues of XX^* . Let $\gamma_1 \geq \gamma_2 \geq \dots \geq \gamma_K$ be the *classical eigenvalue locations* according to ϱ_ϕ (see (3.3)), defined through

$$\int_{\gamma_i}^{\infty} \varrho_\phi(dx) = \frac{i}{N}. \quad (3.13)$$

THEOREM 3.5 (EIGENVALUE RIGIDITY [9]). *Fix $\tau > 0$, and suppose that (2.4), (2.3), and (2.5) hold. Then for $i \in \llbracket 1, M \rrbracket$ we have*

$$|\lambda_i - \gamma_i| \prec (\min\{i, K + 1 - i\})^{-1/3} K^{-2/3} \quad (3.14)$$

if either $i \leq (1 - \tau)K$ or $|\phi - 1| \geq \tau$.

3.2. Link to the semicircle law. It is often convenient to replace the Stieltjes transform $m_\phi(z)$ of $\varrho_\phi(dx)$ with the Stieltjes transform $w_\phi(z)$ of the measure

$$\phi^{1/2} x \varrho_{\phi^{-1}}(dx) = \frac{1}{2\pi} \sqrt{[(x - \gamma_-)(\gamma_+ - x)]_+} dx. \quad (3.15)$$

Note that this is nothing but Wigner's semicircle law centred at $\phi^{1/2} + \phi^{-1/2}$. Thus,

$$w_\phi(z) := \int \frac{\phi^{1/2} x \varrho_{\phi^{-1}}(dx)}{x - z} = \phi^{1/2} (1 + z m_{\phi^{-1}}(z)) = \frac{\phi^{1/2} + \phi^{-1/2} - z + i \sqrt{(z - \gamma_-)(\gamma_+ - z)}}{2}, \quad (3.16)$$

where in the last step we used (3.4). Note that

$$w_\phi = w_{\phi^{-1}}.$$

Using w_ϕ we can write (3.5) as

$$z = (1 - \phi^{-1/2} w_\phi^{-1})(\phi^{1/2} - w_\phi). \quad (3.17)$$

LEMMA 3.6. *For $z \in \mathbf{S}$ and $\phi \geq 1$ we have*

$$|m_\phi(z)| \asymp |w_\phi(z)| \asymp 1, \quad |1 - m_\phi(z)^2| \asymp |1 - w_\phi(z)^2| \asymp \sqrt{\kappa + \eta}, \quad (3.18)$$

as well as

$$\operatorname{Im} m_\phi(z) \asymp \operatorname{Im} w_\phi(z) \asymp \begin{cases} \sqrt{\kappa + \eta} & \text{if } E \in [\gamma_-, \gamma_+] \\ \frac{\eta}{\sqrt{\kappa + \eta}} & \text{if } E \notin [\gamma_-, \gamma_+]. \end{cases} \quad (3.19)$$

Similarly,

$$\operatorname{Re} m_\phi(z) - I(z) \asymp \operatorname{Re} w_\phi(z) - I(z) \asymp \begin{cases} \frac{\eta}{\sqrt{\kappa+\eta}} + \kappa & \text{if } E \in [\gamma_-, \gamma_+] \\ \sqrt{\kappa+\eta} & \text{if } E \notin [\gamma_-, \gamma_+], \end{cases} \quad (3.20)$$

where $I(z) := -1$ for $E \geq \phi^{1/2} + \phi^{-1/2}$ and $I(z) := +1$ for $E < \phi^{1/2} + \phi^{-1/2}$. Finally, for $z \in \mathbf{S}$ we have

$$\operatorname{Im} m_{\phi^{-1}}(z) \asymp \frac{1}{\phi} \operatorname{Im} m_\phi(z) \quad (3.21)$$

(All implicit constants depend on ω in the definition (3.8) of $\mathbf{S}$.)

PROOF. The estimates (3.18) and (3.19) follow from the explicit expressions in (3.4) and (3.16). In fact, these estimates have already appeared in previous works. Indeed, for m_ϕ the estimates (3.18) and (3.19) were proved in [9, Lemma 3.3]. In order to prove them for w_ϕ , we observe that the estimates (3.18) and (3.19) follow from the corresponding ones for the semicircle law, which were proved in [18, Lemma 4.3]. The estimates (3.20) follow from (3.19) and the elementary identity

$$\operatorname{Re} w_\phi = -\frac{E - \phi^{1/2} - \phi^{-1/2}}{1 + \eta / \operatorname{Im} w_\phi},$$

which can be derived from (3.17); the estimates for m_ϕ are derived similarly. Finally, (3.21) follows easily from

$$m_{\phi^{-1}}(z) = \frac{1}{\phi} \left(m_\phi(z) + \frac{1-\phi}{z} \right), \quad (3.22)$$

which may itself be derived from (3.5). $\square$

In analogy to w_ϕ (see (3.16)), we define the matrix-valued function

$$F(z) := \phi^{1/2}(1 + zG(z)).$$

Theorem 3.2 has the following analogue, which compares F with m_ϕ .

LEMMA 3.7. *Suppose that (2.4), (2.3), and (2.5) hold. Then*

$$|\langle \mathbf{v}, F(z)\mathbf{w} \rangle - w_\phi(z)\langle \mathbf{v}, \mathbf{w} \rangle| \prec \sqrt{\frac{\operatorname{Im} w_\phi(z)}{K\eta}} + \frac{1}{K\eta} \quad (3.23)$$

uniformly in $z \in \mathbf{S}$ and any deterministic unit vectors $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$. Moreover,

$$|\langle \mathbf{v}, F(z)\mathbf{w} \rangle - w_\phi(z)\langle \mathbf{v}, \mathbf{w} \rangle| \prec \sqrt{\frac{\operatorname{Im} w_\phi(z)}{K\eta}} \asymp (\kappa + \eta)^{-1/4} K^{-1/2} \quad (3.24)$$

uniformly in $z \in \tilde{\mathbf{S}}$ and any deterministic unit vectors $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$.

PROOF. The proof is an easy consequence of Theorem 3.2 and Lemma 3.6, combined with the fact that for $z \in \mathbf{S}$ or $z \in \tilde{\mathbf{S}}$ we have $|z| \asymp \phi^{1/2}$ for $\phi \geq 1$ and $|z| \asymp \phi^{-1/2}$ for $\phi \leq 1$. $\square$

3.3. Extension of the spectral domain. In this section we extend the spectral domain on which Theorem 3.2 and Lemma 3.7 hold. The argument relies on the Helffer-Sjöstrand functional calculus [14]. Define the domains

$$\hat{\mathbf{S}} \equiv \hat{\mathbf{S}}(\omega, K) := \{z \in \mathbb{C} : E \notin [\gamma_-, \gamma_+], \kappa \geq K^{-2/3+\omega}, \eta > 0\}, \quad \mathbf{B} \equiv \mathbf{B}(\omega) := \{z \in \mathbb{C} : |z| < \omega\}.$$

PROPOSITION 3.8. *Fix $\omega, \tau \in (0, 1)$.*

(i) If $\phi < 1 - \tau$ then

$$|\langle \mathbf{v}, G(z)\mathbf{w} \rangle - m_{\phi^{-1}}(z)\langle \mathbf{v}, \mathbf{w} \rangle| \prec \frac{1}{(\kappa + \eta)^2 + (\kappa + \eta)^{1/4}} K^{-1/2} \quad (3.25)$$

uniformly for $z \in \widehat{\mathbf{S}}$ and any deterministic unit vectors $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$.

(ii) If $|\phi - 1| \leq \tau$ then (3.25) holds uniformly for $z \in \widehat{\mathbf{S}} \setminus \mathbf{B}$ and any deterministic unit vectors $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$.

(iii) If $\phi > 1 + \tau$ then

$$|\langle \mathbf{v}, G(z)\mathbf{w} \rangle - m_{\phi^{-1}}(z)\langle \mathbf{v}, \mathbf{w} \rangle| \prec \frac{1}{\phi^{1/2}|z|((\kappa + \eta) + (\kappa + \eta)^{1/4})} K^{-1/2} \quad (3.26)$$

uniformly for $z \in \widehat{\mathbf{S}} \setminus \{0\}$ and any deterministic unit vectors $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$.

PROOF. By polarization and linearity, we may assume that $\mathbf{w} = \mathbf{v}$. Define the signed measure

$$\rho^\Delta(\mathrm{d}x) := \sum_{i=1}^M \langle \mathbf{v}, \zeta_i \rangle \langle \zeta_i, \mathbf{v} \rangle \delta_{\lambda_i}(\mathrm{d}x) - \varrho_{\phi^{-1}}(\mathrm{d}x), \quad (3.27)$$

so that

$$m^\Delta(z) := \int \frac{\rho^\Delta(\mathrm{d}x)}{x - z} = \langle \mathbf{v}, G(z)\mathbf{v} \rangle - m_{\phi^{-1}}(z).$$

The basic idea of the proof is to apply the Helffer-Sjöstrand formula to the function

$$f_z(x) := \frac{1}{x - z} - \frac{1}{x_0 - z},$$

where x_0 is chosen below. To that end, we need a smooth compactly supported cutoff function χ on the complex plane satisfying $\chi(w) \in [0, 1]$ and $|\partial_{\bar{w}}\chi(w)| \leq C(\omega, \tau)$. We distinguish the three cases $\phi < 1 - \tau$, $|\phi - 1| \leq \tau$, and $\phi > 1 + \tau$.

Let us first focus on the case $\phi < 1 - \tau$. Set $x_0 := \phi^{1/2} + \phi^{-1/2}$ and choose a constant $\omega' = \omega'(\omega, \tau) \in (0, \omega)$ small enough that $\gamma_- \geq 4\omega'$. We require that χ be equal to 1 in the ω' -neighbourhood of $[\gamma_-, \gamma_+]$ and 0 outside of the $2\omega'$ -neighbourhood of $[\gamma_-, \gamma_+]$. By Theorem 3.5 we have $\text{supp } \rho^\Delta \subset \{\chi = 1\}$ with high probability. Now choose z satisfying $\text{dist}(z, [\gamma_-, \gamma_+]) \geq 3\omega'$. Then the Helffer-Sjöstrand formula [14] yields, for $x \in \text{supp } \rho^\Delta$,

$$f_z(x) = \frac{1}{2\pi} \int_{\mathbb{C}} \frac{\partial_{\bar{w}}(f_z(w)\chi(w))}{x - w} \mathrm{d}w \quad (3.28)$$

with high probability, where $\mathrm{d}w$ denotes the two-dimensional Lebesgue measure in the complex plane. Noting that $\int \mathrm{d}\rho^\Delta = 0$, we may therefore write

$$m^\Delta(z) = \int \rho^\Delta(\mathrm{d}x) f_z(x) = \frac{1}{2\pi} \int_{\mathbb{C}} f_z(w) \partial_{\bar{w}}\chi(w) m^\Delta(w) \mathrm{d}w \quad (3.29)$$

with high probability, where in second step we used (3.28) and the fact that f_z is holomorphic away from z . The integral is supported on the set $\{\partial_{\bar{w}}\chi \neq 0\} \subset \{w : \text{dist}(w, [\gamma_-, \gamma_+]) \in [\omega', 2\omega']\}$, on which we have the estimates $|f_z(w)| \leq C(\kappa(z) + \eta(z))^{-2}$ and $|m^\Delta(w)| \prec K^{-1/2}$, as follows from Theorem 3.11 applied to $\mathbf{S}(\omega', K)$ and (3.21). Recalling Remark 3.3, we may plug these estimates into the integral to get

$$|m^\Delta(z)| \prec (\kappa + \eta)^{-2} K^{-1/2},$$

which holds for $\text{dist}(z, [\gamma_-, \gamma_+]) \geq 3\omega'$. (Recall that $|\partial_{\bar{w}}\chi(w)| \leq C$.) Combining this estimate with (3.11), the claim (3.25) follows for $z \in \widehat{\mathbf{S}}$.

Next, we deal with the case $|\phi - 1| \leq \tau$. The argument is similar. We again choose $x_0 := \phi^{1/2} + \phi^{-1/2}$. We require that χ be equal to 1 in the ω -neighbourhood of $[0, \gamma_+]$ and 0 outside of the 2ω -neighbourhood

of $[0, \gamma_+]$. We may now repeat the above argument almost verbatim. For $\text{dist}\{z, [0, \gamma_+]\} \geq 3\omega$ and $w \in \{\partial_{\bar{w}}\chi \neq 0\}$ we find that $|f_z(w)| \leq C(\kappa(z) + \eta(z))^{-2}$ and $|m^\Delta(w)| \prec K^{-1/2}$. Hence, recalling (3.11), we get (3.25) for $z \in \widehat{\mathbf{S}} \setminus \mathbf{B}$.

Finally, suppose that $\phi > 1 + \tau$. Now we set $x_0 := 0$. We choose the same ω' and cutoff function χ as in the case $\phi < 1 - \tau$ above. Suppose that $\text{dist}(z, [\gamma_-, \gamma_+]) \geq 3\omega'$ and $z \neq 0$. Thus, (3.28) holds with high probability for $x \in \text{supp } \rho^\Delta \setminus \{0\}$. Since $f_w(0) = 0$, we therefore find that (3.29) holds. As above, we find that for $w \in \{\partial_{\bar{w}}\chi \neq 0\}$ we have

$$|f_z(w)| \leq \frac{C\phi^{1/2}}{|z|(\kappa(z) + \eta(z))}$$

and $|m^\Delta(w)| \prec \phi^{-1}K^{-1/2}$. Recalling (3.11), we find that (3.26) follows easily. $\square$

Proposition 3.8 yields the follows result for F .

COROLLARY 3.9. *Fix $\omega, \tau \in (0, 1)$.*

(i) *If $\phi < 1 - \tau$ then*

$$|\langle \mathbf{v}, F(z)\mathbf{w} \rangle - w_\phi(z)\langle \mathbf{v}, \mathbf{w} \rangle| \prec \frac{\phi^{1/2}|z|}{(\kappa + \eta)^2 + (\kappa + \eta)^{1/4}} K^{-1/2} \quad (3.30)$$

uniformly for $z \in \widehat{\mathbf{S}}$ and any deterministic unit vectors $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$.

(ii) *If $|\phi - 1| \leq \tau$ then (3.30) holds uniformly for $z \in \widehat{\mathbf{S}} \setminus \mathbf{B}$ and any deterministic unit vectors $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$.*

(iii) *If $\phi > 1 + \tau$ then*

$$|\langle \mathbf{v}, F(z)\mathbf{w} \rangle - w_\phi(z)\langle \mathbf{v}, \mathbf{w} \rangle| \prec \frac{1}{(\kappa + \eta) + (\kappa + \eta)^{1/4}} K^{-1/2} \quad (3.31)$$

uniformly for $z \in \widehat{\mathbf{S}}$ and any deterministic unit vectors $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$.

3.4. Identities for the resolvent and eigenvalues. In this section we derive the identities on which our analysis of the eigenvalues and eigenvectors relies. We write the population covariance matrix Σ from (1.12) as

$$\Sigma = 1 + \phi^{1/2}VDV^*,$$

where $D = \text{diag}(d_i)_{i \in R}$ is an invertible diagonal $|R| \times |R|$ matrix and $V = [\mathbf{v}_i]_{i \in R}$ is the matrix of eigenvectors $\mathbf{v}_i$ of Σ indexed by the set R . Note that V is an $N \times |R|$ isometry, i.e. $V^*V = I_{|R|}$.

We use the definitions

$$G(z) := (H - z)^{-1}, \quad \widetilde{G}(z) := (Q - z)^{-1}, \quad F(z) := \phi^{1/2}(1 + zG(z)).$$

We introduce the $|R| \times |R|$ matrix

$$W(z) := V^*F(z)V.$$

We also denote by $\sigma(A)$ the spectrum of a square matrix A .

LEMMA 3.10. (i) *Suppose that $\mu \notin \sigma(H)$. Then $\mu \in \sigma(Q)$ if and only if*

$$\det(D^{-1} + W(\mu)) = 0. \quad (3.32)$$

(ii) *We have*

$$\Sigma^{1/2}\widetilde{G}(z)\Sigma^{1/2} = G(z) - G(z)V\frac{\phi^{1/2}z}{D^{-1} + W(z)}V^*G(z). \quad (3.33)$$

PROOF. To prove (i), we write the condition $\mu \in \sigma(Q)$ as

$$0 = \det(\Sigma^{1/2}H\Sigma^{1/2} - \mu) = \det(H - \mu\Sigma^{-1})\det(\Sigma) = \det(1 + G(\mu)(1 - \Sigma^{-1})\mu)\det(H - \mu)\det(\Sigma),$$

where we used that $\mu \notin \sigma(H)$. Using

$$1 - \Sigma^{-1} = V \frac{D}{\phi^{-1/2} + D} V^*,$$

the matrix identity $\det(1 + XY) = \det(1 + YX)$, and $\det(\Sigma) \neq 0$, we find

$$0 = \det\left(1 + \frac{D}{\phi^{-1/2} + D} \mu V^* G(\mu) V\right),$$

and the claim follows.

To prove (ii), we write

$$\Sigma^{1/2} \tilde{G}(z) \Sigma^{1/2} = (H - \Sigma^{-1}z)^{-1} = (H - z + (1 - \Sigma^{-1})z)^{-1}.$$

We now use the identity

$$(a + ubv)^{-1} = a^{-1} - a^{-1}u(b^{-1} + va^{-1}u)^{-1}va^{-1}, \quad (3.34)$$

which may be easily verified e.g. using a Neumann series. The claim follows by invoking (3.34) with $a = H - z$, $b = D(\phi^{-1/2} + D)^{-1}$, $u = V$, and $v = zV^*$. $\square$

The result (3.33), when restricted to the range of V , has an alternative form (3.35) which is often easier to work with, since it collects all of the randomness in the single quantity $W(z)$ on its right-hand side.

LEMMA 3.11. *We have*

$$V^* \tilde{G}(z) V = \frac{1}{\phi^{1/2}z} \left(D^{-1} - \frac{\sqrt{1 + \phi^{1/2}D}}{D} \frac{1}{D^{-1} + W(z)} \frac{\sqrt{1 + \phi^{1/2}D}}{D} \right). \quad (3.35)$$

PROOF. From (3.33) we get

$$(1 + \phi^{1/2}D)^{1/2} V^* \tilde{G}V (1 + \phi^{1/2}D)^{1/2} = V^*GV - V^*GV \frac{1}{(D^{-1} + \phi^{1/2})/(z\phi^{1/2}) + V^*GV} V^*GV.$$

Applying the identity

$$a - a(a + b)^{-1}a = b - b(a + b)^{-1}b$$

to the right-hand side yields

$$(1 + \phi^{1/2}D)^{1/2} V^* \tilde{G}V (1 + \phi^{1/2}D)^{1/2} = \frac{1}{\phi^{1/2}z} \left(D^{-1} + \phi^{1/2} - (D^{-1} + \phi^{1/2}) \frac{1}{D^{-1} + W} (D^{-1} + \phi^{1/2}) \right),$$

from which the claim follows. $\square$

4. Eigenvalue locations

In this section we prove Theorems 2.3 and 2.6. The arguments are similar to those of [25, Section 6], and we therefore only sketch the proofs. The proof of [25, Section 6] relies on three main steps: (i) establishing a forbidden region which contains with high probability no eigenvalues of Q ; (ii) a counting estimate for the special case where D does not depend on N , which ensures that each connected component of the allowed region (complement of the forbidden region) contains exactly the right number of eigenvalues of Q ; and (iii) a continuity argument where the counting result of (ii) is extended to arbitrary N -dependent

D using the the gaps established in (i) and the continuity of the eigenvalues as functions of the matrix entries. The steps (ii) and (iii) are exactly the same as in [25], and will not be repeated here. The step (i) differs slightly from that of [25], and in the proofs below we explain these differences.

We need the following eigenvalue interlacing result, which is purely deterministic. It holds for any nonnegative definite $M \times M$ matrix H and any rank-one deformation of the form $Q = (1 + \tilde{d}\mathbf{v}\mathbf{v}^*)^{1/2}H(1 + \tilde{d}\mathbf{v}\mathbf{v}^*)^{1/2}$ with $\tilde{d} \geq -1$ and $\mathbf{v} \in \mathbb{R}^M$.

LEMMA 4.1 (EIGENVALUE INTERLACING). *Let $|R| = 1$ and $D = d \in \mathcal{D}$. For $d > 0$ we have*

$$\mu_1 \geq \lambda_1 \geq \mu_2 \geq \cdots \geq \lambda_{M-1} \geq \mu_M \geq \lambda_M$$

and for $d < 0$ we have

$$\lambda_1 \geq \mu_1 \geq \lambda_2 \geq \cdots \geq \mu_{M-1} \geq \lambda_M \geq \mu_M.$$

PROOF. Using a simple perturbation argument (using that eigenvalues depend continuously on the matrix entries), we may assume without loss of generality that $\lambda_1, \dots, \lambda_M$ are all positive and distinct. Writing $\Sigma = 1 + \phi^{1/2}d\mathbf{v}\mathbf{v}^*$, we get from (3.33) that

$$\tilde{G}_{\mathbf{v}\mathbf{v}}(z) = a^2 G_{\mathbf{v}\mathbf{v}}(z) - a^2 G_{\mathbf{v}\mathbf{v}}(z)^2 \frac{1}{b(z)^{-1} + G_{\mathbf{v}\mathbf{v}}(z)}, \quad a := (\Sigma^{-1/2})_{\mathbf{v}\mathbf{v}}, \quad b(z) := \frac{z}{1 + \phi^{-1/2}d^{-1}}.$$

Note that $a > 0$. Thus we get

$$\frac{1}{G_{\mathbf{v}\mathbf{v}}(z)} + b(z) = \frac{a^2}{\tilde{G}_{\mathbf{v}\mathbf{v}}(z)}.$$

Writing this in spectral decomposition yields

$$\left(\sum_i \frac{\langle \mathbf{v}, \boldsymbol{\xi}_i \rangle^2}{\lambda_i - z} \right)^{-1} = a^2 \left(\sum_i \frac{\langle \mathbf{v}, \boldsymbol{\xi}_i \rangle^2}{\mu_i - z} \right)^{-1} - b(z). \quad (4.1)$$

As above, a simple perturbation argument implies that we may without loss of generality assume that all scalar products in (4.1) are nonzero. Now take $z \in (0, \infty)$. Note that $b(z)$ and d have the same sign.

To conclude the proof, we observe that the left-hand side of (4.1) defines a function of $z \in (0, \infty)$ with $M - 1$ singularities and M zeros, which is smooth and decreasing away from the singularities. Moreover, its zeros are the eigenvalues $\lambda_1, \dots, \lambda_M$. The interlacing property now follows from the fact that z is an eigenvalue of Q if and only if the left-hand side of (4.1) is equal to $-b(z)$. $\square$

COROLLARY 4.2. *For the rank- $|R|$ model (2.1) we have*

$$\mu_i \in [\lambda_{i+r}, \lambda_{i-r}] \quad (i \in \llbracket 1, M \rrbracket),$$

with the convention that $\lambda_i = 0$ for $i > K$ and $\lambda_i = \infty$ for $i < 1$.

We now move on to the proof of Theorem 2.3. Note that the function θ defined in (1.13) may be extended to a biholomorphic function from $\{\zeta \in \mathbb{C} : |\zeta| > 1\}$ to $\{z \in \mathbb{C} : z - (\phi^{1/2} + \phi^{-1/2}) \notin [-2, 2]\}$. Moreover, using (3.17) it is easy to check that for $|\zeta| > 1$ we have

$$w_\phi(z) = -\frac{1}{\zeta} \iff z = \theta(\zeta). \quad (4.2)$$

Throughout the following we shall make use of the subsets of outliers

$$\mathcal{O}_\tau^\pm := \{i : \pm d_i \geq 1 + K^{-1/3+\tau}\}$$

for $\tau \geq 0$. Note that $\mathcal{O} = \mathcal{O}_0^+ \cup \mathcal{O}_0^-$.

PROOF OF THEOREM 2.3. The proof of Proposition 2.3 is similar to that of [25, Equation (2.20)]. We focus first on the outliers to the right of the bulk spectrum. Let $\varepsilon > 0$. We shall prove that there exists an event Ξ of high probability (see Definition 2.1) such that for all $i \in \mathcal{O}_{4\varepsilon}^+$ we have

$$\mathbf{1}(\Xi)|\mu_i - \theta(d_i)| \leq C\Delta(d_i)K^{-1/2+\varepsilon} \quad (4.3)$$

and for $i \in \llbracket |\mathcal{O}_{4\varepsilon}^+| + 1, |\mathcal{O}_{4\varepsilon}^+| + r \rrbracket$ we have

$$\mathbf{1}(\Xi)|\mu_i - \gamma_+| \leq CK^{-2/3+8\varepsilon}. \quad (4.4)$$

Before proving (4.3) and (4.4), we show how they imply (2.9) for $d_i > 0$ and (2.10). From (4.4) we get for i satisfying $K^{-1/3} \leq d_i - 1 \leq K^{-1/3+4\varepsilon}$

$$\mathbf{1}(\Xi)|\mu_i - \theta(d_i)| \leq \mathbf{1}(\Xi)(|\mu_i - \gamma_+| + |\theta(d_i) - \gamma_+|) \leq CK^{-2/3+8\varepsilon} \leq C\Delta(d_i)K^{-1/2+8\varepsilon}. \quad (4.5)$$

Since $\varepsilon > 0$ was arbitrary, (2.9) for $d_i > 0$ and (2.10) follow from (4.3) and (4.5).

What remains is the proof of (4.3) and (4.4). As in [25, Proposition 6.5], the first step is to prove that with high probability there are no eigenvalues outside a neighbourhood of the classical outlier locations $\theta(d_i)$. To that end, we define for each $i \in \mathcal{O}_\varepsilon^+$ the interval

$$I_i(D) := \left[\theta(d_i) - \Delta(d_i)K^{-1/2+\varepsilon}, \theta(d_i) + \Delta(d_i)K^{-1/2+\varepsilon} \right].$$

Moreover, we set $I_0 := [0, \theta(1 + K^{-1/3+2\varepsilon})]$.

We now claim that with high probability the complement of the set $I(D) := I_0 \cup \bigcup_{i \in \mathcal{O}_\varepsilon^+} I_i(D)$ contains no eigenvalues of Q . Indeed, from Theorem 3.5 and Corollary 3.9 combined with Remark 3.3 (with small enough $\omega \equiv \omega(\varepsilon)$), we find that there exists an event Ξ of high probability such that $|\lambda_i - \gamma_+| \leq K^{-2/3+\varepsilon}$ for $i \in \llbracket 1, 2r \rrbracket$ and

$$\mathbf{1}(\Xi)\|W(x) - w_\phi(x)\| \leq \mathcal{E}(x)K^{-1/2+\varepsilon/2}$$

for all $x \notin I_0$, where we defined

$$\mathcal{E}(x) := \begin{cases} \kappa(x)^{-1/4} & \text{if } \kappa(x) \leq 1 \\ \frac{1}{\kappa(x)^2} \left(1 + \frac{\kappa(x)}{1+\phi^{-1/2}}\right) & \text{if } \kappa(x) > 1. \end{cases}$$

In particular, we have $\mathbf{1}(\Xi)\lambda_1 \leq \theta(1 + K^{-1/3+\varepsilon})$. Hence we find from (3.32) that on the event Ξ the value $x \notin I_0$ is an eigenvalue of Q if and only if the matrix

$$\mathbf{1}(\Xi)(D^{-1} + W(x)) = \mathbf{1}(\Xi)(D^{-1} + w_\phi(x) + O(\mathcal{E}(x)K^{-1/2+\varepsilon/2}))$$

is singular. Since $-d_i^{-1} = w_\phi(\theta(d_i))$ for $i \in \mathcal{O}_\varepsilon^+$, we conclude from the definition of $I(D)$ that it suffices to show that if $x \notin I(D)$ then

$$\min_{i \in \mathcal{O}_\varepsilon^+} |w_\phi(x) - w_\phi(\theta(d_i))| \gg \mathcal{E}(x)K^{-1/2+\varepsilon/2}. \quad (4.6)$$

We prove (4.6) using the two following observations. First, w_ϕ is monotone increasing on (γ_+, ∞) and

$$w'_\phi(x) \asymp (d_i^2 - 1)^{-1} \quad (x \in I_i(D)),$$

as follows from (4.2). Second,

$$\Delta(d_i) \asymp \frac{\mathcal{E}(\theta(d_i))}{|w'_\phi(\theta(d_i))|} = (d_i^2 - 1)\mathcal{E}(\theta(d_i)).$$

We omit further details, which may be found e.g. in [25, Section 6]. Thus we conclude that on the event Ξ the complement of $I(D)$ contains no eigenvalues of Q .

The next step of the proof consists in making sure that the allowed neighbourhoods $I_i(D)$ contain exactly the right number of outliers; the counting argument (sketched in the steps (ii) and (iii) at the beginning of this section) follows that of [25, Section 6]. First we consider the case $D = D(0)$ where for all $i \neq j \in \mathcal{O}_\varepsilon^+$ we have $d_i(0), d_j(0) \geq 2$ and $|d_i(0) - d_j(0)| \geq 1$, and show that each interval $\{I_i(D(0)) : i \in \mathcal{O}_\varepsilon^+\}$ contains exactly one eigenvalue of Q (see [25, Proposition 6.6]). We then deduce the general case by a continuity argument, by choosing an appropriate continuous path $(D(t))_{t \in [0,1]}$ joining the initial configuration $D(0)$ to the desired final configuration $D = D(1)$. The continuity argument requires the existence of a gap in the set $I(D)$ to the left of $\bigcup_{i \in \mathcal{O}_{4\varepsilon}^+} I_i(D)$. The existence of such a gap follows easily from the definition of $I(D)$ and the fact that $|R|$ is bounded. The details are the same as in [25, Section 6.5]. Hence (4.3) follows. Moreover, (4.4) follows from the same argument combined with Corollary 4.2 for a lower bound on μ_i . This concludes the analysis of the outliers to the right of the bulk spectrum.

The case of outliers to the left of the bulk spectrum is analogous. Here we assume that $\phi < 1 - \tau$. The argument is exactly the same as for $d_i > 0$, except that we use the bound (3.30) to the left of the bulk spectrum as well as $|\lambda_i - \gamma_-| \leq K^{-2/3+\varepsilon}$ for $i \in \llbracket K - 2r, K \rrbracket$ with high probability. $\square$

PROOF OF THEOREM 2.6. We only give the proof of (2.14); the proof of (2.15) is analogous. Fix $\varepsilon > 0$. By Theorem 2.3, Theorem 3.5, Theorem 3.2, Lemma 3.7, and Remark 3.3, there exists a high-probability event $\Xi \equiv \Xi_N(\varepsilon)$ satisfying the following conditions.

(i) We have

$$\mathbf{1}(\Xi) |\mu_{s_++1} - \gamma_+| \leq K^{-2/3+\varepsilon}, \quad \mathbf{1}(\Xi) |\lambda_i - \gamma_i| \leq i^{-1/3} K^{-2/3+\varepsilon} \quad (i \leq (1 - \tau)K). \quad (4.7)$$

(ii) For $z \in \mathbf{S}(\varepsilon, K)$ we have

$$\mathbf{1}(\Xi) \|W(z) - w_\phi(z)\| \leq K^\varepsilon \left(\sqrt{\frac{\operatorname{Im} w_\phi(z)}{K\eta}} + \frac{1}{K\eta} \right) \quad (4.8)$$

and

$$\max_{i,j} |\langle \mathbf{v}_i, G(z) \mathbf{v}_j \rangle - m_{\phi^{-1}}(z) \delta_{ij}| \leq K^\varepsilon \left(\sqrt{\frac{\operatorname{Im} m_{\phi^{-1}}(z)}{M\eta}} + \frac{1}{M\eta} \right). \quad (4.9)$$

For the following we fix a realization $H \in \Xi$. We suppose first that

$$\alpha_+ \geq K^{-1/3+\varepsilon}, \quad (4.10)$$

and define $\eta := K^{-1+2\varepsilon} \alpha_+^{-1}$. Now suppose that x satisfies

$$x \in [\gamma_+ - 1, \gamma_+ + K^{-2/3+2\varepsilon}], \quad \operatorname{dist}(x, \sigma(H)) > \eta. \quad (4.11)$$

We shall show, using (3.32), that any x satisfying (4.11) cannot be an eigenvalue of Q . First we deduce from (4.8) that

$$\|W(x) - W(x + i\eta)\| \leq C(1 + \phi) \max_i \operatorname{Im} G_{\mathbf{v}_i \mathbf{v}_i}(x + i\eta). \quad (4.12)$$

The estimate (4.12) follows by spectral decomposition of $F(\cdot)$ together with the estimate $2|\lambda_i - x| \geq \sqrt{(\lambda_i - x)^2 + \eta^2}$ for all i . We get from (4.12) and Lemma 3.6 that

$$W(x) = w_\phi(x + i\eta) + O\left(\operatorname{Im} w_\phi(x + i\eta) + \frac{K^\varepsilon}{K\eta}\right) = -1 + O\left(\sqrt{\kappa(x)} + \sqrt{\eta} + K^{-\varepsilon} \alpha_+^{-1}\right).$$

Recalling (3.32), we conclude that on the event Ξ the value x is not an eigenvalue of Q provided

$$\min_i |1/d_i - 1| \geq K^{\varepsilon/2} \left(\sqrt{\kappa(x)} + \sqrt{\eta} + K^{-\varepsilon} \alpha_+^{-1} \right).$$

It is easy to check that this condition is satisfied if

$$\kappa(x) + \eta \leq CK^{-\varepsilon}\alpha_+^2,$$

which holds provided that

$$\kappa(x) \leq CK^{-\varepsilon}\alpha_+^2,$$

where we used (4.10). Recalling (4.7), we therefore conclude that for $i \leq K^{1-2\varepsilon}\alpha_+^3$ the set

$$\left\{x \in [\lambda_{i-r-1}, \gamma_+ + K^{-2/3+2\varepsilon}] : \text{dist}(x, \sigma(H)) > K^{-1+2\varepsilon}\alpha_+^{-1}\right\}$$

contains no eigenvalue of Q .

The next step of the proof is a counting argument (sketched in the steps (ii) and (iii) at the beginning of this section), which uses the eigenvalue interlacing from Lemma 4.1. The details are the same as in [25, Section 6], and hence omitted here. The counting argument implies that for $i \leq K^{1-2\varepsilon}\alpha_+^3$ and assuming (4.10) we have

$$|\mu_{i+s_+} - \lambda_i| \leq CK^{-1+2\varepsilon}\alpha_+^{-1}. \quad (4.13)$$

What remains is to check (4.13) for the cases $\alpha_+ < K^{-1/3+\varepsilon}$ and $i > K^{1-2\varepsilon}\alpha_+^3$.

Suppose first that $\alpha_+ < K^{-1/3+\varepsilon}$. Then using the rigidity from (4.7) and interlacing from Corollary 4.2 we find

$$|\mu_{i+s_+} - \lambda_i| \leq C i^{-1/3} K^{-2/3+\varepsilon} \leq CK^{-1+2\varepsilon}\alpha_+^{-1},$$

where we used the trivial bound $i \geq 1$. Similarly, if $i > K^{1-2\varepsilon}\alpha_+^3$ satisfies $i \leq (1-\tau)K$, we may repeat the same estimate.

We conclude that (4.13) under the sole assumption that $i \leq (1-\tau)K$. Since $\varepsilon > 0$ was arbitrary, (2.14) follows. $\square$

5. Outlier eigenvectors

In this section we focus on the outlier eigenvectors ξ_a , $a \in \mathcal{O}$. Here we in fact prove Theorem 2.9 under the stronger assumption

$$1 + K^{-1/3+\tau} \leq d_i \leq \tau^{-1} \quad (i \in A) \quad (5.1)$$

instead of $1 + K^{-1/3} \leq d_i \leq \tau^{-1}$. How to improve the lower bound from $1 + K^{-1/3+\tau}$ to the claimed $K^{-1/3}$ requires a completely different approach, relying on eigenvector delocalization bounds, and is presented in Section 6 in conjunction with results for the non-outlier eigenvectors ξ_a , $a \notin \mathcal{O}$.

The proof of Theorem 2.14 is similar to that of Theorem 2.9; one has to adapt the proof to cover the range $d_i \in [1 + \tau, \infty)$ instead of $d_i \in [1 + K^{-1/3}, \tau^{-1}]$. The key input is the extension of the spectral domain from Corollary 3.9. For the sake of brevity we omit the details of the proof of Theorem 2.14, and focus solely on Theorem 2.9.

The following proposition is the main result of this section.

PROPOSITION 5.1. *Fix $\tau > 0$. Suppose that A satisfies (5.1). Then for all $i, j = 1, \dots, M$ we have*

$$\begin{aligned} \langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle &= \delta_{ij} \mathbf{1}(i \in A) u(d_i) + O_{\prec} \left[\frac{\mathbf{1}(i, j \in A)}{(d_i - 1)^{1/4} (d_j - 1)^{1/4} M^{1/2}} \right. \\ &\quad \left. + \frac{\sqrt{\sigma_i \sigma_j}}{M} \left(\frac{1}{\nu_i} + \frac{\mathbf{1}(i \in A)}{d_i - 1} \right) \left(\frac{1}{\nu_j} + \frac{\mathbf{1}(j \in A)}{d_j - 1} \right) + \frac{\mathbf{1}(i \in A) \mathbf{1}(j \notin A) (d_i - 1)^{1/2} \sqrt{\sigma_j}}{(1 + \phi)^{1/4} \nu_j M^{1/2}} + (i \leftrightarrow j) \right], \end{aligned} \quad (5.2)$$

where the symbol $(i \leftrightarrow j)$ denotes the preceding terms with i and j interchanged.

Note that, under the assumption (5.1), Theorem 2.9 is an easy consequence of Proposition 5.1. As explained above, the proof of Theorem 2.9 in full generality is given in Section 6, where we give the additional argument required to relax (5.1).

The rest of this section is devoted to the proof of Proposition 5.1. We assume throughout that (5.1) holds.

5.1. Non-overlapping outliers. We first prove a slightly stronger version of (5.2) under the additional *non-overlapping condition*

$$\nu_i(A) \geq (d_i - 1)^{-1/2} K^{-1/2+\delta} \quad (5.3)$$

for all $i \in A$, where $\delta > 0$ is a positive constant. This is a precise version of the second condition of (2.19), whose interpretation was given below (2.19): an outlier indexed by A cannot overlap with an outlier indexed by A^c . Note, however, that there is no restriction on the outliers indexed by A overlapping among themselves. The assumption (5.3) will be removed in Section 5.2. The main estimate for non-overlapping outliers is the following.

PROPOSITION 5.2. *Fix $\tau > 0$ and $\delta > 0$. Suppose that A satisfies (5.1) and (5.3) for all $i \in A$. Then for all $i, j = 1, \dots, M$ we have*

$$\begin{aligned} \langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle &= \delta_{ij} \mathbf{1}(i \in A) u(d_i) + O_{\prec} \left[\frac{\mathbf{1}(i, j \in A)}{(d_i - 1)^{1/4} (d_j - 1)^{1/4} M^{1/2}} \right. \\ &\quad \left. + \frac{\sqrt{\sigma_i \sigma_j}}{M} \left(\frac{1}{\nu_i} + \frac{\mathbf{1}(i \in A)}{d_i - 1} \right) \left(\frac{1}{\nu_j} + \frac{\mathbf{1}(j \in A)}{d_j - 1} \right) + \frac{\mathbf{1}(i \in A) \mathbf{1}(j \notin A) (d_i - 1)^{1/2} \sqrt{\sigma_j}}{(1 + \phi)^{1/4} |d_i - d_j| M^{1/2}} + (i \leftrightarrow j) \right]. \end{aligned} \quad (5.4)$$

REMARK 5.3. The only difference between (5.2) and (5.4) is the term proportional to $\mathbf{1}(j \notin A)$ on the last line. In order to prove (5.2) without the overlapping condition (5.3), it is necessary to start from the stronger bound (5.4); see Section 5.2 below.

The rest of this subsection is devoted to the proof of Proposition 5.2. We begin by defining $\omega := \tau/2$ and letting $\varepsilon < \min\{\tau/3, \delta\}$ be a positive constant to be determined later. We choose a high-probability event $\Xi \equiv \Xi_N(\varepsilon, \tau)$ (see Definition 3.1) satisfying the following conditions.

(i) We have

$$\mathbf{1}(\Xi) |W_{ij}(z) - w_\phi(z) \delta_{ij}| \leq |z - \gamma_+|^{-1/4} K^{-1/2+\varepsilon} \quad (5.5)$$

for $i, j \in R$, large enough K , and all z in the set

$$\{z \in \mathbb{C} : \operatorname{Re} z \geq \gamma_+ + K^{-2/3+\omega}, |z| \leq \omega^{-1}\}. \quad (5.6)$$

(ii) For all i satisfying $1 + K^{-1/3} \leq d_i \leq \omega^{-1}$ we have

$$\mathbf{1}(\Xi) |\mu_i - \theta(d_i)| \leq (d_i - 1)^{1/2} K^{-1/2+\varepsilon} \quad (5.7)$$

(iii) We have

$$\mathbf{1}(\Xi) |\mu_{s_++1} - \gamma_+| \leq K^{-2/3+\varepsilon}. \quad (5.8)$$

Note that such an event Ξ exists. Indeed, (5.7) and (5.8) may be satisfied using Theorem 2.3, and (5.5) using Theorem 3.7 combined with Remark 3.3.

For the sequel we fix a realization $H \in \Xi$ satisfying the conditions (i)–(iii) above. Hence, the rest of the proof of Proposition 5.2 is entirely deterministic, and the randomness only enters in ensuring that Ξ has high probability. Our starting point is a contour integral representation of the projection P_A . In order to construct the contour, we define for each $i \in A$ the radius

$$\rho_i := \frac{\nu_i \wedge (d_i - 1)}{2}. \quad (5.9)$$

We define the contour $\Gamma := \partial \Upsilon$ as the boundary of the union of discs $\Upsilon := \bigcup_{i \in A} B_{\rho_i}(d_i)$, where $B_\rho(d)$ is the open disc of radius ρ around d . We shall sometimes need the decomposition $\Gamma = \bigcup_{i \in A} \Gamma_i$, where $\Gamma_i := \Gamma \cap \partial B_{\rho_i}(d_i)$. See Figure 5.1 for an illustration of Γ .

We shall have to use the estimate (5.5) on the set $\overline{\theta(\Upsilon)}$. Its applicability is an immediate consequence of the following lemma.

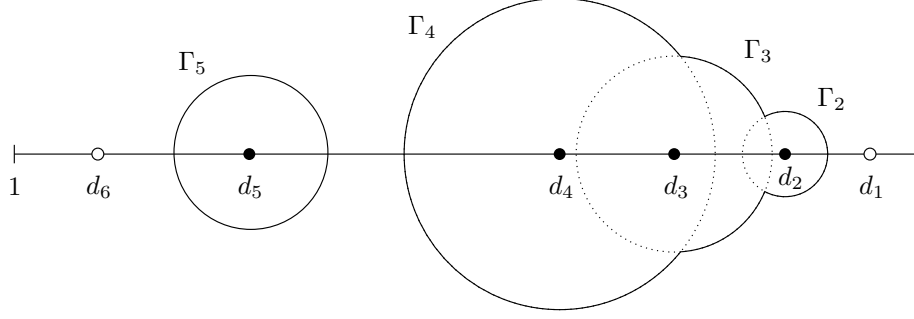


FIGURE 5.1. The integration contour $\Gamma = \bigcup_{i \in A} \Gamma_i$. In this example Γ consists of two components, and we have $|R| = 6$ with $A = \{2, 3, 4, 5\}$. We draw the locations of d_i with $i \in A$ using black dots and the other d_i using white dots. The contour is constructed by drawing circles of radius ρ_i around each d_i for $i \in A$ (depicted with dotted lines). The piece Γ_i consists of the points on the circle centred at d_i that lie outside all other circles.

LEMMA 5.4. *The set $\overline{\theta(\Upsilon)}$ lies in (5.6).*

PROOF. It is easy to check that $\theta(\zeta) \leq \omega^{-1}$ for all $\zeta \in \Upsilon$. In order to check the lower bound on $\operatorname{Re} \theta(\zeta)$, we note that for any $\alpha \in (0, 1)$ there exists a constant $c \equiv c(\alpha, \tau)$ such that

$$\operatorname{Re} \theta(\zeta) \geq \gamma_+ + c(\operatorname{Re} \zeta - 1)^2$$

for $\operatorname{Re} \zeta \geq 1$, $|\operatorname{Im} \zeta| \leq \alpha(\operatorname{Re} \zeta - 1)$, and $|\zeta| \leq \tau^{-1}$. Now the claim follows easily from $\operatorname{Re} \zeta \geq 1 + K^{-1/3+\tau}/2$ for all $\zeta \in \Upsilon$, by choosing $\alpha = 1/\sqrt{3}$. $\square$

LEMMA 5.5. *Each outlier $\{\mu_i\}_{i \in A}$ lies in $\theta(\Upsilon)$, and all other eigenvalues of Q lie in the complement of $\theta(\Upsilon)$.*

PROOF. It suffices to prove that (a) for each $i \in A$ we have $\mu_i \in \theta(B_{\rho_i}(d_i))$ and (b) all the other eigenvalues μ_j satisfy $\mu_j \notin \theta(B_{\rho_i}(d_i))$ for all $i \in A$.

In order to prove (a), we note that

$$\rho_i \geq \frac{1}{2} (d_i - 1)^{-1/2} K^{-1/2+\delta}, \quad (5.10)$$

for $i \in A$, as follows from (5.3) and (5.1). Using

$$|\theta'(\zeta)| \asymp |\zeta - 1| \quad (\operatorname{Re} \zeta \geq 1, |\zeta| \leq \tau^{-1}), \quad (5.11)$$

it is then not hard to get (a) from (5.10) and (5.7).

In order to prove (b), we consider the two cases (i) $1 + K^{-1/3} \leq d_j \leq \omega^{-1}$ with $j \notin A$, and (ii) and $j \geq s_+ + 1$. In the case (i), the claim (b) follows using (5.7), (5.11), and (5.3). In the case (ii), the claim (b) follows from (5.8) and the estimate

$$|\theta(\zeta) - \gamma_+| \asymp |\zeta - 1|^2 \quad (\operatorname{Re} \zeta \geq 1, |\zeta| \leq \tau^{-1}). \quad (5.12)$$

This concludes the proof. $\square$

Using the spectral decomposition of $\tilde{G}(z)$, Lemma 5.5, and the residue theorem, we may write the projection P_A as

$$P_A = -\frac{1}{2\pi i} \oint_{\theta(\Gamma)} \tilde{G}(z) dz = -\frac{1}{2\pi i} \oint_{\Gamma} \tilde{G}(\theta(\zeta)) \theta'(\zeta) d\zeta.$$

Hence we get from (3.35) that

$$V^* P_A V = \phi^{-1/2} \frac{1}{2\pi i} \oint_{\Gamma} \frac{\sqrt{1+\phi^{1/2}D}}{D} \frac{1}{D^{-1}+W(\theta(\zeta))} \frac{\sqrt{1+\phi^{1/2}D}}{D} \frac{\theta'(\zeta)}{\theta(\zeta)} d\zeta. \quad (5.13)$$

This is the desired integral representation of P_A .

We first use (5.13) to compute $\langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle$ in the case $i, j \in R$, where $\mathbf{v}_i$ and $\mathbf{v}_j$ lie in the range of V . In that case we get from (5.13) that

$$\langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle = \frac{\sqrt{\sigma_i \sigma_j}}{\phi^{1/2} d_i d_j} \frac{1}{2\pi i} \oint_{\Gamma} \left(\frac{1}{D^{-1}+W(\theta(\zeta))} \right)_{ij} \frac{\theta'(\zeta)}{\theta(\zeta)} d\zeta.$$

We now perform a resolvent expansion on the denominator

$$D^{-1} + W(\theta) = (D^{-1} + w_{\phi}(\theta)) - \Delta(\theta), \quad \Delta(\theta) := w_{\phi}(\theta) - W(\theta). \quad (5.14)$$

Thus we get

$$\langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle = \frac{\sqrt{\sigma_i \sigma_j}}{\phi^{1/2} d_i d_j} (S_{ij}^{(0)} + S_{ij}^{(1)} + S_{ij}^{(2)}), \quad (5.15)$$

where we defined

$$S_{ij}^{(0)} := \frac{1}{2\pi i} \oint_{\Gamma} \left(\frac{1}{D^{-1} + w_{\phi}(\theta(\zeta))} \right)_{ij} \frac{\theta'(\zeta)}{\theta(\zeta)} d\zeta \quad (5.16)$$

$$S_{ij}^{(1)} := \frac{1}{2\pi i} \oint_{\Gamma} \left(\frac{1}{D^{-1} + w_{\phi}(\theta(\zeta))} \Delta(\theta(\zeta)) \frac{1}{D^{-1} + w_{\phi}(\theta(\zeta))} \right)_{ij} \frac{\theta'(\zeta)}{\theta(\zeta)} d\zeta \quad (5.17)$$

$$S_{ij}^{(2)} := \frac{1}{2\pi i} \oint_{\Gamma} \left(\frac{1}{D^{-1} + w_{\phi}(\theta(\zeta))} \Delta(\theta(\zeta)) \frac{1}{D^{-1} + W(\theta(\zeta))} \Delta(\theta(\zeta)) \frac{1}{D^{-1} + w_{\phi}(\theta(\zeta))} \right)_{ij} \frac{\theta'(\zeta)}{\theta(\zeta)} d\zeta. \quad (5.18)$$

We begin by computing

$$S_{ij}^{(0)} = \delta_{ij} \frac{1}{2\pi i} \oint_{\Gamma} \left(\frac{1}{d_i^{-1} - \zeta^{-1}} \right)_{ij} \frac{\theta'(\zeta)}{\theta(\zeta)} d\zeta = \delta_{ij} \mathbf{1}(i \in A) \frac{d_i^2 - 1}{\theta(d_i)}, \quad (5.19)$$

where we used Cauchy's theorem, (4.2), and the fact that d_i lies in Υ if and only if $i \in A$.

Next, we estimate

$$S_{ij}^{(1)} = d_i d_j \frac{1}{2\pi i} \oint_{\Gamma} \frac{f_{ij}(\zeta)}{(\zeta - d_i)(\zeta - d_j)} d\zeta, \quad f_{ij}(\zeta) := \zeta^2 \Delta(\theta(\zeta)) \frac{\theta'(\zeta)}{\theta(\zeta)}, \quad (5.20)$$

using the fact that f_{ij} is holomorphic inside Γ and satisfies the bounds

$$|f_{ij}(\zeta)| \leq C \phi^{1/2} (1 + \phi)^{-1} |\zeta - 1|^{1/2} K^{-1/2+\varepsilon}, \quad |f'_{ij}(\zeta)| \leq C \phi^{1/2} (1 + \phi)^{-1} |\zeta - 1|^{-1/2} K^{-1/2+\varepsilon}. \quad (5.21)$$

The first bound of (5.21) follows from (5.5), (5.11), and (5.12). The second bound of (5.21) follows by plugging the first one into

$$f'_{ij}(\zeta) = \frac{1}{2\pi i} \oint_{\mathcal{C}} \frac{f_{ij}(\xi)}{(\xi - \zeta)^2} d\xi,$$

where the contour $\mathcal{C}$ is the circle of radius $|\zeta - 1|/2$ centred at ζ . (By assumptions on ε and ω , the function f_{ij} is holomorphic in a neighbourhood of the closed interior of $\mathcal{C}$.)

In order to estimate (5.20), we consider the three cases (i) $i, j \in A$, (ii) $i \in A, j \notin A$, (iii) $i \notin A, j \in A$. Note that (5.20) vanishes if $i, j \notin A$. We start with the case (i). Suppose first that $i \neq j$ and $d_i \neq d_j$. Then we find

$$|S_{ij}^{(1)}| = |d_i d_j| \left| \frac{f_{ij}(d_i) - f_{ij}(d_j)}{d_i - d_j} \right| \leq \frac{|d_i d_j|}{|d_i - d_j|} \left| \int_{d_i}^{d_j} |f'_{ij}(t)| dt \right| \leq \frac{C |d_i d_j| \phi^{1/2}}{(1 + \phi)(d_i - 1)^{1/4} (d_j - 1)^{1/4}} K^{-1/2+\varepsilon}.$$

A simple limiting argument shows that this bound is also valid for $d_i = d_j$ and $i = j$. Next, in the case (ii) we get from (5.21)

$$|S_{ij}^{(1)}| = \frac{|d_i d_j f_{ij}(d_i)|}{|d_i - d_j|} \leq \frac{C|d_i d_j| \phi^{1/2} (d_i - 1)^{1/2}}{(1 + \phi)|d_i - d_j|} K^{-1/2+\varepsilon}.$$

A similar estimate holds for the case (iii). Putting all three cases together, we find

$$|S_{ij}^{(1)}| \leq \frac{C \mathbf{1}(i, j \in A) |d_i d_j| \phi^{1/2}}{(1 + \phi)(d_i - 1)^{1/4} (d_j - 1)^{1/4}} K^{-1/2+\varepsilon} + \frac{C \mathbf{1}(i \in A) \mathbf{1}(j \notin A) |d_i d_j| \phi^{1/2} (d_i - 1)^{1/2}}{(1 + \phi)|d_i - d_j|} K^{-1/2+\varepsilon} + (i \leftrightarrow j). \quad (5.22)$$

What remains is the estimate of $S_{ij}^{(2)}$. Here residue calculations are unavailable, and the precise choice of the contour Γ is crucial. We use the following basic estimate to control the integral.

LEMMA 5.6. *For $k \in A$, $l \in R$, and $\zeta \in \Gamma_k$ we have*

$$|\zeta - d_l| \asymp \rho_k + |d_k - d_l|.$$

PROOF. The upper bound $|\zeta - d_l| \leq \rho_k + |d_k - d_l|$ is trivial, so that we only focus on the lower bound. Suppose first that $l \notin A$. Then we get $|\zeta - d_l| \geq |d_k - d_l| - \rho_k$, from which the claim follows since $|d_k - d_l| \geq 2\rho_k$ by (5.9).

For the remainder of the proof we may therefore suppose that $l \in A$. Define $\delta := |d_k - d_l| - \rho_k - \rho_l$, the distance between the discs $D_{\rho_k}(d_k)$ and $D_{\rho_l}(d_l)$ (see Figure 5.1). We consider the two cases $4\delta \leq |d_k - d_l|$ and $4\delta > |d_k - d_l|$ separately.

Suppose first that $4\delta \leq |d_k - d_l|$. Then by definition of δ we have $|d_k - d_l| \leq \frac{4}{3}(\rho_k + \rho_l)$. Now a simple estimate using the definition of ρ_i yields $\rho_k/5 \leq \rho_l \leq 5\rho_k$, from which we conclude $|d_k - d_l| \leq 8\rho_k$. The claim now follows from the bound $|\zeta - d_l| \geq \rho_l$.

Suppose now that $4\delta > |d_k - d_l|$. Hence $\rho_k + \rho_l \leq \frac{3}{4}|d_k - d_l|$, so that in particular $\rho_k \leq |d_k - d_l|$. Thus we get

$$|\zeta - d_l| \geq |d_k - d_l| - \rho_k - \rho_l \geq \frac{1}{4}|d_k - d_l| \geq \frac{1}{8}(|d_k - d_l| + \rho_k).$$

This concludes the proof. $\square$

From (5.18), (5.5), (5.11), and (5.12) we get

$$|S_{ij}^{(2)}| \leq C \oint_{\Gamma} \frac{\phi^{1/2} |d_i d_j| K^{-1+2\varepsilon}}{(1 + \phi) |\zeta - d_i| |\zeta - d_j|} \left\| \frac{1}{D^{-1} + W(\theta(\zeta))} \right\| |d\zeta|, \quad (5.23)$$

where we also used the estimate $|\theta(\zeta)| \asymp \phi^{-1/2}(1 + \phi)$ for $\zeta \in \Gamma$.

In order to estimate the matrix norm, we observe that for $\zeta \in \Gamma_k$ we have on the one hand

$$\|W(\theta) - w_{\phi}(\theta)\| \leq (d_k - 1)^{-1/2} K^{-1/2+\varepsilon}$$

from (5.5) and on the other hand

$$|w_{\phi}(\theta) - d_l^{-1}| \geq c(|\zeta - d_l| \wedge 1) \geq c|\zeta - d_k| = c\rho_k \geq c(d_k - 1)^{-1/2} K^{-1/2+\delta}$$

for any $l \in R$, where in the last step we used (5.10). Since $\varepsilon < \delta$, these estimates combined with a resolvent expansion give the bound

$$\left\| \frac{1}{D^{-1} + W(\theta(\zeta))} \right\| \leq \frac{1}{\min_{b \in R} |w_{\phi}(\theta) - d_b^{-1}| - \|W(\theta) - w_{\phi}(\theta)\|} \leq \frac{C}{\rho_k}$$

for $\zeta \in \Gamma_k$. Decomposing the integration contour in (5.23) as $\Gamma = \bigcup_{k \in A} \Gamma_k$, and recalling that Γ_k has length bounded by $2\pi\rho_k$, we get from Lemma 5.6

$$|S_{ij}^{(2)}| \leq C \sum_{k \in A} \sup_{\zeta \in \Gamma_k} \frac{\phi^{1/2} |d_i d_j| K^{-1+2\varepsilon}}{(1 + \phi) |\zeta - d_i| |\zeta - d_j|} \leq C \sum_{k \in A} \frac{\phi^{1/2} |d_i d_j| K^{-1+2\varepsilon}}{(1 + \phi)(\rho_k + |d_k - d_i|)(\rho_k + |d_k - d_j|)}. \quad (5.24)$$

We estimate the right-hand side using Cauchy-Schwarz. For $i \notin A$ we find, using (5.9),

$$\sum_{k \in A} \frac{1}{(\rho_k + |d_k - d_i|)^2} \leq \sum_{k \in A} \frac{1}{|d_k - d_i|^2} \leq \frac{C}{\nu_i^2}. \quad (5.25)$$

For $i \in A$ we use (5.9) and the estimate $\rho_k + |d_i - d_k| \geq \rho_i$ for all $k \in A$ to get

$$\sum_{k \in A} \frac{1}{(\rho_k + |d_k - d_i|)^2} \leq \frac{C}{\rho_i^2} \leq \frac{C}{\nu_i^2} + \frac{C}{(d_i - 1)^2}. \quad (5.26)$$

From (5.24), (5.25), and (5.26), we get

$$|S_{ij}^{(2)}| \leq \frac{C\phi^{1/2}|d_i d_j|K^{-1+2\varepsilon}}{1+\phi} \left(\frac{1}{\nu_i} + \frac{\mathbf{1}(i \in A)}{d_i - 1} \right) \left(\frac{1}{\nu_j} + \frac{\mathbf{1}(j \in A)}{d_j - 1} \right). \quad (5.27)$$

Recall that $M \asymp (1+\phi)K$. Hence, plugging (5.19), (5.22), and (5.27) into (5.15), we find

$$\begin{aligned} \langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle &= \delta_{ij} \mathbf{1}(i \in A) u(d_i) + O \left[\frac{\mathbf{1}(i, j \in A) K^\varepsilon}{(d_i - 1)^{1/4} (d_j - 1)^{1/4} M^{1/2}} \right. \\ &\quad \left. + \frac{\sqrt{\sigma_i \sigma_j} K^{2\varepsilon}}{M} \left(\frac{1}{\nu_i} + \frac{\mathbf{1}(i \in A)}{d_i - 1} \right) \left(\frac{1}{\nu_j} + \frac{\mathbf{1}(j \in A)}{d_j - 1} \right) + \frac{\mathbf{1}(i \in A) \mathbf{1}(j \notin A) (d_i - 1)^{1/2} \sqrt{\sigma_j} K^\varepsilon}{(1+\phi)^{1/4} |d_i - d_j| M^{1/2}} + (i \leftrightarrow j) \right]. \end{aligned} \quad (5.28)$$

We have proved (5.28) under the assumption that $i, j \in R$. The general case is an easy corollary. For general $i, j \in \llbracket 1, M \rrbracket$, we define $\hat{R} := R \cup \{i, j\}$ and consider

$$\hat{\Sigma} := 1 + \phi^{1/2} \hat{V} \hat{D} \hat{V}^*, \quad \hat{V} := [\mathbf{v}_k]_{k \in \hat{R}}, \quad \hat{D} := \text{diag}(\hat{d}_k)_{k \in \hat{R}},$$

where $\hat{d}_k := d_k$ for $k \in R$ and $\hat{d}_k \in (0, 1/2)$ for $k \in \hat{R} \setminus R$. Since $|\hat{R}| \leq r + 2$ and $\hat{D}$ is invertible, we may apply the result (5.28) to this modified model. Now taking the limit $\hat{d}_k \rightarrow 0$ for $k \in \hat{R} \setminus R$ in (5.28) concludes the proof in the general case. Now Proposition 5.2 follows since ε may be chosen arbitrarily small. This concludes the proof of Proposition 5.2.

5.2. Removing the non-overlapping assumption. In this subsection we complete the proof of Proposition 5.1 by extending Proposition 5.2 to the case where (5.3) does not hold.

PROOF OF PROPOSITION 5.1. Let $\delta < \tau/4$. We say that $i, j \in \mathcal{O}_{\tau/2}^+$ *overlap* if $|d_i - d_j| \leq (d_i - 1)^{-1/2} K^{-1/2+\delta}$ or $|d_i - d_j| \leq (d_j - 1)^{-1/2} K^{-1/2+\delta}$. For $A \subset \mathcal{O}_\tau^+$ we introduce sets $S(A), L(A) \subset \mathcal{O}_{\tau/2}^+$ satisfying $S(A) \subset A \subset L(A)$. Informally, $S(A) \subset A$ is the largest subset of indices of A that do not overlap with its complement. It is by definition constructed by successively choosing $k \in A$, such that k overlaps with an index of A^c , and removing k from A ; this process is repeated until no such k exists. One can check that the result is independent of the choice of k at each step. Note that $S(A)$ may be empty.

Informally, $L(A) \supset A$ is the smallest subset of indices in $\mathcal{O}_{\tau/2}^+$ that do not overlap with its complement. It is by definition constructed by successively choosing $k \in \mathcal{O}_{\tau/2}^+ \setminus A$, such that k overlaps with an index of A , and adding k to A ; this process is repeated until no such k exists. One can check that the result is independent of the choice of k at each step. See Figure 5.2 for an illustration of $S(A)$ and $L(A)$. Throughout the following we shall repeatedly make use of the fact that, for any $A \subset \mathcal{O}_\tau^+$, Proposition 5.2 is applicable with (τ, A) replaced by $(\tau/2, S(A))$ or $(\tau/2, L(A))$.

After these preparations, we move on to the proof of (5.2). We divide the argument into four steps.

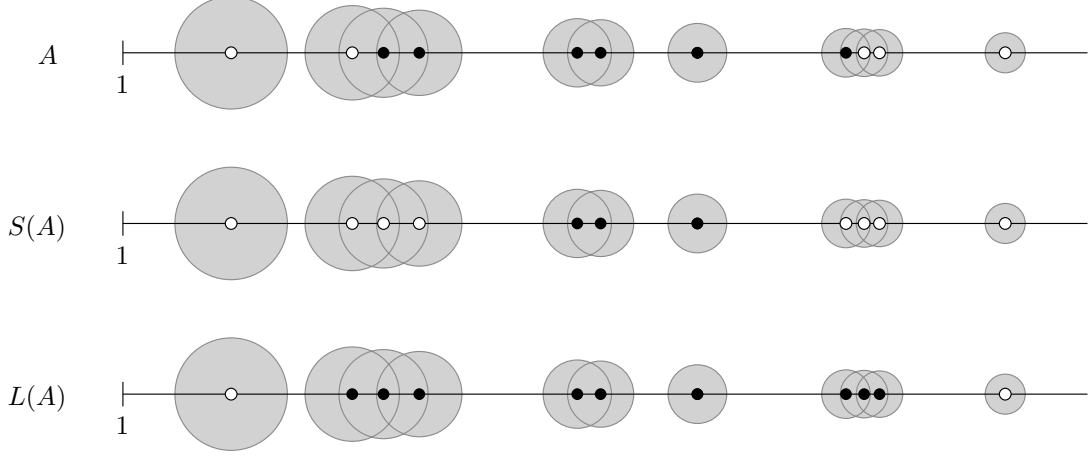


FIGURE 5.2. The construction of the sets $S(A)$ and $L(A)$. The black and white dots are the outlier indices $\{d_i : i \in \mathcal{O}_{\tau/2}^+\}$, contained in the interval $[1, \infty)$. Around each outlier index d_i we draw a grey circle of radius $(d_i - 1)^{-1/2} K^{-1/2+\delta}$. By definition, two dots overlap if one is contained in the grey circle of the other. The three pictures depict (from top to bottom) the sets A , $S(A)$, and $L(A)$, respectively. In each case, the given set is drawn using black dots and its complement using white dots.

(a) $i = j \notin A$. We consider two cases, $i \notin L(A)$ and $i \in L(A)$. Suppose first that $i \notin L(A)$. Using that $|R|$ is bounded, it is not hard to see that $\nu_i(A) \asymp \nu_i(L(A))$. We now invoke Proposition 5.2 and get

$$\langle \mathbf{v}_i, P_A \mathbf{v}_i \rangle \leq \langle \mathbf{v}_i, P_{L(A)} \mathbf{v}_i \rangle \prec \frac{\sigma_i}{M \nu_i(L(A))^2} \leq C \frac{\sigma_i}{M \nu_i(A)^2}. \quad (5.29)$$

In the complementary case, $i \in L(A)$, a simple argument yields

$$\nu_i(A) \leq C(d_i - 1)^{-1/2} K^{-1/2+\delta} \leq C \nu_i(L(A)), \quad (5.30)$$

as well as $\sigma_i \asymp 1 + \phi^{1/2}$. From Proposition 5.2 we therefore get

$$\begin{aligned} \langle \mathbf{v}_i, P_A \mathbf{v}_i \rangle &\leq \langle \mathbf{v}_i, P_{L(A)} \mathbf{v}_i \rangle \prec \frac{d_i - 1}{1 + \phi^{1/2}} + \frac{1}{(d_i - 1)^{1/2} M^{1/2}} + \frac{1 + \phi^{1/2}}{M \nu_i(L(A))^2} + \frac{1 + \phi^{1/2}}{M(d_i - 1)^2} \\ &\leq CK^{2\delta} \frac{d_i - 1}{1 + \phi^{1/2}} \leq CK^{2\delta} \frac{1 + \phi^{1/2}}{M \nu_i(A)^2} \leq CK^{2\delta} \frac{\sigma_i}{M \nu_i(A)^2}, \end{aligned}$$

where we used that $M \asymp (1 + \phi)K$. Recalling (5.29), we conclude

$$\langle \mathbf{v}_i, P_A \mathbf{v}_i \rangle \prec K^{2\delta} \frac{\sigma_i}{M \nu_i(A)^2} \quad (i \notin A). \quad (5.31)$$

(b) $i = j \in A$. We consider the two cases $i \in S(A)$ and $i \notin S(A)$. Suppose first that $i \in S(A)$. We write

$$\langle \mathbf{v}_i, P_A \mathbf{v}_i \rangle = \langle \mathbf{v}_i, P_{S(A)} \mathbf{v}_i \rangle + \langle \mathbf{v}_i, P_{A \setminus S(A)} \mathbf{v}_i \rangle. \quad (5.32)$$

We compute the first term of (5.32) using Proposition 5.2 and the observation that $\nu_i(A) \asymp \nu_i(S(A))$:

$$\langle \mathbf{v}_i, P_{S(A)} \mathbf{v}_i \rangle = u(d_i) + O_{\prec} \left[\frac{1}{(d_i - 1)^{1/2} M^{1/2}} + \frac{\sigma_i}{M} \left(\frac{1}{\nu_i(A)^2} + \frac{1}{(d_i - 1)^2} \right) \right].$$

In order to estimate the second term of (5.32), we note that $\nu_i(A) \asymp \nu_i(A \setminus S(A))$. We therefore apply (5.31) with A replaced by $A \setminus S(A)$ to get

$$\langle \mathbf{v}_i, P_{A \setminus S(A)} \mathbf{v}_i \rangle \prec K^{2\delta} \frac{\sigma_i}{M \nu_i(A)^2}.$$

Going back to (5.32), we have therefore proved that

$$\langle \mathbf{v}_i, P_A \mathbf{v}_i \rangle = u(d_i) + K^{2\delta} O_{\prec} \left[\frac{1}{(d_i - 1)^{1/2} M^{1/2}} + \frac{\sigma_i}{M} \left(\frac{1}{\nu_i(A)^2} + \frac{1}{(d_i - 1)^2} \right) \right] \quad (5.33)$$

for $i \in S(A)$.

Next, we consider the case $i \notin S(A)$. Now we have (5.30), so that Proposition 5.2 yields

$$\langle \mathbf{v}_i, P_A \mathbf{v}_i \rangle \leq \langle \mathbf{v}_i, P_{L(A)} \mathbf{v}_i \rangle \prec u(d_i) + \frac{1}{(d_i - 1)^{1/2} M^{1/2}} + \frac{\sigma_i}{M} \left(\frac{1}{\nu_i(A)^2} + \frac{1}{(d_i - 1)^2} \right).$$

By (5.30) and $M \asymp (1 + \phi)K$, we have

$$u(d_i) = \frac{\sigma_i}{\phi^{1/2} \theta(d_i)} (1 - d_i^{-2}) \leq CK^{2\delta} \frac{1}{(d_i - 1)^{1/2} M^{1/2}},$$

from which we deduce (5.33) also in the case $i \notin S(A)$.

(c) $i \notin A$ or $j \notin A$ or $i = j$. From cases (a) and (b) (i.e. (5.31) and (5.33)), combined with the estimate

$$|\langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle|^2 \leq \langle \mathbf{v}_i, P_A \mathbf{v}_i \rangle \langle \mathbf{v}_j, P_A \mathbf{v}_j \rangle,$$

we find, assuming $i \notin A$, $j \notin A$, or $i = j$, that (5.2) holds with an additional factor $K^{2\delta}$ multiplying the right-hand side.

(d) $i, j \in A$ and $i \neq j$. We now deal with the last remaining case by using the splitting

$$\langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle = \langle \mathbf{v}_i, P_{S(A)} \mathbf{v}_j \rangle + \langle \mathbf{v}_i, P_{A \setminus S(A)} \mathbf{v}_j \rangle. \quad (5.34)$$

The goal is to show that

$$|\langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle| \prec \frac{K^{2\delta}}{(d_i - 1)^{1/4} (d_j - 1)^{1/4} M^{1/2}} + K^{2\delta} \frac{\sqrt{\sigma_i \sigma_j}}{M} \left(\frac{1}{\nu_i(A)} + \frac{1}{d_i - 1} \right) \left(\frac{1}{\nu_j(A)} + \frac{1}{d_j - 1} \right). \quad (5.35)$$

Note that here $\sigma_i \asymp \sigma_j \asymp 1 + \phi^{1/2}$. We consider the four cases (i) $i, j \in S(A)$, (ii) $i \in S(A)$ and $j \notin S(A)$, (iii) $i \notin S(A)$ and $j \in S(A)$, and (iv) $i, j \notin S(A)$.

Consider first the case (i). The first term of (5.34) is bounded using Proposition 5.2 combined with $\nu_i(A) \asymp \nu_i(S(A))$ and $\nu_j(A) \asymp \nu_j(S(A))$. The second term of (5.34) is bounded using (5.2) from case (c) combined with $\nu_i(A) \leq C\nu_i(A \setminus S(A))$ and $\nu_j(A) \leq C\nu_j(A \setminus S(A))$. This yields (5.35) for $\langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle$ in the case (i).

Next, consider the case (ii). For the first term of (5.34) we use the estimates

$$\nu_i(S(A)) \asymp \nu_i(A), \quad \nu_j(A) \leq C(d_j - 1)^{-1/2} K^{-1/2+\delta} \leq C\nu_j(S(A)), \quad \nu_i(A) \leq C|d_i - d_j|.$$

Thus we get from (5.4)

$$\begin{aligned} |\langle \mathbf{v}_i, P_{S(A)} \mathbf{v}_j \rangle| &\prec \frac{1 + \phi^{1/2}}{M\nu_i(S(A))\nu_j(S(A))} + \frac{1 + \phi^{1/2}}{M\nu_j(S(A))(d_i - 1)} + \frac{(d_i - 1)^{1/2}}{M^{1/2}|d_i - d_j|} \\ &\leq \frac{1 + \phi^{1/2}}{M\nu_i(A)\nu_j(A)} + \frac{1 + \phi^{1/2}}{M\nu_j(A)(d_i - 1)} + \frac{(d_i - 1)^{1/2}}{M^{1/2}|d_i - d_j|}. \end{aligned}$$

In order to estimate the last term, we first assume that $d_j \leq d_i$ and $d_i - 1 \leq 2|d_i - d_j|$. Then we find

$$\frac{(d_i - 1)^{1/2}}{M^{1/2}|d_i - d_j|} \leq \frac{2}{M^{1/2}(d_i - 1)^{1/2}} \leq \frac{2}{M^{1/2}(d_i - 1)^{1/4}(d_j - 1)^{1/4}}.$$

Conversely, if $d_i \leq d_j$ or $d_i - 1 \geq 2|d_i - d_j|$, we have $d_i - 1 \leq 2(d_j - 1)$. Thus we get

$$\frac{(d_i - 1)^{1/2}}{M^{1/2}|d_i - d_j|} \leq \frac{C(d_j - 1)^{1/2}}{M^{1/2}\nu_i(A)} \leq K^\delta \frac{1 + \phi^{1/2}}{M\nu_i(A)\nu_j(A)}.$$

Putting both estimates together, we may estimate the first term of (5.34) in the case (ii) as

$$|\langle \mathbf{v}_i, P_{S(A)} \mathbf{v}_j \rangle| \prec \frac{1 + \phi^{1/2}}{M\nu_i(A)\nu_j(A)} + K^\delta \frac{1 + \phi^{1/2}}{M\nu_j(A)(d_i - 1)} + \frac{1}{M^{1/2}(d_i - 1)^{1/4}(d_j - 1)^{1/4}}. \quad (5.36)$$

For the second term of (5.34) in the case (ii) we use the estimates

$$\nu_i(A \setminus S(A)) \asymp \nu_i(A), \quad \nu_j(A) \leq C\nu_j(A \setminus S(A)) \leq C(d_j - 1)^{-1/2}K^{-1/2+\delta}, \quad \nu_i(A) \leq C|d_i - d_j|.$$

Thus we get from case (c) that

$$|\langle \mathbf{v}_i, P_{A \setminus S(A)} \mathbf{v}_j \rangle| \leq \frac{1 + \phi^{1/2}}{M\nu_i(A)} \left(\frac{1}{\nu_j(A)} + \frac{1}{d_j - 1} \right) + \frac{(d_j - 1)^{1/2}}{M^{1/2}|d_i - d_j|},$$

where the last term is bounded by $K^\delta \frac{1 + \phi^{1/2}}{M\nu_i(A)\nu_j(A)}$. Recalling (5.36), we find (5.35) in the case (ii). The case (iii) is dealt with in the same way.

What remains therefore is case (iv). For the first term of (5.34) we use the estimates

$$\nu_i(A) \leq C\nu_i(S(A)), \quad \nu_j(A) \leq C\nu_j(S(A)).$$

Thus we get from (5.4) that

$$\langle \mathbf{v}_i, P_{S(A)} \mathbf{v}_j \rangle \leq \frac{1 + \phi^{1/2}}{M\nu_i(A)\nu_j(A)}.$$

For the second term of (5.34) we use the estimates

$$\nu_i(A) \leq C\nu_i(A \setminus S(A)) \leq C(d_i - 1)^{-1/2}K^{-1/2+\delta}, \quad \nu_j(A) \leq C\nu_j(A \setminus S(A)) \leq C(d_j - 1)^{-1/2}K^{-1/2+\delta}.$$

Therefore we get from case (c) that

$$\langle \mathbf{v}_i, P_{A \setminus S(A)} \mathbf{v}_j \rangle \prec \frac{d_i - 1}{1 + \phi^{1/2}} + \frac{1}{(d_i - 1)^{1/2}M^{1/2}} + \frac{1 + \phi^{1/2}}{M} \left(\frac{1}{\nu_i(A \setminus S(A))^2} + \frac{1}{(d_i - 1)^2} \right) \leq CK^{2\delta} \frac{1 + \phi^{1/2}}{M\nu_i(A)^2},$$

and a similar estimate holds for $\langle \mathbf{v}_j, P_{A \setminus S(A)} \mathbf{v}_j \rangle$. Thus we conclude that

$$|\langle \mathbf{v}_i, P_{A \setminus S(A)} \mathbf{v}_j \rangle| \leq \langle \mathbf{v}_i, P_{A \setminus S(A)} \mathbf{v}_i \rangle^{1/2} \langle \mathbf{v}_j, P_{A \setminus S(A)} \mathbf{v}_j \rangle^{1/2} \leq CK^{2\delta} \frac{1 + \phi^{1/2}}{M\nu_i(A)\nu_j(A)},$$

which is (5.35). This concludes the analysis of case (iv), and hence of case (d).

Conclusion of the proof. Putting the cases (a)–(d) together, we have proved that (5.2) holds for arbitrary i, j with an additional factor $K^{2\delta}$ multiplying the error term on the right-hand side. Since $\delta > 0$ can be chosen arbitrarily small, (5.2) follows. This concludes the proof of Proposition 5.1. $\square$

6. Non-outlier eigenvectors

In this section we focus on the non-outlier eigenvectors ξ_a , $a \notin \mathcal{O}$, as well as outlier eigenvectors close to the bulk spectrum. We derive isotropic delocalization bounds for ξ_a and establish the asymptotic law of the generalized components of ξ_a . We also use the former result to complete the proof of Theorem 2.9 on the outlier eigenvectors.

In Section 6.1 we derive isotropic delocalization bounds on ξ_a for $\text{dist}\{d_a, [-1, 1]\} \leq 1 + K^{-1/3+\tau}$. In Section 6.2 we use these bounds to prove Theorem 2.9 and to complete the proof of Theorem 2.15 started in Section 5. Next, in Section 6.3 we derive the law of the generalized components of ξ_a for $a \notin \mathcal{O}$. This argument requires two tools as input: level repulsion (Proposition 6.4) and quantum unique ergodicity of the eigenvalues ζ_b of H (Proposition 6.7). Both are explained in detail and proved below.

6.1. Bound on the spectral projections in the neighbourhood of the bulk spectrum. We first consider eigenvectors near the right edge of the bulk spectrum.

PROPOSITION 6.1 (EIGENVECTORS NEAR THE RIGHT EDGE). *Fix $\tau \in (0, 1/3)$. For $a \in \llbracket s_+ + 1, (1 - \tau)K \rrbracket$ we have*

$$\langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2 \prec \frac{1}{M} + \frac{\sigma_i}{M(|d_i - 1|^2 + \kappa_a)}. \quad (6.1)$$

Moreover, if $a \in \llbracket 1, s_+ \rrbracket$ satisfies $d_a \leq 1 + K^{-1/3+\tau}$ then

$$\langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2 \prec K^{3\tau} \left(\frac{1}{M} + \frac{\sigma_i}{M(|d_i - 1|^2 + \kappa_a)} \right). \quad (6.2)$$

Proposition 6.1 has the following analogue for the left edge of the bulk spectrum.

PROPOSITION 6.2 (EIGENVECTORS NEAR THE LEFT EDGE). *Fix $\tau \in (0, 1/3)$ and suppose that $|\phi - 1| \geq \tau$. For $a \in \llbracket \tau K, K - s_- \rrbracket$ we have*

$$\langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2 \prec \frac{1}{M} + \frac{\sigma_i}{M(|d_i + 1|^2 + \kappa_a)}. \quad (6.3)$$

Moreover, if $a \in \llbracket K - s_- + 1, K \rrbracket$ satisfies $d_a \geq -1 - K^{-1/3+\tau}$ then

$$\langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2 \prec K^{3\tau} \left(\frac{1}{M} + \frac{\sigma_i}{M(|d_i + 1|^2 + \kappa_a)} \right). \quad (6.4)$$

The proofs of Propositions 6.1 and 6.2 are analogous, and we only prove the former.

PROOF OF PROPOSITION 6.1. Suppose first that $i \in R$. Let $\varepsilon > 0$ and set $\omega := \varepsilon/2$. Using (3.23), Remark 3.3, Theorem 2.3, and (3.14), we choose a high-probability event Ξ satisfying (4.7), (4.8), and

$$\mathbf{1}(\Xi) |\mu_i - \theta(d_i)| \leq (d_i - 1)^{1/2} K^{-1/2+\varepsilon} \quad (1 + K^{-1/3} \leq d_i \leq 1 + K^{-1/3+\tau}). \quad (6.5)$$

For the following we fix a realization $H \in \Xi$. We choose the spectral parameter $z = \mu_a + \eta$, where $\eta > 0$ is the smallest (in fact unique) solution of the equation $\text{Im } w_\phi(\mu_a + i\eta) = K^{-1+6\varepsilon}\eta^{-1}$. Hence (4.8) reads

$$\|W(z) - w_\phi(z)\| \leq \frac{K^{2\varepsilon}}{K\eta}. \quad (6.6)$$

Abbreviating $\kappa \equiv \kappa(\mu_a)$, we find from (3.19) that

$$\eta \asymp \frac{K^{6\varepsilon}}{K\sqrt{\kappa} + K^{2/3+2\varepsilon}} \quad (\mu_a \leq \gamma_+ + K^{-2/3+4\varepsilon}) \quad (6.7)$$

and

$$\eta \asymp K^{-1/2+3\varepsilon} \kappa^{1/4} \quad (\mu_a \geq \gamma_+ + K^{-2/3+4\varepsilon}). \quad (6.8)$$

Armed with these definitions, we may begin the estimate of $\langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2$. The starting point is the bound

$$\langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2 \leq \eta \text{Im } \tilde{G}_{\mathbf{v}_i \mathbf{v}_i}(z), \quad (6.9)$$

which follows easily by spectral decomposition. Since $i \in R$, we get from (3.35), omitting the arguments z for brevity,

$$\begin{aligned} \phi^{1/2} z \tilde{G}_{\mathbf{v}_i \mathbf{v}_i} &= \frac{1}{d_i} - \frac{\sigma_i}{d_i^2} \left(\frac{1}{D^{-1} + W} \right)_{ii} \\ &= \frac{1}{d_i} - \frac{\sigma_i}{d_i^2} \left[\frac{1}{d_i^{-1} + w_\phi} + \frac{1}{(d_i^{-1} + w_\phi)^2} \left((w_\phi - W) + (w_\phi - W) \frac{1}{D^{-1} + W} (w_\phi - W) \right)_{ii} \right], \end{aligned} \quad (6.10)$$

where the last step follows from a resolvent expansion as in (5.14). We estimate the error terms using

$$\min_j |d_j^{-1} + w_\phi| \geq \operatorname{Im} m_\phi = \frac{K^{6\varepsilon}}{K\eta} \gg \frac{K^{2\varepsilon}}{K\eta} \geq \|W - w_\phi\|,$$

where we used the definition of η and (6.6). Hence a resolvent expansion yields

$$\left\| \frac{1}{D^{-1} + W} \right\| \leq \frac{2}{\operatorname{Im} w_\phi} = 2K^{1-6\varepsilon}\eta.$$

We therefore get from (6.10) that

$$\phi^{1/2} z \tilde{G}_{\mathbf{v}_i \mathbf{v}_i} = \frac{w_\phi - \phi^{1/2}}{1 + d_i w_\phi} + O\left(\frac{\sigma_i}{|1 + d_i w_\phi|^2} \frac{K^{2\varepsilon}}{K\eta}\right) \quad (6.11)$$

Next, we claim that for any fixed $\delta \in [0, 1/3 - \varepsilon)$ we have the lower bound

$$|1 + dw_\phi| \geq c(K^{-2\delta}|d-1| + \operatorname{Im} w_\phi) \quad (6.12)$$

whenever $\mu_a \in [\theta(0), \theta(1 + K^{-1/3+\delta+\varepsilon})]$. To prove (6.12), suppose first that $|d-1| \geq 1/2$. By (3.20), there exists a constant $c_0 > 0$ such that for $\kappa \leq c_0$ we have $|\operatorname{Re} w_\phi + 1| \leq 1/4$. Thus we get, for $\kappa \leq c_0$,

$$|1 + dw_\phi| \geq |1 + d \operatorname{Re} w_\phi| \asymp |d-1| + \operatorname{Im} w_\phi,$$

where we used that $\operatorname{Im} w_\phi \leq C$ by (3.18). Moreover, if $\kappa \geq c_0$ we find from (3.19) that $\operatorname{Im} w_\phi \geq c$, from which we get

$$|1 + dw_\phi| \geq |1 + d \operatorname{Re} w_\phi| + |d| \operatorname{Im} w_\phi \geq c(1 + |d|) \asymp |d-1| + \operatorname{Im} w_\phi,$$

where in the second step we used $|\operatorname{Re} w_\phi| \leq C$ as follows from (3.18). This concludes the proof of (6.12) for the case $|d-1| \geq 1/2$.

Suppose now that $|d-1| \leq 1/2$. Then we get

$$|1 + dw_\phi| \asymp |1 + d \operatorname{Re} w_\phi| + |d| \operatorname{Im} w_\phi \geq (|d-1| - |\operatorname{Re} w_\phi + 1|)_+ + \operatorname{Im} w_\phi.$$

We shall estimate this using the elementary bound

$$(x-y)_+ + z \geq \frac{x}{3M} + \frac{z}{3} \quad \text{if } y \leq Mz \text{ for some } M \geq 1. \quad (6.13)$$

For $\mu_a \in [\theta(0), \theta(1)]$ we get from (6.13) with $M = C$, recalling (3.19) and (3.20), that $|1 + dw_\phi| \geq c(|d-1| + \operatorname{Im} w_\phi)$. By a similar argument, for $\mu_a \in [\theta(1), \theta(1 + K^{-1/3+\delta+\varepsilon})]$ we set $M = K^{2\delta}$ and get (6.12) using (6.7) and (6.8). This concludes the proof of (6.12).

Going back to (6.9), we find using (6.11)

$$\begin{aligned} \langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2 &\leq \eta \phi^{-1/2} \operatorname{Im} \left(\frac{w_\phi - \phi^{1/2}}{z(1 + d_i w_\phi)} \right) + \frac{C \sigma_i}{|1 + d_i w_\phi|^2} \frac{K^{2\varepsilon}}{K\eta} \\ &= \frac{\eta^2}{\phi^{1/2}|z|^2} \operatorname{Re} \frac{\phi^{1/2} - w_\phi}{1 + d_i w_\phi} + \frac{\eta \mu_a}{\phi^{1/2}|z|^2} \operatorname{Im} \frac{w_\phi - \phi^{1/2}}{1 + d_i w_\phi} + \frac{C \sigma_i}{\phi^{1/2}|z||1 + d_i w_\phi|^2} \frac{K^{2\varepsilon}}{K}. \end{aligned} \quad (6.14)$$

Using $|z| \asymp \mu_a \asymp \phi^{-1/2} + \phi^{1/2}$ and (6.12), we estimate the first term on the right-hand side of (6.14) as

$$\frac{\eta^2}{\phi^{1/2}|z|^2} \operatorname{Re} \frac{\phi^{1/2} - w_\phi}{1 + d_i w_\phi} \leq \frac{\eta^2}{(1+\phi)|1 + d_i w_\phi|} \leq \frac{\eta^2}{(1+\phi) \operatorname{Im} w_\phi} = \frac{\eta^3 K}{1+\phi} \leq \frac{K^{12\varepsilon+3\delta}}{M},$$

where in the last step we used that $\eta \leq K^{-2/3+4\varepsilon+\delta}$, as follows from (6.7) and (6.8).

Next, we estimate the second term of (6.14) as

$$\frac{\eta\mu_a}{\phi^{1/2}|z|^2} \operatorname{Im} \frac{w_\phi - \phi^{1/2}}{1 + d_i w_\phi} = \frac{\eta\mu_a}{\phi^{1/2}|z|^2} \frac{\sigma_i \operatorname{Im} w_\phi}{|1 + d_i w_\phi|^2} \asymp \frac{\sigma_i \eta \operatorname{Im} w_\phi}{(1 + \phi)|1 + d_i w_\phi|^2} \leq \frac{C\sigma_i K^{6\varepsilon}}{M|1 + d_i w_\phi|^2}.$$

We estimate the last term of (6.14) as

$$\frac{C\sigma_i}{\phi^{1/2}|z||1 + d_i w_\phi|^2} \frac{K^{2\varepsilon}}{K} \leq \frac{C\sigma_i K^{2\varepsilon}}{M|1 + d_i w_\phi|^2}.$$

Putting all three estimates together, we conclude that

$$\langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2 \leq \frac{K^{12\varepsilon+3\delta}}{M} + \frac{C\sigma_i K^{6\varepsilon}}{M|1 + d_i w_\phi|^2}. \quad (6.15)$$

In order to estimate the denominator of (6.15) from below using (6.12), we need a suitable lower bound on $\operatorname{Im} w_\phi(\mu_a + i\eta)$. First, if $a \geq s_+ + 1$ then we get from (4.7), Corollary 4.2, (6.7), and (3.19) that

$$\operatorname{Im} w_\phi(\mu_a + i\eta) \geq c\sqrt{\kappa_a},$$

in which case we get by choosing $\delta = 0$ in (6.12) that

$$\langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2 \leq \frac{K^{12\varepsilon}}{M} + \frac{C\sigma_i K^{6\varepsilon}}{M(|d_i - 1|^2 + \kappa_a)}. \quad (6.16)$$

Next, if $a \leq s_+$ satisfies $d_a \leq K^{-1/3+\tau}$ we get from (6.7), (6.8), and (3.19) that

$$\operatorname{Im} w_\phi(\mu_a + i\eta) \geq c\sqrt{\eta} \geq cK^{-1/3+2\varepsilon}.$$

In this case we have $\mu_a \leq \theta(1 + K^{-1/3+\tau+\varepsilon})$ by (6.5), so that setting $\delta = \tau$ in (6.12) yields

$$\langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2 \leq \frac{K^{12\varepsilon+3\tau}}{M} + \frac{C\sigma_i K^{6\varepsilon+2\tau}}{M(|d_i - 1|^2 + \kappa_a)}. \quad (6.17)$$

Since $\varepsilon > 0$ was arbitrary, (6.1) and (6.2) follow from (6.16) and (6.17) respectively. This concludes the proof of Proposition 6.1 in the case $i \in R$.

Finally, the case $i \notin R$ is handled by replacing R with $\widehat{R} := R \cup \{i\}$ and using a limiting argument, exactly as after (5.28). $\square$

6.2. Proof of Theorems 2.9 and 2.15. We now have all the ingredients needed to prove Theorems 2.9 and 2.15.

PROOF OF THEOREM 2.15. This is an immediate corollary of (6.1) and (6.3) from Propositions 6.1 and 6.2. $\square$

PROOF OF THEOREM 2.9. We prove Theorem 2.9 using Propositions 5.1, 5.2, and 6.1. First we remark that it suffices to prove that (5.2) holds for $A \subset \mathcal{O}$ satisfying $1 + K^{-1/3} \leq d_k \leq \tau^{-1}$ for all $k \in A$. Indeed, supposing this is done, we get the estimate

$$\begin{aligned} \langle \mathbf{w}, P_A \mathbf{w} \rangle &= \langle \mathbf{w}, T_A \mathbf{w} \rangle + O_{\prec} \left[\sum_{i \in A} \frac{w_i^2}{M^{1/2}(d_i - 1)^{1/2}} + \sum_{i \in A} \frac{\sigma_i w_i^2}{M(d_i - 1)^2} \right. \\ &\quad \left. + \sum_{i=1}^M \frac{\sigma_i w_i^2}{M\nu_i^2} + \langle \mathbf{w}, T_A \mathbf{w} \rangle^{1/2} \left(\sum_{i \notin A} \frac{\sigma_i w_i^2}{M\nu_i^2} \right)^{1/2} \right], \end{aligned}$$

from which Theorem 2.9 follows by noting that the second error term may be absorbed into the first, recalling that $\sigma_i \asymp 1 + \phi^{1/2}$ for $i \in A$, that $M \asymp (1 + \phi)K$, and that $d_i - 1 \geq K^{-1/3}$.

Fix $\varepsilon > 0$. Note that there exists some $s \in [1, |R|]$ satisfying the following gap condition: for all k such that $d_k > 1 + sK^{-1/3+\varepsilon}$ we have $d_k \geq 1 + (s+1)K^{-1/3+\varepsilon}$. The idea of the proof is to split $A = A_0 \sqcup A_1$, such that $d_k \leq 1 + sK^{-1/3+\varepsilon}$ for $k \in A_0$ and $d_k \geq 1 + (s+1)K^{-1/3+\varepsilon}$ for $k \in A_1$. Note that such a splitting exists by the above gap property. Without loss of generality, we assume that $A_0 \neq \emptyset$ (for otherwise the claim follows from Proposition 5.1).

It suffices to consider the six cases (a) $i, j \in A_0$, (b) $i \in A_0$ and $j \in A_1$, (c) $i \in A_0$ and $j \notin A$, (d) $i, j \in A_1$, (e) $i \in A_1$ and $j \notin A$, (f) $i, j \notin A$.

(a) $i, j \in A_0$. We split

$$\langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle = \langle \mathbf{v}_i, P_{A_0} \mathbf{v}_j \rangle + \langle \mathbf{v}_i, P_{A_1} \mathbf{v}_j \rangle. \quad (6.18)$$

We apply Cauchy-Schwarz and Proposition 6.1 to the first term, and Proposition 5.1 to the second term. Using the above gap condition, we find

$$\begin{aligned} |\langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle| &\prec \frac{K^{3\varepsilon} \sqrt{\sigma_i \sigma_j}}{M|d_i - 1||d_j - 1|} + \frac{\sqrt{\sigma_i \sigma_j}}{M} \left(\frac{1}{\nu_i} + \frac{1}{d_i - 1} \right) \left(\frac{1}{\nu_j} + \frac{1}{d_j - 1} \right) \\ &= \delta_{ij} u(d_i) + K^{3\varepsilon} O \left[\frac{1}{(d_i - 1)^{1/4} (d_j - 1)^{1/4} M^{1/2}} + \frac{\sqrt{\sigma_i \sigma_j}}{M} \left(\frac{1}{\nu_i} + \frac{1}{d_i - 1} \right) \left(\frac{1}{\nu_j} + \frac{1}{d_j - 1} \right) \right], \end{aligned}$$

where the last step follows from $d_i - 1 \leq CK^{-1/3+\varepsilon}$.

(b) $i \in A_0$ and $j \in A_1$. For this case it is crucial to use the stronger bound (5.4) and not (5.2). Hence, we need the non-overlapping condition (5.3). To that end, we assume first that (5.3) holds with $\delta := \varepsilon$. Thus, by the above gap assumption (5.3) also holds for A_1 . In this case we get from (6.18) and Propositions 5.2 and 6.1 that

$$|\langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle| \prec \frac{K^{3\varepsilon} \sqrt{\sigma_i \sigma_j}}{M|d_i - 1||d_j - 1|} + \frac{\sqrt{\sigma_i \sigma_j}}{M} \left(\frac{1}{\nu_i} + \frac{1}{d_i - 1} \right) \left(\frac{1}{\nu_j} + \frac{1}{d_j - 1} \right) + \frac{(d_j - 1)^{1/2} \sqrt{\sigma_i}}{(1 + \phi^{1/4})|d_i - d_j|M^{1/2}}.$$

Clearly, the first two terms are bounded by the right-hand side of (5.2) times $K^{3\varepsilon}$. The last term is estimated as

$$\frac{(d_j - 1)^{1/2} \sqrt{\sigma_i}}{(1 + \phi^{1/4})|d_i - d_j|M^{1/2}} \asymp \frac{(d_j - 1)^{1/2}}{|d_i - d_j|M^{1/2}} \leq \frac{1}{(d_i - 1)^{1/4} (d_j - 1)^{1/4} M^{1/2}},$$

where we used that $d_i - 1 \leq d_j - 1 \leq C|d_i - d_j|$ by the above gap condition. This concludes the proof in the case where the non-overlapping condition (5.3) holds.

If (5.3) does not hold, we replace A_1 with the smaller set $S(A_1)$ defined in Section 5.2. Then we proceed as above, except that we have to deal in addition with the term $\langle \mathbf{v}_i, P_{A_1 \setminus S(A_1)} \mathbf{v}_j \rangle$. The details are analogous to those of Section 5.2, and we omit them here.

(c), (e), (f) $j \notin A$. We use the splitting (6.18) and apply Cauchy-Schwarz and Proposition 6.1 to the first term, and Proposition 5.1 to the second term. Since $\nu_j(A_1) \leq |d_j - 1|$ in all cases, it is easy to prove that (6.18) is bounded by $K^{3\varepsilon}$ times the right-hand side of (5.2).

(d) $i, j \in A_1$. From (6.18) and Propositions 6.1 and 5.1 we get

$$\begin{aligned} \langle \mathbf{v}_i, P_A \mathbf{v}_j \rangle &= \delta_{ij} u(d_i) \\ &+ O \prec \left[\frac{K^{3\varepsilon} \sqrt{\sigma_i \sigma_j}}{M|d_i - 1||d_j - 1|} + \frac{1}{(d_i - 1)^{1/4} (d_j - 1)^{1/4} M^{1/2}} + \frac{\sqrt{\sigma_i \sigma_j}}{M} \left(\frac{1}{\nu_i} + \frac{1}{d_i - 1} \right) \left(\frac{1}{\nu_j} + \frac{1}{d_j - 1} \right) \right], \end{aligned}$$

From which we get (5.2) with the error term multiplied by $K^{3\varepsilon}$.

Conclusion of the proof. We have proved that, for all $i, j \in \llbracket 1, M \rrbracket$ and A satisfying the assumptions of Theorem 2.9, the estimate (5.2) holds with an additional factor $K^{3\varepsilon}$ multiplying the error term. Since ε was arbitrary, we get (5.2). This concludes the proof. $\square$

6.3. The law of the non-outlier eigenvectors. For $a \leq K/2$ define

$$\Delta_a := K^{-2/3} a^{-1/3}, \quad (6.19)$$

the typical distance between λ_{a+1} and λ_a . More precisely, the classical locations γ_a defined in (3.13) satisfy $\gamma_a - \gamma_{a+1} \asymp \Delta_a$ for $a \leq K/2$.

We may now state the main result behind the proof of Theorem 2.18. Recall the definitions (2.13) of α_+ and (2.8) of s_+ , the number of outliers to the right of the bulk spectrum.

PROPOSITION 6.3. *Let $s_+ + 1 \leq a \leq K^{1-\tau} \alpha_+^3$ and define $b := a - s_+$. Define the event*

$$\Omega \equiv \Omega_{a,b,\tau} := \left\{ |\mu_{b'+s_+} - \lambda_{b'}| \leq K^{-\tau/4} \Delta_a \text{ for } |b' - b| \leq 1 \right\} \cap \left\{ |\lambda_{b'} - \lambda_b| \geq K^{-\tau/6} \Delta_a \text{ for } |b' - b| = 1 \right\}.$$

Then

$$\mathbf{1}(\Omega) \langle \mathbf{w}, \boldsymbol{\xi}_a \rangle^2 = \mathbf{1}(\Omega) \left| \left\langle \sum_i \frac{\sqrt{\sigma_i} w_i}{d_i - 1} \mathbf{v}_i, \boldsymbol{\zeta}_b \right\rangle \right|^2 + O_{\prec} \left[\frac{K^{-1/3+\tau/5} a^{1/3}}{\alpha_+} \sum_i \frac{\sigma_i w_i^2}{M(d_i - 1)^2} \right]. \quad (6.20)$$

Informally, Proposition 6.3 expresses generalized components of the eigenvectors of Q in terms of generalized components of eigenvectors of H , under the assumption that Ω has high probability. We first show how Proposition 6.3 implies Theorem 2.18. This argument requires two key tools. The first one is level repulsion, which, together with the eigenvalue sticking from Theorem 2.6, will imply that Ω indeed has high probability. The second tool is quantum unique ergodicity of the eigenvectors of H , which establishes the law of the generalized components of the eigenvectors of H .

The precise statement of level repulsion sufficient for our needs is as follows.

PROPOSITION 6.4 (LEVEL REPULSION). *Fix $\tau \in (0, 1)$. For any $\varepsilon > 0$ there exists a $\delta > 0$ such that for all $a \leq K^{1-\tau}$ we have*

$$\mathbb{P}(|\lambda_a - \lambda_{a+1}| \leq \Delta_a K^{-\varepsilon}) \leq K^{-\delta}. \quad (6.21)$$

The proof of Proposition 6.4 consists of two steps: (i) establishing (6.21) for the case of Gaussian X and (ii) a comparison argument showing that if $X^{(1)}$ and $X^{(2)}$ are two matrix ensembles satisfying (2.4) and (2.5), and if (6.21) holds for $X^{(1)}$, then (6.21) also holds for $X^{(2)}$. Both steps have already appeared, in a somewhat different form, in the literature. Step (i) is performed in Lemma 6.5 below, and step (ii) in Lemma 6.6 below. Together, Lemmas 6.5 and 6.6 immediately yield Proposition 6.4.

LEMMA 6.5 (LEVEL REPULSION FOR THE GAUSSIAN CASE). *Proposition 6.4 holds if X is Gaussian.*

PROOF. We mimic the proof of Theorem 3.2 in [12]. Indeed, the proof from [12, Appendix D] carries over almost verbatim. The key input is the eigenvalue rigidity from Theorem 3.5, which for the model of [12] was established using a different method than Theorem 3.5. As in [12], we condition on the eigenvalues $\{\lambda_i : i > K^{1-\tau}\}$. On the conditioned measure, level repulsion follows as in [12]. Finally, thanks to Theorem 3.5 we know that the frozen eigenvalues $\{\lambda_i : i > K^{1-\tau}\}$ are with high probability near their classical locations. Note that for $\phi \approx 1$, the rigidity estimate (3.14) only holds for indices $i \leq (1 - \tau)K$; however, this is enough for the argument of [12, Appendix D], which is insensitive to the locations of eigenvalues at a distance of order one from the right edge γ_+ . We omit the full details. $\square$

LEMMA 6.6 (STABILITY OF LEVEL REPULSION). *Let $X^{(1)}$ and $X^{(2)}$ be two matrix ensembles satisfying (2.4) and (2.5). Suppose that Proposition 6.4 holds for $X^{(1)}$. Then Proposition 6.4 also holds for $X^{(2)}$.*

The proof of Lemma 6.6 relies on Green function comparison, and is given in Section 7.4.

The second tool behind the proof of Theorem 2.18 is the quantum unique ergodicity of the eigenvectors $\boldsymbol{\zeta}_a$ of the matrix $H = XX^*$, stated in Proposition 6.7 below. The first result on quantum unique ergodicity of Wigner matrices is [27], where the quantum unique ergodicity of eigenvectors near the spectral edge was established. Under an additional four-moment matching condition, this result was extended to the bulk. Subsequently, this second result was derived using a different method in [38]. Recently, a new approach

to the quantum unique ergodicity was developed in [13], where quantum unique ergodicity is established for all eigenvectors of generalized Wigner matrices. In this paper, we adopt the approach of [27], based on Green function comparison. As compared to the method of [13], its first advantage is that it is completely local in the spectrum, and in particular when applied near the right-hand edge of the spectrum it is insensitive to the presence of a hard edge at the origin. The second advantage of the current method is that it is very robust and may be used to establish the asymptotic joint distribution of an arbitrary family of generalized components of eigenvectors, as in Remark 6.8 below; we remark that such joint laws cannot currently be analysed using the method of [13]. On the other hand, our results only hold for eigenvector indices a satisfying $a \leq K^{1-\tau}$ for some $\tau > 0$, while those of [13] admit $\tau = 0$.

Our proof of quantum unique ergodicity generalizes that of [27] in three directions. First, we extend the method of [27] to sample covariance matrices (in fact to general sample covariance matrices of the form (2.1) with $\Sigma = TT^* = I_M$; see Section 8). Second, we consider generalized components $\langle \mathbf{w}, \zeta_a \rangle$ of the eigenvectors instead of the cartesian components $\zeta_a(i)$. The third and deepest generalization is that we establish quantum unique ergodicity much further into the bulk, requiring only that $a \leq K^{1-\tau}$ instead of the assumption $a \leq (\log K)^{C \log \log K}$ from [27].

PROPOSITION 6.7 (QUANTUM UNIQUE ERGODICITY). *Fix $\tau \in (0, 1)$. Then for any $a \leq K^{1-\tau}$ and deterministic unit vector $\mathbf{w} \in \mathbb{R}^M$ we have*

$$M \langle \mathbf{w}, \zeta_a \rangle^2 \rightarrow \chi_1^2, \quad (6.22)$$

in the sense of moments, uniformly in a and $\mathbf{w}$.

REMARK 6.8. For simplicity, and bearing the application to Theorem 2.18 in mind, in Proposition 6.7 we establish the convergence of a single generalized component of a single eigenvector. However, our method may be easily extended to yield

$$\left(M \langle \mathbf{v}_1, \zeta_{a_1} \rangle \langle \zeta_{a_1}, \mathbf{w}_1 \rangle, \dots, M \langle \mathbf{v}_k, \zeta_{a_k} \rangle \langle \zeta_{a_k}, \mathbf{w}_k \rangle \right) \stackrel{d}{\sim} (Z_1, \dots, Z_k)$$

for any deterministic unit vectors $\mathbf{v}_1, \dots, \mathbf{v}_l, \mathbf{w}_1, \dots, \mathbf{w}_k \in \mathbb{R}^M$ and $a_1, \dots, a_k \leq K^{1-\tau}$, whereby we use the notation $A_N \stackrel{d}{\sim} B_N$ to mean that A_N and B_N are tight, and $\lim_{N \rightarrow \infty} \mathbb{E}(f(A_N) - f(B_N)) = 0$ for all polynomially bounded and continuous f . Here $(Z_1, \dots, Z_k)$ is a family of independent random variables defined by $Z_i = A_i B_i$, where A_i and B_i are jointly Gaussian with covariance matrix

$$\begin{pmatrix} 1 & \langle \mathbf{v}_i, \mathbf{w}_i \rangle \\ \langle \mathbf{v}_i, \mathbf{w}_i \rangle & 1 \end{pmatrix}.$$

The proof of this generalization of Proposition 6.7 follows that of Proposition 6.7 presented in Section 7, requiring only heavier notation. In fact, our method may also be used to prove the universality of the joint eigenvalue-eigenvector distribution for any matrix Q of the form (2.1) with $\Sigma = TT^* = I_M$; see Theorem 8.3 below for a precise statement.

The proof of Proposition 6.7 is postponed to Section 7.

Supposing Proposition 6.3 holds, together with Propositions 6.4 and 6.7, we may complete the proof of Theorem 2.18.

PROOF OF THEOREM 2.18. Abbreviating $b := a - s_+$ and

$$\mathbf{u} := \frac{1}{\sqrt{M}} \sum_i \frac{\sqrt{\sigma_i} w_i}{d_i - 1} \mathbf{v}_i,$$

we define

$$\widehat{\Theta}(a, \mathbf{w}) := M \frac{\langle \mathbf{u}, \zeta_b \rangle^2}{|\mathbf{u}|^2}.$$

Then, by assumption on a , we may rewrite (6.20) as

$$\mathbf{1}(\Omega) \langle \mathbf{w}, \xi_a \rangle^2 = \mathbf{1}(\Omega) |\mathbf{u}|^2 \widehat{\Theta}(a, \mathbf{w}) + O_{\prec}(K^{-2\tau/15} |\mathbf{u}|^2).$$

Moreover, by Theorem 2.6 and Proposition 6.4, we have $\mathbb{P}(\Omega) \geq 1 - K^{-c}$ for some constant $c > 0$. Finally, by Proposition 6.7 we have $\widehat{\Theta}(a, \mathbf{w}) \rightarrow \chi_1^2$ in distribution (even in the sense of moments). The claim now follows easily. $\square$

The remainder of this section is devoted to the proof of Proposition 6.3.

PROOF OF PROPOSITION 6.3. We define the contour Γ_a as the positively oriented circle of radius $K^{-\tau/5}\Delta_a$ with centre λ_b . Let $\varepsilon > 0$ and $\tau := 1/2$, and choose a high-probability event Ξ such that (4.7), (4.8), and (4.9) hold. For the following we fix a realization $H \in \Omega \cap \Xi$. Define

$$Y_a(\mathbf{w}) := -\frac{1}{2\pi i} \oint_{\Gamma_a} \phi^{1/2} z \widetilde{G}_{\mathbf{w}\mathbf{w}}(z) dz.$$

By the residue theorem and the definition of Ω , we find

$$Y_a(\mathbf{w}) = \phi^{1/2} \mu_a \langle \mathbf{w}, \boldsymbol{\xi}_a \rangle^2. \quad (6.23)$$

To simplify notation, suppose now that $i \in R$ and consider $\mathbf{w} = \mathbf{v}_i$. From (3.35) we find that

$$Y_a(\mathbf{v}_i) = \frac{\sigma_i}{d_i^2} \frac{1}{2\pi i} \oint_{\Gamma_a} \left(\frac{1}{D^{-1} + W(z)} \right)_{ii} dz. \quad (6.24)$$

In order to compute (6.24), we need precise estimates for W on Γ_a . Because the contour Γ_a crosses the branch cut of w_ϕ , we should not compare $W(z)$ to $w_\phi(z)$ for $z \in \Gamma_a$. Instead, we compare $W(z)$ to $w_\phi(z_0)$, where

$$z_0 := \lambda_b + i\eta, \quad \eta := K^{-\tau/5}\Delta_a.$$

We claim that

$$\|W(z) - w_\phi(z_0)\| \leq CK^{-1+\varepsilon}\eta^{-1}. \quad (6.25)$$

for all $z \in \Gamma_a$. To see this, we split

$$\|W(z) - w_\phi(z_0)\| \leq \|W(z) - W(z_0)\| + \|W(z_0) - w_\phi(z_0)\|. \quad (6.26)$$

We estimate the first term of (6.26) by spectral decomposition, using that $\text{dist}(z, \sigma(H)) \geq c\eta$, similarly to (4.12). The result is

$$\begin{aligned} \|W(z) - W(z_0)\| &= C(1 + \phi) \max_i \text{Im } G_{\mathbf{v}_i \mathbf{v}_i}(z_0) \\ &\leq C(1 + \phi) \left(\text{Im } m_{\phi^{-1}}(z_0) + \frac{K^\varepsilon}{1 + \phi} \frac{1}{K\eta} \right) \\ &\leq C \text{Im } w_\phi(z_0) + CK^{-1+\varepsilon}\eta^{-1} \\ &\leq CK^{-1+\varepsilon}\eta^{-1}, \end{aligned}$$

where we used (4.7), (4.9), and Lemma 3.6. Moreover, we estimate the second term of (6.26) using (4.8) as

$$\|W(z_0) - w_\phi(z_0)\| \leq K^{-1+\varepsilon}\eta^{-1}.$$

This concludes the proof of (6.25).

Next, we claim that

$$|1 + d_i w_\phi(z_0)| \geq c|d_i - 1|. \quad (6.27)$$

The proof of (6.27) is analogous to that of (6.12), using (4.7) and the assumption on a ; we omit the details.

Armed with (6.25) and (6.27), we may analyse (6.24). A resolvent expansion in the matrix $w_\phi(z_0) - W(z)$ yields

$$\begin{aligned} &Y_a(\mathbf{v}_i) \\ &= \frac{\sigma_i}{d_i^2} \frac{1}{2\pi i} \oint_{\Gamma_a} \left(\frac{1}{d_i^{-1} + w_\phi(z_0)} + \frac{w_\phi(z_0) - W_{ii}(z)}{(d_i^{-1} + w_\phi(z_0))^2} + \left(\frac{w_\phi(z_0) - W(z)}{d_i^{-1} + w_\phi(z_0)} \frac{1}{D^{-1} + W(z)} \frac{w_\phi(z_0) - W(z)}{d_i^{-1} + w_\phi(z_0)} \right)_{ii} \right) dz. \end{aligned} \quad (6.28)$$

We estimate the third term using the bound

$$\left\| \frac{1}{D^{-1} + W(z)} \right\| \leq \frac{C}{\alpha_+}. \quad (6.29)$$

To prove (6.29), we note first that by (6.27) we have

$$\min_i |d_i^{-1} + w_\phi(z_0)| \geq \min_i \frac{|1 + d_i w_\phi(z_0)|}{|d_i|} \geq c \min_i \frac{|d_i - 1|}{|d_i|} \geq c \alpha_+.$$

By (6.25) and assumption on a , it is easy to check that

$$\min_i |d_i^{-1} + w_\phi(z_0)| \geq K^{\tau/5} \|W(z) - w_\phi(z_0)\|,$$

from which (6.29) follows.

We may now return to (6.28). The first term vanishes, the second is computed by spectral decomposition of W , and the third is estimated using (6.29). This gives

$$Y_a(\mathbf{v}_i) = \frac{\phi^{1/2} \sigma_i \lambda_b}{(1 + d_i w_\phi(z_0))^2} \langle \mathbf{v}_i, \boldsymbol{\zeta}_b \rangle^2 + O\left(\frac{\sigma_i}{|d_i - 1|^2 K} \frac{K^{2\varepsilon}}{K \eta \alpha_+}\right),$$

where we also used (6.27).

Recalling (6.23) and (4.7), we therefore get

$$\langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2 = \frac{\sigma_i \lambda_b / \mu_a}{(1 + d_i w_\phi(z_0))^2} \langle \mathbf{v}_i, \boldsymbol{\zeta}_b \rangle^2 + O\left(\frac{\sigma_i}{|d_i - 1|^2 M} \frac{K^{-1/3+2\varepsilon+\tau/5} a^{1/3}}{\alpha_+}\right),$$

where we used $\phi^{1/2} \mu_a \asymp 1 + \phi$.

In order to simplify the leading term, we use

$$\frac{-1}{1 + d_i w_\phi(z_0)} = \frac{1}{d_i - 1} + O\left(\frac{K^{-1/3+\varepsilon} a^{1/3}}{|d_i - 1| \alpha_+}\right),$$

as follows from

$$|w_\phi(z_0) + 1| \leq C \sqrt{\kappa(z_0)} + \eta \leq K^{-1/3+\varepsilon} a^{1/3},$$

where we used Lemma 3.6. Moreover, we use that

$$\lambda_b / \mu_a = 1 + O(K^{-2/3}).$$

Using that Ξ has high probability for all $\varepsilon > 0$ and recalling the isotropic delocalization bound (3.12), we therefore get for any random H that

$$\mathbf{1}(\Omega) \langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle^2 = \mathbf{1}(\Omega) \frac{\sigma_i}{(d_i - 1)^2} \langle \mathbf{v}_i, \boldsymbol{\zeta}_b \rangle^2 + O_{\prec}\left(\frac{\sigma_i}{|d_i - 1|^2 M} \frac{K^{-1/3+\tau/5} a^{1/3}}{\alpha_+}\right), \quad (6.30)$$

We proved (6.30) under the assumption that $i \in R$, but a continuity argument analogous to that given after (5.28) implies that (6.30) holds for all $i \in \llbracket 1, M \rrbracket$. The above argument may be repeated verbatim to yield

$$\mathbf{1}(\Omega) \langle \mathbf{v}_i, \boldsymbol{\xi}_a \rangle \langle \boldsymbol{\xi}_a, \mathbf{v}_j \rangle = \mathbf{1}(\Omega) \frac{\sqrt{\sigma_i \sigma_j}}{(d_i - 1)(d_j - 1)} \langle \mathbf{v}_i, \boldsymbol{\zeta}_b \rangle \langle \boldsymbol{\zeta}_b, \mathbf{v}_j \rangle + O_{\prec}\left(\frac{\sqrt{\sigma_i \sigma_j}}{|d_i - 1| |d_j - 1| M} \frac{K^{-1/3+\tau/5} a^{1/3}}{\alpha_+}\right).$$

Since we may always choose the basis $\{\mathbf{v}_i\}_{i=1}^M$ so that at most $|R| + 1$ components of $(w_1, \dots, w_M)$ are nonzero, the claim now follows easily. $\square$

7. Quantum unique ergodicity near the soft edge of H

This section is devoted to the proof of Proposition 6.7. In particular, this section only concerns the undeformed matrix $H = XX^*$ and not Q . In this section, therefore, we sometimes use symbols such as Q and R , which have nothing to do with their counterparts pertaining to the deformed matrix Q from the previous sections.

LEMMA 7.1. *Fix $\tau \in (0, 1)$. Let h be a smooth function satisfying*

$$|h'(x)| \leq C(1 + |x|)^C \quad (7.1)$$

for some positive constant C . Let $a \leq K^{1-\tau}$ and suppose that λ_a satisfies (6.21) with some constants ε and δ . Then for small enough $\delta_1 = \delta_1(\varepsilon, \delta)$ and $\delta_2 = \delta_2(\varepsilon, \delta, \delta_1)$ the following holds. Defining

$$\eta := \Delta_a K^{-2\varepsilon}, \quad E^\pm := E \pm K^{\delta_1} \eta, \quad I := [\gamma_a - K^{\delta_2} \Delta_a, \gamma_a + K^{\delta_2} \Delta_a], \quad (7.2)$$

we have

$$\mathbb{E} h(M \langle \mathbf{w}, \zeta_a \rangle^2) - \mathbb{E} h\left(\frac{M}{\pi} \int_I \operatorname{Im} G_{\mathbf{w}\mathbf{w}}(E + i\eta) \chi(E) dE\right) = O(K^{-\delta_2/2}), \quad (7.3)$$

where we defined $\chi(E) := \mathbf{1}(\lambda_{a+1} \leq E^- \leq \lambda_a)$ and introduced the convention $\lambda_0 := +\infty$.

PROOF. By the assumption (7.1) on h , rigidity (3.14), and delocalization (3.12), we can write

$$\mathbb{E} h(M \langle \mathbf{w}, \zeta_a \rangle^2) = \mathbb{E} h\left(\frac{M\eta}{\pi} \int_{I \cap [\alpha, \beta]} \frac{\langle \mathbf{w}, \zeta_a \rangle^2}{(E - \lambda_a)^2 + \eta^2} dE\right) + O(K^{-\delta_1/2}) \quad (7.4)$$

provided that

$$\alpha \leq \lambda_a^-, \quad \beta \geq \lambda_a^+, \quad (7.5)$$

where defined $\lambda_a^\pm := \lambda_a \pm N^{\delta_1} \eta$. For the following we choose

$$\alpha := \min\{\lambda_a^-, \lambda_{a+1}^+\}, \quad \beta := \lambda_a^+.$$

Now from (6.21) we get $\mathbb{P}(\lambda_{a+1}^+ \geq \lambda_a^-) \leq K^{-\delta}$ for $\delta_1 < \varepsilon$. For $\delta_1 < \delta$, we therefore get

$$\mathbb{E} h(M |\langle \mathbf{w}, \zeta_a \rangle|^2) = \mathbb{E} h\left(\frac{M\eta}{\pi} \int_I \frac{|\langle \mathbf{w}, \zeta_a \rangle|^2}{(E - \lambda_a)^2 + \eta^2} \chi(E) dE\right) + O(K^{-\delta_1/2}). \quad (7.6)$$

Next, we replace the integrand in (7.6) by $\operatorname{Im} G_{\mathbf{w}\mathbf{w}}(E + i\eta)$. By spectral decomposition, we have

$$\operatorname{Im} G_{\mathbf{w}\mathbf{w}}(E + i\eta) = \sum_{b \neq a} \frac{\eta \langle \mathbf{w}, \zeta_b \rangle^2}{(E - \lambda_b)^2 + \eta^2} + \frac{\eta \langle \mathbf{w}, \zeta_a \rangle^2}{(E - \lambda_a)^2 + \eta^2}. \quad (7.7)$$

Next, we claim that for $\delta_2 < \delta_1$ we have

$$M \int_I \operatorname{Im} G_{\mathbf{w}\mathbf{w}}(E + i\eta) dE \prec K^{\delta_2}. \quad (7.8)$$

The proof of (7.8) follows by using the spectral decomposition from (7.7) with the delocalization bound (3.12); for $|b - a| \geq K^{\delta_2}$ we use the rigidity bound (3.14), and for $|b - a| \leq K^{\delta_2}$ we estimate the integral using that $\int \frac{\eta}{e^2 + \eta^2} de = \pi$. We omit the full details.

Using (7.1), the bound (7.8), and that all terms on the right-hand side of (7.7) are nonnegative, we get from the mean-value theorem and (7.6) that the left-hand side of (7.3) is bounded by

$$K^{C\delta_2} \mathbb{E} \sum_{b \neq a} \frac{M\eta}{\pi} \int_I \frac{\langle \mathbf{w}, \zeta_b \rangle^2}{(E - \lambda_b)^2 + \eta^2} \chi(E) dE + CK^{-\delta_1/2}. \quad (7.9)$$

Using the eigenvalue rigidity from (3.14), it is not hard to see that there exists a constant C_1 such that the contribution of $|b - a| \geq K^{C_1 \delta_2}$ to (7.9) is bounded by $K^{-\delta_2}$. In order to prove (7.3), therefore, it suffices to prove

$$\mathbb{E} \sum_{b: |b-a| \leq K^{C_1 \delta_2}} \frac{M\eta}{\pi} \int_I \frac{\langle \mathbf{w}, \boldsymbol{\zeta}_b \rangle^2}{(E - \lambda_b)^2 + \eta^2} \chi(E) dE = O(K^{-\delta_2/2}) \quad (7.10)$$

For $b > a$, we get using (3.12) that

$$\sum_{b > a, |b-a| \leq K^{C_1 \delta_2}} \frac{M\eta}{\pi} \mathbb{E} \int_I \frac{\langle \mathbf{w}, \boldsymbol{\zeta}_b \rangle^2}{(E - \lambda_b)^2 + \eta^2} \chi(E) dE \leq K^{C_1 \delta_2} \mathbb{E} \int_{\lambda_{a+1}^+}^{\infty} \frac{\eta}{(E - \lambda_{a+1})^2 + \eta^2} dE \quad (7.11)$$

$$\leq CK^{C_1 \delta_2 - \delta_1/2}, \quad (7.12)$$

which is the right-hand side of (7.10) provided δ_2 is chosen small enough.

For $b < a$, we partition $I = I_1 \cup I_2$ with $I_1 \cap I_2 = \emptyset$ and

$$I_1 := \left\{ E \in I : \exists b < a, |b - a| \leq K^{C_1 \delta_2}, |E - \lambda_b| \leq \eta K^{\delta_1} \right\}. \quad (7.13)$$

As above, we find

$$\sum_{b < a, |b-a| \leq K^{C_1 \delta_2}} \frac{M\eta}{\pi} \mathbb{E} \int_{I_2} \frac{\langle \mathbf{w}, \boldsymbol{\zeta}_b \rangle^2}{(E - \lambda_b)^2 + \eta^2} \chi(E) dE \leq K^{C_2 \delta_2 - \delta_1/2}.$$

Let us therefore consider the integral over I_1 . One readily finds, for $\lambda_a \leq \lambda_{a-1} \leq \lambda_b$, that

$$\frac{1}{(E - \lambda_b)^2 + \eta^2} \mathbf{1}(E^- \leq \lambda_a) \leq \frac{K^{2\delta_1}}{(\lambda_b - \lambda_a)^2 + \eta^2} \leq \frac{K^{2\delta_1}}{(\lambda_{a+1} - \lambda_a)^2 + \eta^2}. \quad (7.14)$$

Using delocalization (3.12) we therefore find that

$$\sum_{b < a, |b-a| \leq K^{C_1 \delta_2}} \frac{M\eta}{\pi} \mathbb{E} \int_{I_1} \frac{\langle \mathbf{w}, \boldsymbol{\zeta}_b \rangle^2}{(E - \lambda_b)^2 + \eta^2} \chi(E) dE \leq K^{C_1 \delta_2 + 2\delta_1} \mathbb{E} \frac{\eta^2}{(\lambda_{a-1} - \lambda_a)^2 + \eta^2} \quad (7.15)$$

The expectation $\mathbb{E} \frac{\eta^2}{(\lambda_{a-1} - \lambda_a)^2 + \eta^2}$ in (7.15) is bounded by $\mathbb{P}(|\lambda_{a-1} - \lambda_a| \leq \Delta_a K^{-\varepsilon}) + O(K^{-\varepsilon})$. Using (6.21), we therefore obtain

$$\sum_{\substack{b < a \\ |b-a| \leq K^{C_1 \delta_2}}} \frac{M\eta}{\pi} \mathbb{E} \int_{I_1} \frac{\langle \mathbf{w}, \boldsymbol{\zeta}_b \rangle^2}{(E - \lambda_b)^2 + \eta^2} \chi(E) dE \leq K^{C_1 \delta_2 + 2\delta_1 - \delta} \quad (7.16)$$

This concludes the proof. $\square$

In the next step, stated in Lemma 7.2 below, we replace the sharp cutoff function χ in (7.3) with a smooth function of H . Note first that from Lemma 7.1 and the rigidity (3.14), we get

$$\mathbb{E} h(M \langle \mathbf{w}, \boldsymbol{\zeta}_a \rangle^2) - \mathbb{E} h \left(\frac{M}{\pi} \int_I \operatorname{Im} G_{\mathbf{w}\mathbf{w}}(E + i\eta) \mathbf{1}(\mathcal{N}(E^-, E_\delta) = a) dE \right) = O(K^{-\delta_2/2}), \quad (7.17)$$

where $E_\delta := \gamma_+ + 2N^{-2/3+\delta}$ and $\mathcal{N}(E^-, E_\delta) := |\{i : E^- < \lambda_i < E_\delta\}|$ is an eigenvalue counting function.

Next, for any $E_1, E_2 \in [\gamma_- - 1, \gamma_+ + 1]$ and $\tilde{\eta} > 0$ we define $f(\lambda) \equiv f_{E_1, E_2, \tilde{\eta}}(\lambda)$ to be the characteristic function of $[E_1, E_2]$ smoothed on scale $\tilde{\eta}$: $f = 1$ on $[E_1, E_2]$, $f = 0$ on $\mathbb{R} \setminus [E_1 - \tilde{\eta}, E_2 + \tilde{\eta}]$ and $|f'| \leq C \tilde{\eta}^{-1}$, $|f''| \leq C \tilde{\eta}^{-2}$. Moreover, let $q \equiv q_a : \mathbb{R} \rightarrow \mathbb{R}_+$ be a smooth cutoff function concentrated around a , satisfying

$$q(x) = q_a(x) = 1 \quad \text{if } |x - a| \leq 1/3, \quad q(x) = 0 \quad \text{if } |x - a| \geq 2/3, \quad |q'| \leq 6. \quad (7.18)$$

The following result is the appropriate smoothed version of (7.17). It is a simple extension of Lemma 3.2 and Equation (5.8) from [27], and its proof is omitted.

LEMMA 7.2. Let $E_\delta := \gamma_+ + 2K^{-2/3+\delta}$ and

$$\tilde{\eta} := \eta K^{-\varepsilon} = \Delta_a K^{-3\varepsilon},$$

and abbreviate $q \equiv q_a$ and $f_E \equiv f_{E^-, E_\delta, \tilde{\eta}}$. Then under the assumptions of Lemma 7.1 we have

$$\mathbb{E} h(M\langle \mathbf{w}, \zeta_a \rangle^2) - \mathbb{E} h\left(\frac{M}{\pi} \int_I \operatorname{Im} G_{\mathbf{w}\mathbf{w}}(E + i\eta) q(\operatorname{Tr} f_E(H)) dE\right) = O(K^{-\delta_2/2}). \quad (7.19)$$

We may now conclude the proof of Proposition 6.7.

PROOF OF PROPOSITION 6.7. The basic strategy of the proof is to compare the distribution of $\langle \mathbf{w}, \zeta_a \rangle^2$ under a general X to that under a Gaussian X . In the latter case, by unitary invariance of $H = XX^*$, we know that ζ_a is uniformly distributed on the unit sphere of $\mathbb{R}^M$, so that $M\langle \mathbf{w}, \zeta_a \rangle^2 \rightarrow \chi_1^2$ in distribution.

For the comparison argument, we use the Green function comparison method applied to the Helffer-Sjöstrand representation of $f(H)$. Using Lemma 7.2 it suffices to estimate

$$(\mathbb{E} - \mathbb{E}^{\text{Gauss}}) h\left(\frac{M}{\pi} \int_I \operatorname{Im} G_{\mathbf{w}\mathbf{w}}(E + i\eta) q(\operatorname{Tr} f_E(H)) dE\right) \quad (7.20)$$

where $\mathbb{E}^{\text{Gauss}}$ denotes the expectation with respect to Gaussian X . Now we express $f(H)$ in terms of Green functions using Helffer-Sjöstrand functional calculus. Recall the definition of κ_a from (2.23). Let $g(y)$ be a smooth cutoff function with support in $[-\kappa_a, \kappa_a]$, with $g(y) = 1$ for $|y| \leq \kappa_a/2$ and $\|g^{(n)}\|_\infty \leq C\kappa_a^{-n}$, where $g^{(n)}$ denotes the n -th derivative of g . Then, similarly to (3.28), we have (see e.g. Equation (B.12) of [19])

$$f_E(\lambda) = \frac{1}{2\pi} \int_{\mathbb{R}^2} \frac{i\sigma f_E''(e)g(\sigma) + if_E(e)g'(\sigma) - \sigma f_E'(e)g'(\sigma)}{\lambda - e - i\sigma} de d\sigma. \quad (7.21)$$

Thus we get the functional calculus, with $G(z) = (H - z)^{-1}$,

$$\begin{aligned} \operatorname{Tr} f_E(H) &= \frac{1}{2\pi} \int_{\mathbb{R}^2} \left(i\sigma f_E''(e)g(\sigma) + if_E(e)g'(\sigma) - \sigma f_E'(e)g'(\sigma) \right) \operatorname{Tr} G(e + i\sigma) de d\sigma \\ &= \frac{1}{2\pi} \int_{\mathbb{R}^2} \left(if_E(e)g'(\sigma) - \sigma f_E'(e)g'(\sigma) \right) \operatorname{Tr} G(e + i\sigma) de d\sigma \\ &\quad + \frac{i}{2\pi} \int_{|\sigma| > \tilde{\eta}K^{-d\varepsilon}} d\sigma g(\sigma) \int de f_E''(e) \sigma \operatorname{Tr} G(e + i\sigma) \\ &\quad + \frac{i}{2\pi} \int_{-\tilde{\eta}K^{-d\varepsilon}}^{\tilde{\eta}K^{-d\varepsilon}} d\sigma \int de f_E''(e) \sigma \operatorname{Tr} G(e + i\sigma). \end{aligned} \quad (7.22)$$

As in Lemma 5.1 of [27], one can easily extend (3.10) to η satisfying the lower bound $\eta > 0$ instead of $\eta \geq K^{-1+\omega}$ in (3.8); the proof is identical to that of [27, Lemma 5.1]. Thus we have, for $e \in [\gamma_+ - 1, \gamma_+ + 1]$ and $\sigma \in (0, 1)$,

$$\sigma \operatorname{Tr} G(e + i\sigma) = O_{\prec}(1). \quad (7.23)$$

Therefore, by the trivial symmetry $\sigma \mapsto -\sigma$ combined with complex conjugation, the third term on the right-hand side of (7.22) is bounded by

$$\frac{i}{2\pi} \int_{-\tilde{\eta}K^{-d\varepsilon}}^{\tilde{\eta}K^{-d\varepsilon}} d\sigma \int de f_E''(e) \sigma \operatorname{Tr} G(e + i\sigma) = O_{\prec}(K^{-d\varepsilon}), \quad (7.24)$$

where we used that $\int |f_E''(e)| de = O(\tilde{\eta}^{-1})$. Next, we note that (3.10) and Lemma 3.6 imply

$$\frac{M}{\pi} \int_I \operatorname{Im} G_{\mathbf{w}\mathbf{w}}(E + i\eta) dE = O_{\prec}(K^{3\varepsilon}). \quad (7.25)$$

Recalling (7.1) and using the mean value theorem, we find from (7.19), (7.24), and (7.25) that for large enough d , in order to estimate (7.20), and hence prove (6.22), it suffices to prove the following lemma. Note that in it we choose $X^{(1)}$ to be the original ensemble and $X^{(2)}$ to be the Gaussian ensemble. $\square$

LEMMA 7.3. Suppose that the two $M \times N$ matrix ensembles $X^{(1)}$ and $X^{(2)}$ satisfy (2.4) and (2.5). Suppose that the assumptions of Lemma 7.1 hold, and recall the notations

$$\eta := \Delta_a K^{-2\varepsilon}, \quad \tilde{\eta} := \Delta_a K^{-3\varepsilon}, \quad E^\pm := E \pm K^{\delta_1} \eta, \quad E_\delta := \gamma_+ + 2K^{-2/3+\delta}, \quad (7.26)$$

as well as

$$f_E \equiv f_{E^-, E_\delta, \tilde{\eta}}, \quad I := [\gamma_a - K^{\delta_2} \Delta_a, \gamma_a + K^{\delta_2} \Delta_a], \quad (7.27)$$

where $f_{E^-, E_\delta, \tilde{\eta}}$ was defined above (7.18). Recall $q \equiv q_a$ from (7.18). Finally, suppose that $\varepsilon > 4\delta_1$ and $\delta_1 > 4\delta_2$.

Then for any $d > 1$ and for small enough $\varepsilon \equiv \varepsilon(\tau, d) > 0$ and $\delta_2 \equiv \delta_2(\varepsilon, \delta, \delta_1)$ we have

$$[\mathbb{E}^{(1)} - \mathbb{E}^{(2)}] h \left[\int_I x(E) q(y(E)) dE \right] = O(K^{-\delta_2}), \quad (7.28)$$

where we defined

$$x(E) := \frac{M}{\pi} \operatorname{Im} G_{\mathbf{w}\mathbf{w}}(E + i\eta) \quad (7.29)$$

and

$$\begin{aligned} y(E) &:= \frac{1}{2\pi} \int_{\mathbb{R}^2} i\sigma f_E''(e) g(\sigma) \operatorname{Tr} G(e + i\sigma) \mathbf{1}(|\sigma| > \tilde{\eta} K^{-d\varepsilon}) de d\sigma \\ &+ \frac{1}{2\pi} \int_{\mathbb{R}^2} (if_E(e) g'(\sigma) - \sigma f_E'(e) g'(\sigma)) \operatorname{Tr} G(e + i\sigma) de d\sigma. \end{aligned} \quad (7.30)$$

The rest of this section is devoted to the proof of Lemma 7.3.

7.1. Proof of Lemma 7.3 I: preparations. We shall use the Green function comparison method [20, 21, 27] to prove Lemma 7.3. For definiteness, we assume throughout the remainder of Section 7 that $\phi \geq 1$. The case $\phi < 1$ is dealt with similarly, and we omit the details.

We first collect some basic identities and estimates that serve as a starting point for the Green function comparison argument. We work on the product probability space of the ensembles $X^{(1)}$ and $X^{(2)}$. We fix a bijective ordering map Φ on the index set of the matrix entries,

$$\Phi : \{(i, \mu) : 1 \leq i \leq M, 1 \leq \mu \leq N\} \longrightarrow \llbracket 1, MN \rrbracket,$$

and define the interpolating matrix X_γ , $\gamma \in \llbracket 1, MN \rrbracket$, through

$$(X_\gamma)_{i\mu} := \begin{cases} X_{i\mu}^{(1)} & \text{if } \Phi(i, \mu) > \gamma \\ X_{i\mu}^{(2)} & \text{if } \Phi(i, \mu) \leq \gamma. \end{cases}$$

In particular, $X_0 = X^{(1)}$ and $X_{MN} = X^{(2)}$. Hence we have the telescopic sum

$$[\mathbb{E}^{(1)} - \mathbb{E}^{(2)}] h \left[\int_I x(E) q(y(E)) dE \right] = \sum_{\gamma=1}^{MN} [\mathbb{E}^{X_{\gamma-1}} - \mathbb{E}^{X_\gamma}] h \left[\int_I x(E) q(y(E)) dE \right] \quad (7.31)$$

(in self-explanatory notation).

Let us now fix a γ and let (b, β) be determined by $\Phi(b, \beta) = \gamma$. Throughout the following we consider b, β to be arbitrary but fixed and often omit dependence on them from the notation. Our strategy is to compare $X_{\gamma-1}$ with X_γ for each γ . In the end we shall sum up the differences in the telescopic sum (7.31).

Note that $X_{\gamma-1}$ and X_γ differ only in the matrix entry indexed by (b, β) . Thus we may write

$$\begin{aligned} X_{\gamma-1} &= Q + \tilde{U}, & \tilde{U}_{i\mu} &:= \delta_{ib} \delta_{\mu\beta} X_{b\beta}^{(1)}, \\ X_\gamma &= Q + U, & U_{i\mu} &:= \delta_{ib} \delta_{\mu\beta} X_{b\beta}^{(2)}. \end{aligned} \quad (7.32)$$

Here Q is the matrix obtained from X_γ (or, equivalently, from $X_{\gamma-1}$) by setting the entry indexed by (b, β) to zero. Next, we define the resolvents

$$T(z) := (QQ^* - z)^{-1}, \quad S(z) := (X_\gamma X_\gamma^* - z)^{-1}. \quad (7.33)$$

We shall show that the difference between the expectations $\mathbb{E}^{X_\gamma}$ and $\mathbb{E}^Q$ depends only on the first two moments of $X_{b\beta}^{(2)}$, up to an error term that is negligible even after summation over γ . Together with same argument applied to $\mathbb{E}^{X_{\gamma-1}}$, and the fact that the second moments of $X_{b\beta}^{(1)}$ and $X_{b\beta}^{(2)}$ coincide, this will prove Lemma 7.3.

We define $x_T(E)$ and $y_T(E)$ as in (7.29) and (7.30) with G replaced by T , and similarly $x_S(E)$ and $y_S(E)$ with G replaced by S . Throughout the following we use the notation $\mathbf{w} = (w(i))_{i=1}^M$ for the components of $\mathbf{w}$. In order to prove (7.28) using (7.31), it is enough to prove that for some constant $c > 0$ we have

$$\mathbb{E}h \left[\int_I x_S(E) q(y_S(E)) dE \right] - \mathbb{E}h \left[\int_I x_T(E) q(y_T(E)) dE \right] = \mathbb{E}\mathcal{A} + O(K^{-c})(\phi^{-1}K^{-2} + K^{-1}|w(b)|^2), \quad (7.34)$$

where $\mathcal{A}$ is polynomial of degree two in $U_{b\beta}$ whose coefficients are Q -measurable.

The rest of this section is therefore devoted to the proof of (7.34). Recall that we assume throughout that $\phi \geq 1$ for definiteness; in particular, $K = N$.

We begin by collecting some basic identities from linear algebra. In addition to $G(z) := (XX^* - z)^{-1}$ we introduce the auxiliary resolvent $R(z) := (X^*X - z)^{-1}$. Moreover, for $\mu \in \llbracket 1, M \rrbracket$ we split

$$X = X_{[\mu]} + X^{[\mu]}, \quad (X_{[\mu]})_{i\nu} := \mathbf{1}(\nu = \mu)X_{i\nu}, \quad (X^{[\mu]})_{i\nu} := \mathbf{1}(\nu \neq \mu)X_{i\nu}.$$

We also define the resolvent $G^{[\mu]} := (X^{[\mu]}(X^{[\mu]})^* - z)^{-1}$. A simple Neumann series yields the identity

$$G = G^{[\mu]} - \frac{G^{[\mu]}X_{[\mu]}X_{[\mu]}^*G^{[\mu]}}{1 + (X^*G^{[\mu]}X)_{\mu\mu}}. \quad (7.35)$$

Moreover, from [9, Equation (3.11)], we find

$$zR_{\mu\mu} = -\frac{1}{1 + (X^*G^{[\mu]}X)_{\mu\mu}}. \quad (7.36)$$

From (7.35) and (7.36) we easily get

$$GX_{[\mu]} = -zR_{\mu\mu}G^{[\mu]}X_{[\mu]}, \quad X_\mu^*G = -zR_{\mu\mu}X_{[\mu]}^*G^{[\mu]}. \quad (7.37)$$

Throughout the following we shall make use of the fundamental error parameter

$$\Psi(z) := \sqrt{\frac{\operatorname{Im} m_\phi(z)}{N\eta}} + \frac{1}{N\eta}, \quad (7.38)$$

which is analogous to the right-hand side of (3.10) and will play a similar role. We record the following estimate, which is analogous to Theorem 3.2.

LEMMA 7.4. *Under the assumptions of Theorem 3.2 we have, for $z \in \mathbf{S}$,*

$$|(GX)_{\mathbf{w}\mu}| \prec \phi^{-1/4}\Psi \quad (7.39)$$

and

$$|(X^*GX)_{\mu\nu} - \delta_{\mu\nu}(1 + zm_\phi)| \prec \phi^{1/2}\Psi. \quad (7.40)$$

PROOF. This result is a generalization of (5.22) in [33]. The key identity is (7.37). Since $G^{[\mu]}$ is independent of $(X_{i\mu})_{i=1}^N$, we may apply the large deviation estimate [9, Lemma 3.1] to $G^{[\mu]}X_{[\mu]}$. Moreover, $|R_{\mu\mu}| \prec 1$, as follows from Theorem 3.2 applied to X^* , and Lemma 3.6. Thus we get

$$\begin{aligned} |(GX)_{\mathbf{w}\mu}| &\prec \phi^{1/2} \left((MN)^{-1/2} \sum_{i=1}^M |G_{\mathbf{w}i}^{[\mu]}|^2 \right)^{1/2} = \phi^{1/4} \left(\frac{1}{N\eta} \operatorname{Im} G_{\mathbf{w}\mathbf{w}}^{[\mu]} \right)^{1/2} \\ &\prec \phi^{1/4} \left(\frac{1}{N\eta} \left(\operatorname{Im} m_{\phi^{-1}} + \phi^{-1}\Psi \right) \right)^{1/2} \leq C\phi^{-1/4}\Psi, \end{aligned}$$

where the second step follows by spectral decomposition, the third step from Theorem 3.2 applied to $X^{[\mu]}$ as well as (3.21), and the last step by definition of Ψ . This concludes the proof of (7.39).

Finally, (7.40) follows easily from Theorem 3.2 applied to the identity $X^*GX = 1 + zR$. $\square$

After these preparations, we continue the proof of (7.34). We first expand the difference between S and T in terms of V (see (7.32)). We use the resolvent expansion: for any $m \in \mathbb{N}$ we have

$$S = T + \sum_{k=1}^m (-1)^k [T(QU^* + UQ^* + UU^*)]^k T + (-1)^{m+1} [T(QU^* + UQ^* + UU^*)]^{m+1} S \quad (7.41)$$

and

$$T = S + \sum_{k=1}^m [S(XU^* + UX^* + UU^*)]^k S + [S(XU^* + UX^* + UU^*)]^{m+1} T. \quad (7.42)$$

Note that Theorem 3.2 and Lemma 7.4 immediately yield for $z \in \mathbf{S}$

$$|S_{\mathbf{v}\mathbf{w}} - \langle \mathbf{v}, \mathbf{w} \rangle m_{\phi^{-1}}| \prec \phi^{-1}\Psi, \quad |(SX_\gamma)_{\mathbf{v}i}| \prec \phi^{-1/4}\Psi, \quad |(X_\gamma^* SX_\gamma)_{\mu\nu} - \delta_{\mu\nu}(1 + zm_\phi)| \prec \phi^{1/2}\Psi$$

Using (7.42), we may extend these estimates to analogous ones on Q and T instead of X_γ and S . Indeed, using the facts $\|R\| \leq \eta^{-1}$, $\Psi \geq N^{-1/2}$, and $|U_{b\beta}| \prec \phi^{-1/4}N^{-1/2}$ (which are easily derived from the definitions of the objects on the left-hand sides) combined with (7.42), we get the following result.

LEMMA 7.5. *For $A \in \{S, T\}$ and $B \in \{X_\gamma, Q\}$ we have*

$$|A_{\mathbf{v}\mathbf{w}} - \langle \mathbf{v}, \mathbf{w} \rangle m_{\phi^{-1}}| \prec \phi^{-1}\Psi, \quad |(AB)_{\mathbf{v}i}| \prec \phi^{-1/4}\Psi, \quad |(B^*AB)_{\mu\nu} - \delta_{\mu\nu}(1 + zm_\phi(z))| \prec \phi^{1/2}\Psi. \quad (7.43)$$

The final tool that we shall need is the following lemma, which collects basic algebraic properties of stochastic domination $\prec$. We shall use them tacitly throughout the following. Their proof is an elementary exercise using union bounds and Cauchy-Schwarz. See [9, Lemma 3.2] for a more general statement.

LEMMA 7.6. (i) *Suppose that $A(v) \prec B(v)$ uniformly in $v \in V$. If $|V| \leq N^C$ for some constant C then $\sum_{v \in V} A(v) \prec \sum_{v \in V} B(v)$.*

(ii) *Suppose that $A_1 \prec B_1$ and $A_2 \prec B_2$. Then $A_1 A_2 \prec B_1 B_2$.*

(iii) *Suppose that $\Psi \geq N^{-C}$ is deterministic and A is a nonnegative random variable satisfying $\mathbb{E}A^2 \leq N^C$. Then $A \prec \Psi$ implies that $\mathbb{E}A \prec \Psi$.*

If the above random variables depend on an additional parameter u and all hypotheses are uniform in u then so are the conclusions.

7.2. Proof of Lemma 7.3 II: the main expansion. Lemma 7.5 contains the a-priori estimates needed to control the resolvent expansion (7.41). The precise form that we shall need is contained in the following lemma, which is our main expansion. Define the control parameter

$$\Psi_b(z) := \phi^{1/2}|w(b)| + \Psi.$$

LEMMA 7.7 (RESOLVENT EXPANSION OF $x(E)$ AND $y(E)$). *The following results hold for $E \in I$. (Recall the definition (7.27). For brevity, we omit E from our notation.)*

(i) *We have the expansion*

$$x_S - x_T = \sum_{l=1}^3 x_l U_{b\beta}^l + O_{\prec}(\phi^{-1}N^{-1}(\phi^{1/4}|w(b)| + \Psi)^2), \quad (7.44)$$

where x_l is a polynomial, with constant number of terms, in the variables

$$\{T_{bb}, T_{\mathbf{w}b}, T_{b\mathbf{w}}, (TQ)_{\mathbf{w}\beta}, (Q^*T)_{\beta\mathbf{w}}, (Q^*TQ)_{\beta\beta}\}.$$

In each term of x_l , the index $\mathbf{w}$ appears exactly twice, while the indices b and β each appear exactly l times.

Moreover, we have the estimates

$$|x_1| + |x_3| \prec \phi^{-1/4}N\Psi\Psi_b, \quad |x_2| \prec \phi^{-1/2}N\Psi_b^2 + N\Psi^2, \quad (7.45)$$

where the spectral parameter on the right-hand side is $z = E + i\eta$.

(ii) *We have the expansion*

$$\mathrm{Tr} S - \mathrm{Tr} T = \sum_{l=1}^3 J_l U_{b\beta}^l + O_{\prec}(\phi^{-1}N^{-1}\Psi^2), \quad (7.46)$$

where J_l is a polynomial, with constant number of terms, in the variables

$$\{T_{bb}, (T^2)_{bb}, (T^2Q)_{b\beta}, (Q^*T^2)_{\beta b}, (Q^*TQ)_{\beta\beta}, (Q^*T^2Q)_{\beta\beta}\}.$$

In each term of J_l , T^2 appears exactly once, while the indices b and β each appear exactly l times.

Moreover, for $z \in \mathbf{S}$ we have the estimates

$$|J_1| + |J_3| \prec \phi^{-1/4}N\Psi^2, \quad |J_2| \prec N\Psi^2. \quad (7.47)$$

(iii) *Defining*

$$y_l := \frac{1}{2\pi} \int_{\mathbb{R}^2} J_l \left(i\sigma f_E''(e)g(\sigma) \mathbf{1}(|\sigma| > \tilde{\eta}N^{-d\varepsilon}) + i f_E(e)g'(\sigma) - \sigma f_E'(e)g'(\sigma) \right) \mathrm{d}e \mathrm{d}\sigma,$$

we have the expansion

$$y_S - y_T = \sum_{l=1}^3 y_l U_{a\beta}^l + O_{\prec}(N^{C\varepsilon}\phi^{-1}N^{-2}\kappa_a^{1/2}) \quad (7.48)$$

together with the bounds

$$|y_1| + |y_3| \prec \phi^{-1/4}N^{C\varepsilon}\kappa_a^{1/2}, \quad |y_2| \prec N^{C\varepsilon}\kappa_a^{1/2}. \quad (7.49)$$

Here all constants C depend on the fixed parameter d .

PROOF. The proof is an application of the resolvent expansion (7.41) with $m = 3$ to the definitions of x and y .

We begin with part (i). The expansion (7.44) is obtained by expanding the resolvent $S_{\mathbf{w}\mathbf{w}}$ in the definition of x_S using (7.41) with $m = 3$. The terms are regrouped according the power, l , of $U_{b\beta}$. The error term of (7.44) contains all terms with $l \geq 4$. It is a simple matter to check that the polynomials x_l , for $l = 1, 2, 3$, have the claimed algebraic properties. In order to establish the claimed bounds on the terms of the expansion, we use Lemma 7.5 to derive the estimates

$$|T_{\mathbf{w}b}| \prec \phi^{-1}\Psi_b, \quad |(TQ)_{\mathbf{w}\beta}| \prec \phi^{-1/4}\Psi, \quad |T_{bb}| \prec C\phi^{-1/2}, \quad |(Q^*TQ)_{\beta\beta}| \prec \phi^{1/2}, \quad (7.50)$$

and the same estimates hold if T is replaced by S . Note that in (7.50) we used the bound $|m_{\phi^{-1}}| \asymp \phi^{-1/2}$, which follows from the identity

$$m_{\phi^{-1}}(z) = \frac{1}{\phi} \left(m_{\phi}(z) + \frac{1-\phi}{z} \right)$$

and Lemma 3.6. Using (7.50), it is not hard to conclude the proof of part (i).

Part (ii) is proved in the same way as part (i), simply by setting $\mathbf{w} = \mathbf{e}_i$ and summing over $i = 1, \dots, M$.

What remains is to prove the bounds in part (iii). To that end, we integrate by parts, first in e and then in σ , in the term containing $f_E''(e)$, and obtain

$$\begin{aligned} \int_{\mathbb{R}^2} i\sigma f_E''(e)g(\sigma)J_l(e+i\sigma)\mathbf{1}(|\sigma| > \tilde{\eta}_d) \, d\sigma &= \sum_{\pm} \mp \int \tilde{\eta}_d f_E'(e)g(\pm\tilde{\eta}_d)J_l(e \pm i\tilde{\eta}_d) \, de \\ &+ \int_{\mathbb{R}^2} (\sigma g'(\sigma) + g(\sigma))f_E'(e)J_l(e+i\sigma)\mathbf{1}(|\sigma| > \tilde{\eta}_d) \, d\sigma, \end{aligned}$$

where we abbreviated $\tilde{\eta}_d := \tilde{\eta}N^{-d\varepsilon}$. Thus we get the bound

$$\begin{aligned} |y_l(E)| &\leq \int de \tilde{\eta}_d |f_E'(e)| |J_l(e+i\tilde{\eta}_d)| + \int d\sigma \, d\sigma' (|f_E(e)g'(\sigma)| + |\sigma f_E'(e)g'(\sigma)|) |J_l(e+i\sigma)| \\ &+ \int de \int_{\tilde{\eta}_d}^{\infty} d\sigma (\sigma |g'(\sigma)| + g(\sigma)) |f_E'(e)| |J_l(e+i\sigma)|. \end{aligned} \quad (7.51)$$

Using (7.51), the conclusion of the proof of part (iii) follows by a careful estimate of each term on the right-hand side, using part (ii) as input. The ingredients are the definitions of $\tilde{\eta}_d$ and κ_a , as well as the estimate

$$\Psi^2(z) \leq C \frac{\sqrt{\kappa} + \sqrt{\eta}}{N\eta} + \frac{C}{N^2\eta^2}.$$

The same argument yields the error bound in (7.48). This concludes the proof. $\square$

Armed with the expansion from Lemma 7.7, we may do a Taylor expansion of q . To that end, we record the estimate $\int_I |x_T(E)| \, dE \prec N^{C\varepsilon}$, as follows from Lemma 7.5. Hence using Lemma 7.7 and expanding $q(y_S(E))$ around $q(y_T(E))$ with a fourth order rest term, we get

$$\begin{aligned} \int_I x_S(E) q(y_S(E)) \, dE - \int_I x_T(E) q(y_T(E)) \, dE \\ = \sum_{\mathbf{l} \in \mathcal{L}} A_{\mathbf{l}} U_{b\beta}^{|\mathbf{l}|} + O_{\prec} \left(\phi^{-1} N^{-2+C\varepsilon} \kappa_a^{1/2} + \phi^{-1/2} N^{-1+C\varepsilon} \Delta_a |w(b)|^2 \right), \end{aligned} \quad (7.52)$$

where we defined

$$\mathcal{L} := \{ \mathbf{l} = (l_0, \dots, l_m) \in \llbracket 0, 3 \rrbracket \times \llbracket 1, 3 \rrbracket^m : m \in \mathbb{N}, 1 \leq |\mathbf{l}| \leq 3 \}, \quad |\mathbf{l}| := \sum_{i=0}^m l_i,$$

as well as the polynomial

$$A_{\mathbf{l}} := \int_I \frac{q^{(m)}(y_T)}{m!} x_{l_0} y_{l_1} \cdots y_{l_m} dE, \quad (7.53)$$

where we abbreviated $m \equiv m(\mathbf{l})$. Here we use the convention that $x_0 := x_T$. Note that $\mathcal{L}$ is a finite set (it has 14 elements), and for each $\mathbf{l} \in \mathcal{L}$ the polynomial $A_{\mathbf{l}}$ is independent of $U_{b\beta}$. In the estimate of the error term on the right-hand side of (7.52) we also used the fact that for $E \in I$ we have $\Psi(E + i\eta) \leq N^{C\varepsilon} \kappa_a^{1/2}$

Next, using lemma 7.7 and $N\Delta_a \asymp \kappa_a^{-1/2}$, we find

$$|A_{\mathbf{l}}| \prec N^{C\varepsilon} \begin{cases} \phi^{-1/4}(\kappa_a^{1/2} + \phi^{1/2}|w(b)|) & \text{if } |\mathbf{l}| = 1 \\ \kappa_a^{1/2} + \kappa_a^{-1/2} \phi^{1/2}|w(b)|^2 & \text{if } |\mathbf{l}| = 2 \\ \phi^{-1/4}(\kappa_a^{1/2} + \phi^{1/2}|w(b)| + \phi|w(b)|^2) & \text{if } |\mathbf{l}| = 3. \end{cases} \quad (7.54)$$

Using (7.52) and (7.54), we may do a Taylor expansion of h on the left-hand side of (7.34). This yields

$$\begin{aligned} h \left[\int_I x_S q(y_S) dE \right] - h \left[\int_I x_T q(y_T) dE \right] &= \sum_{k=1}^3 \frac{1}{k!} h^{(k)}(A_0) \left(\sum_{\mathbf{l} \in \mathcal{L}} A_{\mathbf{l}} U_{b\beta}^{|\mathbf{l}|} \right)^k \\ &+ O_{\prec} \left(\phi^{-1} N^{-2+C\varepsilon} \kappa_a^{1/2} + N^{-2+C\varepsilon} \kappa_a^{-1/2} |w(b)|^2 \right), \end{aligned} \quad (7.55)$$

where we abbreviated $A_0 := \int_I x_0 dE$. Since $a \leq N^{1-\tau}$, it is easy to see that, by choosing ε small enough depending on τ , the error term in (7.55) is bounded by $N^{-c}(\phi^{-1} N^{-2} + N^{-1}|w(b)|^2)$ for some positive constant c . Taking the expectation and recalling that $|U_{b\beta}| \prec \phi^{-1/4} N^{-1/2}$, we therefore get

$$\begin{aligned} \mathbb{E} h \left[\int_I x_S q(y_S) dE \right] - \mathbb{E} h \left[\int_I x_T q(y_T) dE \right] &= \mathbb{E} \mathcal{A} + O_{\prec} \left(N^{-c}(\phi^{-1} N^{-1} + N^{-1}|w(b)|^2) \right) \\ &+ \mathbb{E} U_{b\beta}^3 \mathbb{E} \sum_{k=1}^3 \frac{1}{k!} h^{(k)}(A_0) \sum_{\mathbf{l}_1, \dots, \mathbf{l}_k \in \mathcal{L}} \mathbf{1} \left(\sum_{i=1}^k |\mathbf{l}_i| = 3 \right) \prod_{i=1}^k A_{\mathbf{l}_i}, \end{aligned} \quad (7.56)$$

where $\mathbb{E} \mathcal{A}$ is as described after (7.34), i.e. it depends on the random variable $U_{b\beta}$ only through its first two moments.

At this point we note that if we make the stronger assumption that the first *three* moments of $X^{(1)}$ and $X^{(2)}$ match (which, in the ultimate application to the proof of Proposition 6.7, means that $\mathbb{E} X_{i\mu}^3 = 0$), the proof is now complete. Indeed, in that case we may allow $\mathcal{A}$ to be a polynomial of degree three in $U_{b\beta}$ with Q -measurable coefficients, and we may absorb the last line of (7.56) into $\mathbb{E} \mathcal{A}$. This completes the proof of (7.34), and hence of Lemma 7.3, for the special case that the third moments of $X^{(1)}$ and $X^{(2)}$ match.

For the general case, we still have to estimate the last line of (7.56). The terms that we need to analyse are

$$h^{(3)}(A_0) A_{(1)}^m A_{(0,1)}^n \quad (m+n=3) \quad (7.57a)$$

$$h^{(2)}(A_0) (A_{(2)} + A_{(0,2)} + A_{(1,1)} + A_{(0,1,1)}) (A_{(1)} + A_{(0,1)}) \quad (7.57b)$$

$$h^{(1)}(A_0) (A_{(3)} + A_{(0,3)} + A_{(1,2)} + A_{(2,1)} + A_{(0,1,2)} + A_{(1,1,1)} + A_{(0,1,1,1)}). \quad (7.57c)$$

These terms are dealt with in the following lemma.

LEMMA 7.8. *Let Y denote any term of (7.57). Then there is a constant $c > 0$ such that*

$$|\mathbb{E} Y| \leq N^{-c} (\phi^{-1/4} N^{-1/2} + \phi^{3/4} |w(b)|^2). \quad (7.58)$$

Plugging the estimate of Lemma 7.8 into (7.56), and recalling that $\mathbb{E} U_{b\beta}^3 \leq C \phi^{-3/4} N^{-3/2}$, it is easy to complete the proof of (7.34), and hence of Lemma 7.3. Lemma 7.8 is proved in the next subsection.

7.3. Proof of Lemma 7.3 III: the terms of order three and proof of Lemma 7.8. Recall that we assume $\phi \geq 1$, i.e. $K = N$; the case $\phi \leq 1$ is dealt with analogously, and we omit the details.

We first remark that using the bounds (7.54) we find

$$|\mathbb{E}Y| \leq N^{C\varepsilon} (\kappa_a^{-1/2} \phi^{3/4} |w(b)|^2 + \phi^{1/4} |w(b)| + \phi^{-1/4} \kappa_a^{1/2}). \quad (7.59)$$

Comparing this to (7.58), we see that we need to gain an additional factor $N^{-1/2}$. How to do so is the content of this subsection.

The basic idea behind the additional factor $N^{-1/2}$ is that the expectation $\mathbb{E}Y$ is smaller than the typical size $\sqrt{\mathbb{E}|Y|^2}$ of Y by a factor $N^{-1/2}$. This is a rather general property of random variables which can be written, up to a negligible error term, as a polynomial of odd degree in the entries $\{Q_{i\beta}\}_{i=1}^m$. A systematic representation of a large family of random variables in terms of polynomials was first given in [16], and was combined with a parity argument in [9]. Subsequently, an analogous parity argument for more singular functions was developed in [41]. Following [41], we refer to the process of representing a random variable Y as a polynomial in $\{Q_{i\beta}\}_{i=1}^M$ up to a negligible error term as the *polynomialization* of Y .

We shall develop a new approach to the polynomialization of the variables (7.57). The main reason is that these variables have a complicated algebraic structure, which needs to be combined with the Helffer-Sjöstrand representation (7.30). These difficulties lead us to define a family of graded polynomials (given in Definitions 7.10–7.12), which is general enough to cover the polynomialization of all terms from (7.57) and imposes conditions on the coefficients that ensure the gain of $N^{-1/2}$. The basic structure behind these polynomials is a classification based on the ℓ^2 - and ℓ^3 -norms of their coefficients.

Let us outline the rough idea of the parity argument. We use the notations $Q = Q_{[\beta]} + Q^{[\beta]}$ and $T^{[\beta]}(z) := (Q^{[\beta]}(Q^{[\beta]})^* - z)^{-1}$, in analogy to those introduced before (7.35). A simple example of a polynomial is

$$\mathcal{P}_2 = (Q^* T^{[\beta]} Q)_{\beta\beta} = \sum_{i,j} T_{ij}^{[\beta]} Q_{i\beta} Q_{j\beta}.$$

This is a polynomial of degree two. Note that the coefficients $T_{ij}^{[\beta]}$ are $Q^{[\beta]}$ -measurable, i.e. independent of $Q_{[\beta]}$. It is not hard to see that $\mathbb{E}\mathcal{P}_2$ is of the same order as $\sqrt{\mathbb{E}|\mathcal{P}_2|^2}$, so that taking the expectation of $\mathcal{P}_2$ does not yield better bounds. The situation changes drastically if the polynomial is *odd* degree. Consider for instance the polynomial

$$\mathcal{P}_3 := (Q^* T^{[\beta]} Q)_{\beta\beta} (T^{[\beta]} Q)_{\mathbf{w}\beta} = \sum_{i,j,k} T_{ij}^{[\beta]} T_{\mathbf{w}k}^{[\beta]} Q_{i\beta} Q_{j\beta} Q_{k\beta}.$$

Now we have $|\mathbb{E}\mathcal{P}_3| \lesssim N^{-1/2} \sqrt{\mathbb{E}|\mathcal{P}_3|^2}$. The reason for this gain of a factor $N^{-1/2}$ is clear: taking the expectation forces all three summation indices i, j, k to coincide.

In the following we define a large family of $\mathbb{Z}_2$ -graded polynomials that is sufficiently general to cover the polynomializations of the terms (7.57). We shall introduce a notation $O_{\prec,*}(A)$, which generalizes the notation $O_{\prec}(A)$ from (2.1); here $*$ $\in$ {even, odd} denotes the parity of the polynomial, and A its size. We always have the trivial bound $O_{\prec,*}(A) = O_{\prec}(A)$. In addition, we roughly have the estimates

$$\mathbb{E}O_{\prec,\text{even}}(A) \lesssim A, \quad \mathbb{E}O_{\prec,\text{odd}}(A) \lesssim N^{-1/2} A.$$

The need to gain an additional factor $N^{-1/2}$ from odd polynomials imposes nontrivial constraints on the polynomial coefficients, which are carefully stated in Definitions 7.10–7.12; they have been tailored to the class of polynomials generated by the terms (7.57).

We now move on to the proof of Lemma 7.8. We recall that we assume throughout that $\phi \geq 1$. We first introduce a family of graded polynomials suitable for our purposes. It depends on a constant C_0 , which we shall fix during the proof to be some large but fixed number.

DEFINITION 7.9 (ADMISSIBLE WEIGHTS). Let $\varrho = (\varrho_i : i \in \llbracket 1, \phi N \rrbracket)$ be a family of deterministic nonnegative weights. We say that ϱ is an admissible weight if

$$\frac{1}{N^{1/2} \phi^{1/4}} \left(\sum_i \varrho_i^2 \right)^{1/2} \leq 1, \quad \frac{1}{N^{1/2} \phi^{1/4}} \left(\sum_i \varrho_i^3 \right)^{1/3} \leq N^{-1/6}. \quad (7.60)$$

DEFINITION 7.10 ($O_{\prec,d}(\cdot)$). For a given degree $d \in \mathbb{N}$ let

$$\mathcal{P} = \sum_{i_1, \dots, i_d=1}^{\phi N} V_{i_1 \dots i_d} Q_{i_1 \beta} \cdots Q_{i_d \beta} \quad (7.61)$$

be a polynomial in Q . Analogously to the notation $O_{\prec}(\cdot)$ introduced in Definition 2.1, we write $\mathcal{P} = O_{\prec,d}(A)$ if the following conditions are satisfied.

- (i) A is deterministic and $V_{i_1 \dots i_d}$ is $Q^{[\beta]}$ -measurable.
- (ii) There exist admissible weights $\varrho^{(1)}, \dots, \varrho^{(d)}$ such that

$$|V_{i_1 \dots i_d}| \prec A \varrho_{i_1}^{(1)} \cdots \varrho_{i_d}^{(d)}. \quad (7.62)$$

- (iii) We have the deterministic bound $|V_{i_1 \dots i_d}| \leq N^{C_0}$.

DEFINITION 7.11 ($O_{\prec,\diamond}(\cdot)$). Let $\mathcal{P}$ be a polynomial of the form

$$\mathcal{P} = \sum_{i=1}^{\phi N} V_i \left(Q_{i\beta}^2 - \frac{1}{N\phi^{1/2}} \right). \quad (7.63)$$

We write $\mathcal{P} = O_{\prec,\diamond}(A)$ if V_i is $Q^{[\beta]}$ -measurable, $|V_i| \leq N^{C_0}$, and $|V_i| \prec A$ for some deterministic A .

DEFINITION 7.12 (GRADED POLYNOMIALS). We write $\mathcal{P} = O_{\prec,\text{even}}(A)$ if $\mathcal{P}$ is a sum of at most C_0 terms of the form

$$A \mathcal{P}_0 \prod_{s=1}^m \mathcal{P}_i, \quad \mathcal{P}_0 = O_{\prec,2n}(1), \quad \mathcal{P}_i = O_{\prec,\diamond}(1),$$

where $n, m \leq C_0$ and A is deterministic.

Moreover, we write $\mathcal{P} = O_{\prec,\text{odd}}(A)$ if $\mathcal{P} = \widehat{\mathcal{P}} \mathcal{P}_{\text{even}}$, where $\widehat{\mathcal{P}} = O_{\prec,1}(1)$ and $\mathcal{P}_{\text{even}} = O_{\prec,\text{even}}(A)$.

Definitions 7.10–7.12 refine Definition 2.1 in the sense that

$$\mathcal{P} = O_{\prec,d}(A) \quad \text{or} \quad \mathcal{P} = O_{\prec,\diamond}(A) \quad \implies \quad \mathcal{P} = P_{\prec}(A). \quad (7.64)$$

Indeed, let $\mathcal{P} = O_{\prec,d}(A)$ be of the form (7.61). Then a simple large deviation estimate (e.g. a trivial extension of [17, Theorem B.1 (iii)]) yields

$$|\mathcal{P}| \prec \left((N\phi^{1/2})^{-d} \sum_{i_1, \dots, i_d} |V_{i_1 \dots i_d}|^2 \right)^{1/2} \prec A,$$

where the last step follows from the definition of admissible weights. Similarly, if $\mathcal{P} = O_{\prec,\diamond}(A)$ is of the form (7.63), a large deviation estimate (e.g. [17, Theorem B.1 (i)]) yields

$$|\mathcal{P}| \prec \left(N^{-2}\phi^{-1} \sum_i |V_i|^2 \right)^{1/2} \prec N^{-1/2} A \leq A.$$

Note that terms of the form $O_{\prec,\cdot}(A)$ satisfy simple algebraic rules. For instance, we have

$$O_{\prec,\text{even}}(A_1) + O_{\prec,\text{even}}(A_2) = O_{\prec,\text{even}}(A_1 + A_2),$$

and

$$O_{\prec,\text{odd}}(A_1) O_{\prec,\text{even}}(A_2) = O_{\prec,\text{odd}}(A_1 A_2)$$

after possibly increasing C_0 . (As with the standard big O notation, such expressions are to be read from left to right.) We stress that such operations may be performed an arbitrary, but bounded, number of times. It is a triviality that all of the following arguments will involve at most C_0 such algebraic operations on graded polynomials, for large enough C_0 .

The point of the graded polynomials is that bounds of the form (7.64) are improved if d is odd and we take the expectation. The precise statement is the following.

LEMMA 7.13. Let $\mathcal{P} = O_{\prec, \text{odd}}(A)$ for some deterministic $A \leq N^C$. Then for any fixed $D > 0$ we have

$$|\mathbb{E}\mathcal{P}| \prec N^{-1/2}A + N^{-D}.$$

PROOF. It suffices to consider the case $A = 1$ and

$$\mathcal{P} = \sum_{i_0} W_{i_0} Q_{i_0\beta} \sum_{i_1, \dots, i_d} V_{i_1 \dots i_d} Q_{i_1\beta} \dots Q_{i_d\beta} \prod_{l=d+1}^{d+m} \left(\sum_{i_l} V_{i_l}^{(l)} \left(Q_{i_l\beta}^2 - \frac{1}{N\phi^{1/2}} \right) \right),$$

where d is even. We suppose that $|W_{i_0}| \prec \varrho_{i_0}^{(0)}$, $|V_{i_1 \dots i_d}| \prec \varrho_{i_1}^{(1)} \dots \varrho_{i_d}^{(d)}$, and $|V_{i_{d+l}}^{(d+l)}| \prec 1$ for $l = d+1, \dots, d+m$. Here $\varrho_{i_k}^{(k)}$ denotes an admissible weight (see Definition 7.9). Thus we have

$$|\mathbb{E}\mathcal{P}| \prec \sum_{i_0, \dots, i_{d+m}} \varrho_{i_0}^{(0)} \dots \varrho_{i_d}^{(d)} \left| \mathbb{E} \left(Q_{i_0\beta} \dots Q_{i_d\beta} \prod_{l=d+1}^{d+m} \left(Q_{i_l\beta}^2 - \frac{1}{N\phi^{1/2}} \right) \right) \right| + N^{-D},$$

where the term N^{-D} comes from the trivial deterministic bound $|V_{i_1 \dots i_d}| \leq N^C$ on the low-probability event of $\prec$ in (7.62), and analogous bounds for the other $Q^{[\beta]}$ -measurable coefficients.

The expectation imposes that each summation index $i_0, \dots, i_{d+m}$ coincide with at least one other one. Thus we get

$$|\mathbb{E}\mathcal{P}| \prec \sum_{i_0, \dots, i_{d+m}} \tilde{\varrho}_{i_0}^{(0)} \dots \tilde{\varrho}_{i_d}^{(d)} I(i_0, \dots, i_{d+m}) \frac{1}{(N\phi^{1/2})^m} + N^{-D}, \quad (7.65)$$

where the indicator function $I(\cdot)$ imposes the condition that each summation index must coincide with at least another one, and we introduced the weight $\tilde{\varrho}_i^{(k)} := N^{-1/2}\phi^{-1/4}\varrho_i$. Note that

$$\sum_i \tilde{\varrho}_i^{(k)} \leq N^{1/2}\phi^{1/2}, \quad \sum_i (\tilde{\varrho}_i^{(k)})^2 \leq 1, \quad \sum_i (\tilde{\varrho}_i^{(k)})^q \leq N^{-q/6} \quad (q \geq 3). \quad (7.66)$$

Here for $q > 3$ we used the inequality $\|\tilde{\varrho}^{(k)}\|_{\ell^q} \leq \|\tilde{\varrho}^{(k)}\|_{\ell^p}$ for $q \geq p$. The indicator function I on the right-hand side of (7.65) imposes a reduction in the number of independent summation indices. We may write $I = \sum_P I_P$ as a sum over all partitions P of the set $\llbracket 0, d+m \rrbracket$ with blocks of size at least two, whereby

$$I_P(i_1, \dots, i_{d+m}) = \prod_{p \in P} \mathbf{1}(i_k = i_l \text{ for all } k, l \in p).$$

Hence the summation over $i_1, \dots, i_{d+m}$ factors into a product over the blocks of P . We shall show that the contribution of each block is at most one, and that there is a block whose contribution is at most $N^{-1/2}$.

Fix $p \in P$ and denote by S_p the contribution of the block p to the summation in the main term of (7.65). Define $s := |p \cap \llbracket 0, d \rrbracket|$ and $t := |p \cap \llbracket d+1, d+m \rrbracket|$. By definition of P , we have $s+t \geq 2$. By the inequality of arithmetic and geometric means, we have

$$S_p \leq \max_k \sum_i (\tilde{\varrho}_i^{(k)})^s \frac{1}{(N\phi^{1/2})^t}.$$

Using (7.66) it is easy to conclude that

$$S_p \leq \begin{cases} 1 & \text{if } (s, t) = (2, 0) \\ N^{-1/2} & \text{if } (s, t) \neq (2, 0). \end{cases}$$

Moreover, since d is even, at least one block of P satisfies $(s, t) \neq (2, 0)$.

Thus we find that

$$|\mathbb{E}\mathcal{P}| \prec \sum_P \prod_{p \in P} S_p + N^{-D} \leq C_{d+m} N^{-1/2} + N^{-D}.$$

Since $d+m \leq 2C_0$, the proof is complete. $\square$

In order to apply Lemma 7.13 to the terms Y from (7.57), we need to expand Y in terms of graded polynomials. This expansion is summarized in the following result, which gives the polynomializations of the coefficients of the terms from (7.57). For an arbitrary unit vector $\mathbf{v} \in \mathbb{R}^M$ we define the control parameter

$$\Psi^{\mathbf{v}} := \Psi + (N^{-1}\|\mathbf{v}\|_{\infty})^{1/3}, \quad \|\mathbf{v}\|_{\infty} := \max_i |v(i)|.$$

LEMMA 7.14. *Fix $D > 0$. Then there exists $C_0 = C_0(D)$ such that for any unit vector $\mathbf{v} \in \mathbb{R}^M$ we have*

$$T_{\mathbf{v}\mathbf{v}} = T_{\mathbf{v}\mathbf{v}}^{[\beta]} + O_{\prec, \text{even}}(\phi^{-1}(\Psi^{\mathbf{v}})^2) + O_{\prec}(N^{-D}), \quad (7.67)$$

$$T_{bb} = O_{\prec, \text{even}}(\phi^{-1/2}) + O_{\prec}(N^{-D}), \quad (7.68)$$

$$T_{\mathbf{w}b} = O_{\prec, \text{even}}(\phi^{-1}\Psi_b) + O_{\prec}(N^{-D}), \quad (7.69)$$

$$(TQ)_{\mathbf{v}\beta} = O_{\prec, \text{odd}}(\phi^{-1/4}\Psi^{\mathbf{v}}) + O_{\prec}(N^{-D}), \quad (7.70)$$

$$(Q^*TQ)_{\beta\beta} = O_{\prec, \text{even}}(\phi^{1/2}) + O_{\prec}(N^{-D}), \quad (7.71)$$

uniformly for $z \in \mathbf{S}$.

PROOF. We begin by noting that (3.10) applied to $X^{[\mu]}$ and (3.21) combined with a large deviation estimate (see [17, Theorem B.1]) yields

$$(X^*G^{[\beta]}X)_{\beta\beta} = \phi^{1/2}m_{\phi^{-1}} + O_{\prec}(\phi^{-1/2}\Psi).$$

Using (7.42) and Lemma 7.5, it is not hard to deduce that

$$(Q^*T^{[\beta]}Q)_{\beta\beta} = \phi^{1/2}m_{\phi^{-1}} + O_{\prec}(\phi^{-1/2}\Psi).$$

Thus for any fixed n we may expand

$$\begin{aligned} -\frac{1}{1 + (Q^*T^{[\beta]}Q)_{\beta\beta}} &= -\frac{1}{1 + \phi^{1/2}m_{\phi^{-1}} - (\phi^{1/2}m_{\phi^{-1}} - (Q^*T^{[\beta]}Q)_{\beta\beta})} \\ &= -\sum_{k=0}^n (zm_{\phi})^{k+1} (\phi^{1/2}m_{\phi^{-1}} - (Q^*T^{[\beta]}Q)_{\beta\beta})^k + O_{\prec}(\phi^{1/2}\Psi^{n+1}), \end{aligned}$$

where in the second step we used (3.5) and (3.22). Now we split

$$\begin{aligned} (Q^*T^{[\beta]}Q)_{\beta\beta} - \phi^{1/2}m_{\phi^{-1}} &= \sum_{i \neq j} T_{ij}^{[\beta]} Q_{i\beta} Q_{j\beta} + \sum_i T_{ii}^{[\beta]} \left(Q_{i\beta} - \frac{1}{N\phi^{1/2}} \right) + \sum_i \frac{1}{N\phi^{1/2}} (T_{ii}^{[\beta]} - m_{\phi^{-1}}) \\ &= O_{\prec, 2}(\phi^{-1/2}\Psi) + O_{\prec, \diamond}(\phi^{-1/2}) + O_{\prec, 0}(\phi^{-1/2}\Psi) \\ &= O_{\prec, \text{even}}(\phi^{-1/2}), \end{aligned}$$

where in the second step we used the estimates $|T_{ij}^{[\beta]} - \delta_{ij}m_{\phi^{-1}}| \prec \phi^{-1}\Psi$ and $|m_{\phi^{-1}}| \leq C\phi^{-1/2}$. Since $|zm_{\phi}| \leq C\phi^{1/2}$, we therefore conclude that

$$-\frac{1}{1 + (Q^*T^{[\beta]}Q)_{\beta\beta}} = O_{\prec, \text{even}}(n\phi^{1/2}) + O_{\prec}(\phi^{1/2}\Psi^{n+1}).$$

From (7.38) and the definition of η , we readily find that $\Psi \leq N^{-c\tau}$ for some constant c . Therefore choosing $n \equiv n(\tau, D)$ large enough yields

$$-\frac{1}{1 + (Q^*T^{[\beta]}Q)_{\beta\beta}} = O_{\prec, \text{even}}(\phi^{1/2}) + O_{\prec}(\phi^{1/2}N^{-D}). \quad (7.72)$$

Having established (7.72), the remainder of the proof is relatively straightforward. From (7.36) and (7.37) we get

$$(TQ)_{\mathbf{v}\beta} = \frac{1}{1 + (Q^*T^{[\beta]}Q)_{\beta\beta}} (T^{[\beta]}Q)_{\mathbf{v}\beta}.$$

Moreover, using that $T_{\mathbf{v}i}^{[\beta]} = v(i)m_{\phi^{-1}} + O_{\prec}(\phi^{-1}\Psi) = O_{\prec}(\phi^{-1/2}|v(i)| + \phi^{-1}\Psi)$, we find

$$\frac{1}{N\phi^{1/2}} \sum_i |T_{\mathbf{v}i}^{[\beta]}|^2 \prec \phi^{-3/2}\Psi^2, \quad \frac{1}{N^{3/2}\phi^{3/4}} \sum_i |T_{\mathbf{v}i}^{[\beta]}|^3 \prec N^{-1/2}\phi^{-9/4}(\Psi^{\mathbf{v}})^3.$$

We conclude that

$$(T^{[\beta]}Q)_{\mathbf{v}\beta} = \sum_i T_{\mathbf{v}i}^{[\beta]} Q_{i\beta} = O_{\prec,1}(\phi^{-3/4}\Psi^{\mathbf{v}}). \quad (7.73)$$

Now (7.70) follows easily from (7.73) and (7.72).

Moreover, (7.67) and (7.69) follow from (7.35) combined with (7.73) and (7.72). For (7.69) we estimate the second term in (7.35) by

$$O_{\prec,\text{even}}(\phi^{1/2}\phi^{-3/2}\Psi^{\mathbf{w}}\Psi^{\mathbf{e}_b}) = O_{\prec,\text{even}}(\phi^{-1}(\Psi + N^{-1/3})^2) = O_{\prec,\text{even}}(\phi^{-1}\Psi),$$

where in the last step we used that $\Psi \geq N^{-1/2}$. Moreover, (7.68) is a trivial consequence of (7.67). Finally, (7.71) follows from (7.35) and (7.72) combined with

$$(Q^*T^{[\beta]}Q)_{\beta\beta} = O_{\prec,2}(1).$$

This concludes the proof. $\square$

Note that the upper bounds in Lemma 7.14 are the same as those of (7.50), except that Ψ is replaced with the larger quantity $\Psi^{\mathbf{v}}$. In order to get back to Ψ from $\Psi^{\mathbf{v}}$, we use the following trivial result.

LEMMA 7.15. *We have*

$$\Psi^{\mathbf{v}} \prec \Psi$$

if

$$\Psi \geq N^{-1/3} \quad \text{or} \quad \|\mathbf{v}\|_{\infty} \prec N^{-1/2}. \quad (7.74)$$

PROOF. The claim follows immediately from the upper bound $\Psi \geq N^{-1/2}$, valid for all $z \in \mathbf{S}$. $\square$

In each application of Lemma 7.14, we shall verify one of the conditions of (7.74). The first condition is verified for $\eta \leq N^{-2/3}$, which always holds for the coefficients of x_1, x_2 , and x_3 (recall (7.2)).

The second condition of (7.74) will be verified when computing the coefficients of y_1, y_2 , and y_3 . To that end, we make use of the freedom of the choice of basis when computing the trace in the definition of J_1, J_2 , and J_3 . We shall choose a basis that is completely delocalized. The following simple result guarantees the existence of such a basis.

LEMMA 7.16. *There exists an orthonormal basis $\mathbf{w}_1, \dots, \mathbf{w}_M$ of $\mathbb{R}^M$ satisfying*

$$|w_i(j)| \prec M^{-1/2} \quad (7.75)$$

uniformly in i and j .

PROOF. Let the matrix $[\mathbf{w}_1 \cdots \mathbf{w}_M]$ of orthonormal basis vectors be uniformly distributed on the orthogonal group $O(M)$. Then each $\mathbf{w}_i$ is uniformly distributed on the unit sphere, and by standard Gaussian concentration arguments one finds that $|w_i(j)| \prec M^{-1/2}$. In particular, there exists an orthonormal basis $\mathbf{w}_1, \dots, \mathbf{w}_M$ satisfying (7.75). In fact, a slightly more careful analysis shows that one can choose $|w_i(j)| \leq (2 + \varepsilon)(\log M)^{1/2}M^{-1/2}$ for any fixed $\varepsilon > 0$ and large enough M . $\square$

We may now derive estimates on the matrix T^2 by writing $(T^2)_{jk} = \sum_i T_{j\mathbf{w}_i} T_{\mathbf{w}_i k}$, where $\{\mathbf{w}_i\}$ is a basis satisfying (7.75). From Lemmas 7.14 and 7.15 we get the following result.

LEMMA 7.17. Fix $D > 0$. Then there exists $C_0 = C_0(D)$ such that

$$\mathrm{Tr} T = \mathrm{Tr} T^{[\beta]} + O_{\prec, \text{even}}(\phi^{-1} N \Psi^2) + O_{\prec}(N^{-D}), \quad (7.76)$$

$$(T^2)_{bb} = O_{\prec, \text{even}}(N \phi^{-1} \Psi^2) + O_{\prec}(N^{-D}), \quad (7.77)$$

$$(T^2 Q)_{b\beta} = O_{\prec, \text{odd}}(\phi^{-1/4} N \Psi^2) + O_{\prec}(N^{-D}), \quad (7.78)$$

$$(Q^* T^2 Q)_{\beta\beta} = O_{\prec, \text{even}}(\phi^{1/2} N \Psi^2) + O_{\prec}(N^{-D}), \quad (7.79)$$

uniformly for $z \in \mathbf{S}$.

PROOF. We prove (7.78); the other estimates are proved similarly. We choose a basis $\mathbf{w}_1, \dots, \mathbf{w}_M$ as in Lemma 7.16, and write

$$(T^2 Q)_{b\beta} = \sum_i T_{b\mathbf{w}_i}(TQ)_{\mathbf{w}_i\beta} = \sum_{i=1}^{N\phi} O_{\prec, \text{even}}(\phi^{-1} \Psi + \phi^{-1/2} |w_i(b)|) O_{\prec, \text{odd}}(\phi^{-1/4} \Psi) + O_{\prec}(N^{-D}),$$

where we used (7.69) with $\mathbf{w}$ replaced by $\mathbf{w}_i$, (7.70), and Lemma 7.15. Summing over i , and recalling that $\Psi \geq N^{-1/2}$, it is easy to conclude (7.78). $\square$

In particular, as in (7.47) we find

$$J_1, J_3 = O_{\prec, \text{odd}}(\phi^{-1/4} N \Psi^2) + O_{\prec}(N^{-D}), \quad J_2 = O_{\prec, \text{even}}(N \Psi^2) + O_{\prec}(N^{-D}), \quad (7.80)$$

where the parity of J_i follows easily from its definition.

The estimates from Lemma 7.14 are compatible with integration in the following sense. Suppose that $\mathcal{P}(s)$ depends on a parameter $s \in S$, where $S \subset \mathbb{R}^k$ has bounded volume, and that $\mathcal{P}(s) = O_{\prec, *}(A(s)) + O_{\prec}(N^{-D})$ uniformly in $s \in S$, where $A(s)$ is a deterministic function of s and $*$ $\in \{\text{even}, \text{odd}\}$ denotes the parity of $\mathcal{P}$. Suppose in addition that $\mathcal{P}(s)$ is Lipschitz continuous with Lipschitz constant N^C . Then, analogously to Remark 3.3, we have

$$\int_S \mathcal{P}(s) ds = O_{\prec, *} \left(\int_S A(s) ds \right) + O_{\prec} \left(\int_S A(s) ds N^{-D} \right).$$

Lemmas 7.14 and 7.17 are the key estimates of the coefficients appearing in (7.57). We claim that all estimates of Lemma 7.7, along with (7.50), remain valid, in the sense that an estimate of the form $|u| \prec v$ is to be replaced with

$$u = O_{\prec, *}(v) + O_{\prec}(N^{-D}),$$

where $*$ $\in \{\text{even}, \text{odd}\}$ denotes the parity of polynomialization of u . Indeed, for the estimates (7.45) on x_i , we always have $\mathrm{Im} z = \eta \leq N^{-2/3}$, so that by Lemma 7.15 we have $\Psi^{\mathbf{v}} \prec \Psi$. Thus we get from Lemma 7.14 that

$$x_1, x_3 = O_{\prec, \text{odd}}(\phi^{-1/4} N \Psi \Psi_b) + O_{\prec}(N^{-D}), \quad x_2 = O_{\prec, \text{even}}(\phi^{-1/2} N \Psi_b^2 + N \Psi^2) + O_{\prec}(N^{-D}),$$

where the parity of x_i may be easily deduced from their definitions. Moreover, for the estimates (7.49) we use (7.80) to get

$$y_1, y_3 = O_{\prec, \text{odd}}(\phi^{-1/4} N^{C\varepsilon} \kappa_a^{1/2}) + O_{\prec}(N^{-D}), \quad y_2 = O_{\prec, \text{even}}(N^{C\varepsilon} \kappa_a^{1/2}) + O_{\prec}(N^{-D}).$$

Note that, thanks to Lemmas 7.14 and 7.17, we have obtained exactly the same upper bounds on the coefficients x_i and y_i as the ones obtained in Lemma 7.7, but we have in addition expressed them, up to a negligible error, as graded polynomials, to which Lemma 7.13 is applicable.

In addition to the coefficients x_i and y_i , we have to control the coefficient $q^{(m)}(y_T)$ in the definition (7.53) of A_1 . We in fact claim that

$$q^{(m)}(y_T) = O_{\prec, \text{even}}(N^{C\varepsilon}) + O_{\prec}(N^{-D}). \quad (7.81)$$

This follows from the estimate

$$y_T = y_{T^{[\beta]}} + O_{\prec, \text{even}}(N^{C\varepsilon} \kappa_a) + O_{\prec}(N^{-D}) = O_{\prec, \text{even}}(N^{C\varepsilon}) + O_{\prec}(N^{-D}),$$

which may be derived from (7.76), combined with a Taylor expansion of $q^{(m)}$. Similarly, we find that

$$h^{(k)}(A_0) = O_{\prec, \text{even}}(N^{C\varepsilon}) + O_{\prec}(N^{-D}) \quad (k = 1, 2, 3). \quad (7.82)$$

We may now put everything together. Noting that the degree of the polynomializations of the expressions (7.57) is always odd, we obtain, in analogy to (7.59) that

$$Y = O_{\prec, \text{odd}}\left(N^{C\varepsilon}\left(\kappa_a^{-1/2}\phi^{3/4}|w(b)|^2 + \phi^{1/4}|w(b)| + \phi^{-1/4}\kappa_a^{1/2}\right)\right) + O_{\prec}(N^{-D})$$

for Y being any term of (7.57). Hence Lemma 7.8 follows from Lemma 7.13 and Young's inequality.

7.4. Stability of level repulsion: proof of Lemma 6.6. This is a Green function comparison argument, using the machinery introduced in Section 7.1. A similar comparison argument was given in Propositions 2.4 and 2.5 of [27]. The details in the sample covariance case and for indices a satisfying $a \leq K^{1-\tau}$ follow an argument very similar to (in fact simpler than) the one from Sections 7.1–7.3. As in the proofs of Propositions 2.4 and 2.5 of [27], one writes the level repulsion condition in terms of resolvents. In our case, one uses the representation (7.22) as the starting point. Then the machinery of Sections 7.1–7.3 may be applied with minor modifications. We omit the details.

8. Extension to general T and universality for the uncorrelated case

In this section we relax the assumption (3.1), and hence extend all arguments of Sections 3–7 to cover general T . We also prove the fixed-index joint eigenvector-eigenvalue universality of the matrix H defined in (2.12), for indices bounded by $K^{1-\tau}$ for some $\tau > 0$.

Bearing the applications in the current paper in mind, we state the results of this section for the matrix H from (2.12), but it is a triviality that all results and their proofs carry over to case of arbitrary Q from (2.1) provided that $\Sigma = TT^* = I_M$.

8.1. The isotropic Marchenko-Pastur law for YY^* . We start with the singular value decomposition of T , which we write as

$$T = O'(\Lambda, 0)O'' = O'\Lambda(I_M, 0)O'',$$

where $O' \in O(M)$ and $O'' \in O(M+r)$ are orthogonal matrices, 0 is the $M \times r$ zero matrix, and Λ is an $M \times M$ diagonal matrix containing the singular values of T . Setting

$$\Sigma^{1/2} = O'\Lambda(O')^*, \quad O := \begin{pmatrix} O' & 0 \\ 0 & I_r \end{pmatrix} O'', \quad (8.1)$$

we have

$$T = \Sigma^{1/2}(I_M, 0)O.$$

We conclude that

$$Q = \Sigma^{1/2}H\Sigma^{1/2},$$

where $H := YY^*$ and $Y := (I_M, 0)OX$ were defined in (2.12). Comparing this to (3.2), we find that to relax the assumption (3.1) we have to generalize the arguments of Sections 3–7 by replacing XX^* with $H = YY^*$.

The generalization of $G = (XX^* - z)^{-1}$ is the resolvent of YY^* ,

$$\widehat{G}(z) := (YY^* - z)^{-1}.$$

We also abbreviate

$$G' := (OXX^*O^* - z)^{-1}.$$

Throughout the following we identify $\mathbf{w} \in \mathbb{R}^M$ with its natural embedding $\begin{pmatrix} \mathbf{w} \\ 0 \end{pmatrix} \in \mathbb{R}^{M+r}$. Thus, for example, for $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$ we may write $G'_{\mathbf{vw}}$.

THEOREM 8.1 (LOCAL LAWS FOR YY^*). *Theorem 3.2 remains valid with G replaced by $\widehat{G}$. Moreover, Theorems 3.4 and 3.5 remain valid for ζ_i and λ_i denoting the eigenvectors and eigenvalues of YY^* .*

PROOF. It suffices to prove the first sentence, since all claims in the second sentence follow from the isotropic law (see [9] for more details). We only prove (3.10) for $\widehat{G}$; the other bound, (3.11) for $\widehat{G}$, is proved similarly. To simplify the presentation, we suppose that $r = 1$; the case $r \geq 2$ is a trivial extension. Abbreviate $\bar{M} := M + 1$. Noting that $Y_{i\mu} = \mathbf{1}(i \neq \bar{M})(OX)_{i\mu}$, we find from [9, Definition 3.5 and Equation (3.7)] that

$$\widehat{G}_{\mathbf{vw}} = G'_{\mathbf{vw}} - \frac{G'_{\mathbf{v}\bar{M}}G'_{\bar{M}\mathbf{w}}}{G'_{\bar{M}\bar{M}}}. \quad (8.2)$$

For definiteness, we focus on (3.10) for $\widehat{G}$; the proof of (3.11) for $\widehat{G}$ is similar. Since $G' = OGO^*$, we have $G'_{\mathbf{vw}} = G_{O^*\mathbf{v}O^*\mathbf{w}}$. Hence, using (3.10) and (8.2), the proof will be complete provided we can show that

$$\left| \frac{G'_{\mathbf{v}\bar{M}}G'_{\bar{M}\mathbf{w}}}{G'_{\bar{M}\bar{M}}} \right| \prec \Phi, \quad \Phi := \sqrt{\frac{\operatorname{Im} m_{\phi^{-1}}(z)}{M\eta}} + \frac{1}{M\eta} \prec \phi^{-1}\Psi, \quad (8.3)$$

where we recall the definition (7.38) of Ψ . In fact, from Lemma 3.6 and (3.22) we find that $\Phi/|m_{\phi^{-1}}| \leq N^{-c}$ for some positive constant c depending on τ . Hence (3.10) yields

$$\left| \frac{G'_{\mathbf{v}\bar{M}}G'_{\bar{M}\mathbf{w}}}{G'_{\bar{M}\bar{M}}} \right| \prec \Phi^2/|m_{\phi^{-1}}| \leq \Phi.$$

This concludes the proof. $\square$

Having established Theorem 8.1, all arguments from Sections 3–6 that use it as input may be taken over verbatim, after replacing G by $\widehat{G}$. More precisely, all results from Sections 3–6 remain valid for a general Q , with the exception of Proposition 6.4, Lemmas 6.5 and 6.6, and Proposition 6.7. Therefore we have completed the proofs of all of our main results except Theorem 2.18.

In order to prove Theorem 2.18, we still have to prove Lemmas 6.5 and 6.6 and Proposition 6.7 for YY^* instead of XX^* . Lemma 6.5 is easy: for Gaussian X we have $Y \stackrel{d}{=} (I_M, 0)X = \widetilde{X}$, where $\widetilde{X}$ is the $M \times N$ matrix obtained from X by deleting its bottom r rows.

The proofs of Lemma 6.6 and Proposition 6.7 rely on Green function comparison. What remains, therefore, is to extend the argument of Section 7 from $H = XX^*$ to $H = YY^*$.

8.2. Quantum unique ergodicity for YY^* . In this subsection we prove Proposition 6.7 for the eigenvectors ζ_a of $H = YY^*$. As explained in Section 7.4, the proof of Lemma 6.6 is analogous and therefore omitted. We proceed exactly as in Section 7, replacing G with $\widehat{G}$. It suffices to prove the following result.

LEMMA 8.2. *Lemma 7.3 remains valid if $x(E)$ and $y(E)$ are replaced with $\widehat{x}(E)$ and $\widehat{y}(E)$, obtained from the definitions (7.29) and (7.30) by replacing G with $\widehat{G}$.*

PROOF. We take over the notation from the proof of Theorem 8.1, and to simplify notation again assume that $r = 1$. As in Section 7, we suppose for definiteness that $\phi \geq 1$. Defining $\mathbf{u} := O\mathbf{w}$ and $\mathbf{r} := O\mathbf{e}_{M+1}$, we have $\langle \mathbf{u}, \mathbf{r} \rangle = 0$ and, using (8.2),

$$\widehat{G}_{\mathbf{ww}} = G_{\mathbf{uu}} - \frac{G_{\mathbf{ur}}G_{\mathbf{ru}}}{G_{\mathbf{rr}}}, \quad \operatorname{Tr} \widehat{G} = \operatorname{Tr} G - \frac{(G^2)_{\mathbf{rr}}}{G_{\mathbf{rr}}}.$$

We conclude that

$$\widehat{x}(E) = x(E) - \frac{M}{\pi} \operatorname{Im} \left(\frac{G_{\mathbf{ur}}G_{\mathbf{ru}}}{G_{\mathbf{rr}}} \right) (E + i\eta).$$

Recalling (3.10) and (7.38), we find that the second term is stochastically dominated by

$$M \frac{\phi^{-2}\Psi^2}{\phi^{-1/2}} \leq N\Psi^2 \leq CN \frac{1}{N^2\eta^2} = \frac{CN^{4\varepsilon}}{N\Delta_a} \frac{1}{\Delta_a} \leq CN^{4\varepsilon} \kappa_a^{1/2} \frac{1}{\Delta_a},$$

where in the second step we used that $\Psi \leq C(N\eta)^{-1}$, as follows from Lemma 3.6 and the definition of η in (7.2). Recalling the definitions from (7.2), we therefore conclude that for small enough $\varepsilon \equiv \varepsilon(\tau)$ we have

$$|I| \sup_{E \in I} |\hat{x}(E) - x(E)| \prec N^{-c} \quad (8.4)$$

for some positive constant c depending on τ .

Similarly, we have for any $z \in \mathbf{S}$

$$|\mathrm{Tr} G - \mathrm{Tr} \hat{G}| \prec N\Psi^2 \leq N^{-c}\eta^{-1} \quad (8.5)$$

for some positive constant c depending on τ . Plugging (8.5) into the definition of $\hat{y}(E)$ and estimating the error term using integration by parts, as in (7.51), we get

$$\sup_{E \in I} |\hat{y}(E) - y(E)| \prec N^{-c}.$$

Using the mean value theorem and the bound $|y(E)| \prec 1$, we therefore get

$$h \left[\int_I \hat{x}(E) q(\hat{y}(E)) \mathrm{d}E \right] = h \left[\int_I x(E) q(y(E)) \mathrm{d}E \right] + O_{\prec}(N^{-c}).$$

Combined with (7.28), this concludes the proof. $\square$

This concludes the proof of Theorem 2.18 for the case of general T .

8.3. The joint eigenvalue-eigenvector universality of YY^* near the spectral edges. In this section we observe that the technology developed in Section 7 allows us to establish the universality of the joint eigenvalue-eigenvector distribution of Q provided that $\Sigma = I_M$. Without loss of generality, we consider the case where Q is given by $H = YY^*$ defined in (2.12). This result applies to arbitrary eigenvalue and eigenvector indices which are bounded by $K^{1-\tau}$, and does in particular not need to invoke eigenvalue correlation functions.

This result generalizes the quantum unique ergodicity from Proposition 6.7 and its extension from Remark 6.8 by also including the distribution of the eigenvalues. The universality of both the eigenvalues and the eigenvectors is formulated in the sense of fixed indices. A result in a similar spirit was given in [27, Theorem 1.6], except that the upper bound on the eigenvalue and eigenvector indices $(\log K)^{C \log \log N}$ from [27] is improved all the way to $K^{1-\tau}$, for any $\tau > 0$. A result covering all eigenvalue and eigenvector indices, i.e. with an index upper bound K , was given in [27, Theorem 1.10] and [38, Theorem 8], but under the assumption of a four-moment matching assumption. Theorem 8.3 is a true universality result in that it does not require any moment matching assumptions, but it does require an index upper bound of $K^{1-\tau}$ instead of K on the eigenvalue and eigenvector indices.

In addition, Theorem 8.3 extends the previous results from [27] and [38] by considering arbitrary generalized components $\langle \zeta_a, \mathbf{v} \rangle$ of the eigenvectors. Finally, Theorem 8.3 holds for the general class of covariance matrices defined in (2.12).

THEOREM 8.3 (UNIVERSALITY FOR THE UNCORRELATED CASE). *Fix $\tau > 0$, $k = 1, 2, 3, \dots$, and $r = 0, 1, 2, \dots$. Choose an observable $h \in C^4(\mathbb{R}^{2k})$ satisfying*

$$|\partial^\alpha h(x)| \leq C(1 + |x|)^C$$

for some constant $C > 0$ and for all $\alpha \in \mathbb{N}^{2k}$ satisfying $|\alpha| \leq 4$. Let X be an $(M + r) \times N$ matrix, and define H through (2.12) for some orthogonal $O \in \mathrm{O}(M + r)$. Denote by $\lambda_1 \geq \dots \geq \lambda_M$ the eigenvalues of H and by $\zeta_1, \dots, \zeta_M$ the associated unit eigenvectors. Let $\mathbb{E}^{(1)}$ and $\mathbb{E}^{(2)}$ denote the expectations with respect to two laws on X , both of which satisfy (2.4) and (2.5). Recall the definition (6.19) of Δ_a , the typical distance between λ_a and λ_{a+1} , and (3.13) of the classical location γ_a .

Then for any indices $a_1, \dots, a_k, b_1, \dots, b_k \in \llbracket 1, K^{1-\tau} \rrbracket$ and deterministic unit vectors $\mathbf{u}_1, \mathbf{w}_1, \dots, \mathbf{u}_k, \mathbf{w}_k \in \mathbb{R}^M$ we have

$$(\mathbb{E}^{(1)} - \mathbb{E}^{(2)}) h \left(\frac{\lambda_{a_1} - \gamma_{a_1}}{\Delta_{a_1}}, \dots, \frac{\lambda_{a_k} - \gamma_{a_k}}{\Delta_{a_k}}, M \langle \mathbf{u}_1, \boldsymbol{\zeta}_{b_1} \rangle \langle \boldsymbol{\zeta}_{b_1}, \mathbf{w}_1 \rangle, \dots, M \langle \mathbf{u}_k, \boldsymbol{\zeta}_{b_k} \rangle \langle \boldsymbol{\zeta}_{b_k}, \mathbf{w}_k \rangle \right) = O(N^{-c})$$

for some constant $c \equiv c(\tau, k, r, h) > 0$.

PROOF. The proof is a Green function comparison argument, a minor modification of that developed in Section 7. We write the distribution of $\lambda_a - \gamma_a$ in terms of the resolvent $\widehat{G}$, starting from the Helffer-Sjöstrand representation (7.22), exactly as in [27, Sections 4 and 5]. We omit further details. $\square$

REMARK 8.4. In particular, Theorem 8.3 establishes the fixed-index universality of eigenvalues with indices bounded by $K^{1-\tau}$. Indeed, we may choose $\mathbb{E}^{(2)}$ to be the expectation with respect to a Gaussian law, in which case $H \stackrel{d}{=} \widetilde{X} \widetilde{X}^*$, where $\widetilde{X}$ is a $M \times N$ and Gaussian. (For example, the top eigenvalue of H is distributed according to the Tracy-Widom-1 distribution, etc.)

We note that even this fixed-index universality of eigenvalues is a new result, having previously only been established under the four-moment matching condition [27, 38] (in the context of Wigner matrices).

REMARK 8.5. We formulated Theorem 8.3 for the real symmetric covariance matrices of the form (2.12), but it and its proof remain valid for complex Hermitian covariance matrices, as well as Wigner matrices (both real symmetric and complex Hermitian).

REMARK 8.6. Assuming $|\phi - 1| > \tau$, the condition $a \leq K^{1-\tau}$ on the indices in Theorem 8.3 may be replaced with $a \notin \llbracket K^{1-\tau}, K - K^{1-\tau} \rrbracket$.

REMARK 8.7. Combining Theorems 8.3 and 2.6, we get the following universality result for Q . Fix $\tau > 0$, $k = 1, 2, 3, \dots$, and $r = 0, 1, 2, \dots$. For any continuous and bounded function h on $\mathbb{R}^k$ we have

$$\lim_{N \rightarrow \infty} \left[\mathbb{E} h \left(\frac{\mu_{s_+ + a_1} - \gamma_{a_1}}{\Delta_{a_1}}, \dots, \frac{\mu_{s_+ + a_k} - \gamma_{a_k}}{\Delta_{a_k}} \right) - \mathbb{E}^{\text{Wish}} h \left(\frac{\lambda_{a_1} - \gamma_{a_1}}{\Delta_{a_1}}, \dots, \frac{\lambda_{a_k} - \gamma_{a_k}}{\Delta_{a_k}} \right) \right] = 0$$

for any indices $a_1, \dots, a_k \leq K^{1-\tau} \alpha_+^3$. Here $\mathbb{E}^{\text{Wish}}$ denotes expectation with respect to the Wishart case, where $r = 0$, $T = I_M$, and X is Gaussian. A similar result holds near the left edge provided that $|\phi - 1| \geq \tau$.

9. Extension to $\dot{Q}$ and proof of Theorem 2.22

In this section we explain how to extend our analysis from Q defined in (2.1) to $\dot{Q}$ defined in (2.2), hence proving Theorem 2.22. We define the resolvent

$$\dot{G}(z) := (X(1 - \mathbf{e}\mathbf{e}^*)X^* - z)^{-1},$$

which will replace $G(z) = (XX^* - z)^{-1}$ when analysing with $\dot{Q}$ instead of Q . We begin by noting that the isotropic local laws hold for also for $\dot{G}$.

THEOREM 9.1 (LOCAL LAWS FOR $X(1 - \mathbf{e}\mathbf{e}^*)X^*$). *Theorem 3.2 remains valid with G replaced by $\dot{G}$. Moreover, Theorems 3.4 and 3.5 remain valid for ζ_i and λ_i denoting the eigenvectors and eigenvalues of $X(1 - \mathbf{e}\mathbf{e}^*)X^*$.*

PROOF. As in the proof of Theorem 8.1, we only prove (3.10) for $\dot{G}$. Using the identity (3.34) we get

$$\dot{G} = (XX^* - z - X\mathbf{e}\mathbf{e}^*X^*)^{-1} = G + \frac{1}{1 - (X^*GX)_{\mathbf{e}\mathbf{e}}} GX\mathbf{e}\mathbf{e}^*X^*G. \quad (9.1)$$

Using (3.10), the proof will be complete provided we can show that

$$\left| \frac{(GX)_{\mathbf{ve}}(X^*G)_{\mathbf{ew}}}{1 - (X^*GX)_{\mathbf{ee}}} \right| \prec \Phi \quad (9.2)$$

for unit vectors $\mathbf{v}, \mathbf{w} \in \mathbb{R}^M$, where Φ was defined in (8.3). Recall the definition $R(z) := (X^*X - z)^{-1}$. From the elementary identity $X^*GX = 1 + zR$ and Theorem 3.2 applied to X^* instead of X , we get

$$\left| \frac{1}{1 - (X^*GX)_{\mathbf{ee}}} \right| \prec \frac{1}{|z||m_\phi|} \leq \frac{C}{1 + \phi^{1/2}},$$

where in the last step we used (3.18) and (3.22). Using Lemma 9.3 below with $\mathbf{x} = \mathbf{e}$, (9.2) therefore follows provided we can prove that

$$\phi^{1/2}(1 + \phi^{1/2})\Phi^2 \leq C\Phi.$$

This is an immediate consequence of the estimate $(1 + \phi)\Phi \leq C$, which itself easily follows from the definition (3.8) of $\mathbf{S}$ and (3.21). $\square$

Next, we deal with the quantum unique ergodicity of $X(1 - \mathbf{ee}^*)X^*$. As explained in Section 8.2, it suffices to prove the following result.

LEMMA 9.2. *Lemma 7.3 remains valid if $x(E)$ and $y(E)$ are replaced with $\dot{x}(E)$ and $\dot{y}(E)$, obtained from the definitions (7.29) and (7.30) by replacing G with $\dot{G}$.*

PROOF. The proof mirrors closely that of Lemma 8.2, using the identity (9.1) instead of (8.2) as input. We omit the details. $\square$

Using Theorem 9.1 and Lemma 9.2, combined with the results of Section 8.2, we conclude the proof of Theorem 2.22. To be precise, the arguments of Sections 8 and 9 have to be successively combined so as to obtain the isotropic local laws and quantum unique ergodicity of the matrix $Y(1 - \mathbf{ee}^*)Y^*$. This has to be done in the following order. First, using Theorem 9.1 and Lemma 9.2, one establishes the local laws and quantum unique ergodicity for $X(1 - \mathbf{ee}^*)X^*$. Second, using these results as input, one repeats the arguments of Section 8, except that XX^* is replaced with $X(1 - \mathbf{ee}^*)X^*$; this is a trivial modification of the arguments presented in Section 8. Thus we get the local laws and quantum unique ergodicity for the matrix

$$Y(1 - \mathbf{ee}^*)Y^* = (I_M, 0)OX(I_N - \mathbf{ee}^*)X^*O^*(I_M, 0)^*.$$

Moreover, we find that Theorem 8.3 also holds if $H = YY^*$ from (2.12) is replaced with $Y(1 - \mathbf{ee}^*)Y^*$.

All that remains is the proof of the following estimate, which generalizes (7.39).

LEMMA 9.3. *For $z \in \mathbf{S}$ and deterministic unit vectors $\mathbf{v} \in \mathbb{R}^M$ and $\mathbf{x} \in \mathbb{R}^N$, we have*

$$|(GX)_{\mathbf{vx}}| \prec \phi^{1/4}(1 + \phi^{1/2})\Phi, \quad (9.3)$$

where Φ was defined in (8.3).

PROOF. In the case where $\mathbf{x} = \mathbf{e}_\mu$ is a standard unit vector of $\mathbb{R}^N$, (9.3) is a trivial extension of (7.39) (which was proved under the assumption that $\phi \geq 1$). For general $\mathbf{x}$, the proof requires more work. Indeed, writing $(GX)_{\mathbf{vx}} = \sum_\mu (GX)_{\mathbf{v}\mu} u(\mu)$ and estimating $|(GX)_{\mathbf{v}\mu}|$ by $O(\phi^{1/4}(1 + \phi^{1/2})\Phi)$ leads to a bound proportional to the ℓ^1 -norm of $\mathbf{x}$ instead of its ℓ^2 -norm. In order to obtain the sharp bound, which is proportional to the ℓ^2 -norm, we need to exploit cancellations among the summands. This phenomenon is related to the *fluctuation averaging* from [16], and was previously exploited in [9] to obtain the isotropic laws from Theorem 3.2. It is best made use of by estimating the p -th moment for an even integer p ,

$$\mathbb{E}|(GX)_{\mathbf{vx}}|^p = |z|^p \sum_{\mu_1, \dots, \mu_p} x(\mu_1) \cdots x(\mu_p) \mathbb{E} \left(R_{\mu_1 \mu_1} (G^{[\mu_1]} X)_{\mathbf{v} \mu_1} \cdots \overline{R_{\mu_p \mu_p} (G^{[\mu_p]} X)_{\mathbf{v} \mu_p}} \right); \quad (9.4)$$

here we used the first identity of (7.37). A similar argument was given in [9, Section 5]. The basic idea is to make all resolvents on the right-hand side of (9.4) independent of the columns of X indexed by $\{\mu_1, \dots, \mu_p\}$ (see [9, Definition 3.7]). As in [9, Section 5], we do this using the identities from [9, Lemma 3.8] for the entries of R . In addition, for the entries of G we use the identity (in the notation of [9, Definition 3.7])

$$G_{\mathbf{v}\mathbf{w}}^{[T]} = G_{\mathbf{v}\mathbf{w}}^{[T\mu]} + zR_{\mu\mu}^{[T]} \sum_{k,l=1}^M G_{\mathbf{v}\mathbf{k}}^{[T\mu]} G_{\mathbf{l}\mathbf{w}}^{[T\mu]} X_{k\mu} X_{l\mu}, \quad (9.5)$$

which follows from (7.35) and (7.36). As in [9, Section 5], the resulting expansion may be conveniently organized using graphs, and brutally truncated after a number of steps that depends only on p and ω (here ω is the constant from $\mathbf{S}$ in (3.8)). The key observation is that, once the expansion is performed, we may take the pairing among the variables $\{X_{k\mu_i} : k \in \llbracket 1, N \rrbracket, i \in \llbracket 1, p \rrbracket\}$; we find that each independent summation index μ_i comes with a weight bounded by $x(\mu_i)^2 + N^{-1}$, which sums to $O(1)$. We refer to [9] for the full details of the method, and leave the modifications outlined above to the reader. $\square$

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