

MULTIPLICATIVE FREE CONVOLUTION AND INFORMATION-PLUS-NOISE TYPE MATRICES

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Free probability and random matrix theory has shown to be a fruitful combination in many fields of research, such as digital communications, nuclear physics and mathematical finance. The link between free probability and eigenvalue distributions of random matrices will be strengthened further in this paper. It will be shown how the concept of multiplicative free convolution can be used to express known results for eigenvalue distributions of a type of random matrices called Information-Plus-Noise matrices. The result is proved in a free probability framework, and some new results, useful for problems related to free probability, are presented in this context. The connection between free probability and estimators for covariance matrices is also made through the notion of free deconvolution.

1. Introduction. Applications of free probability have been growing rapidly over the last years. Random matrices and their limit eigenvalue distributions is an area where free probability has proved to be useful [7]. Random matrices are a useful tool for modelling systems, for instance in digital communications [19, 20], nuclear physics [6, 8] and mathematical finance [2]. This paper is a contribution to the random matrix facet of free probability, in that the connection between certain random matrices and free probability is clarified further. We will focus on what we call *Information-Plus-Noise Type Matrices*, i.e. random matrices on the form

$$(1.1) \quad W_n = \frac{1}{N}(R_n + \sigma X_n)(R_n + \sigma X_n)^*,$$

where R_n and X_n are independent random matrices of dimension $n \times N$. These can be thought of as sample covariance matrices of random vectors $r_n + \sigma x_n$, where r_n can be interpreted as a vector carrying the information in a system, and x_n additive noise, with σ the strength of the noise. We impose no assumption on independence between samples. We will use some

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common restrictions on the noise: X_n will contain i.i.d. complex entries of unit variance. n and N will be increased so that

$$(1.2) \quad \lim_{n \rightarrow \infty} \frac{n}{N} = c.$$

In [3], Dozier and Silverstein explain how the limit eigenvalue distribution μ_W of the matrix W_n can be found, based on knowledge of the limit eigenvalue distribution μ_Γ of the matrix $\Gamma_n = \frac{1}{N}R_nR_n^*$. The result is expressed in terms of a solution to a function equation (equation (4.1)). We will show that there is an equivalent way of expressing this solution, using the concept of *multiplicative free deconvolution*, denoted by \boxminus (multiplicative free convolution, as well as freeness and asymptotic freeness are defined in section 2). The following is the main result of the paper:

THEOREM 1.1. *Assume that the entries X_{ij}^n of X_n are Gaussian, independent and identically distributed with expectation 0 and variance 1. Assume also that the empirical eigenvalue distribution of $\Gamma_n = \frac{1}{N}R_nR_n^*$ converges in distribution almost surely to a compactly supported probability measure μ_Γ . Then we have that the empirical eigenvalue distribution of W_n also converges in distribution almost surely to a compactly supported probability measure μ_W uniquely identified by*

$$(1.3) \quad \mu_W \boxminus \mu_c = (\mu_\Gamma \boxminus \mu_c) \boxplus \mu_{\sigma^2 I}.$$

Some remarks are needed to explain theorem 1.1. By the *empirical eigenvalue distribution* of an $n \times n$ random matrix X we mean the (random) atomic measure

$$\frac{1}{n} (\delta(\lambda_1(X)) + \cdots + \delta(\lambda_n(X))),$$

where $\lambda_1(X), \dots, \lambda_n(X)$ are the (random) eigenvalues of X . That μ_n converges in distribution to μ means that the moments of μ_n converge to the moments of μ . Theorem 1.1 requires compactly supported measures, and these have moments of all orders.

The conditions in theorem 1.1 are somewhat stronger than those in [3] due to the restriction to measures with compact support. Contrary to [3], we also restrict to noise-matrices with Gaussian entries.

Theorem 1.1 yields a short expression for μ_W , removing the need for solving the equation in [3] directly. It essentially says that the connection between μ_W and μ_Γ can be expressed compactly in the *deconvolved domain*, where the connection can be viewed as a shift of the spectrum with the noise variance σ^2 . The proof of theorem 1.1 is based on methods from free

probability, with some new results established along the way. Some of these deserve extra attention, in particular theorem 3.4. This can be thought of as a version of theorem 1.1 where μ_W and μ_Γ are interpreted as distributions of free random variables.

Theorems 3.2 and 3.3 also deserve some extra attention. These address asymptotic freeness almost everywhere [7] for two random matrices where

1. both converge in distribution almost everywhere to compactly supported limits, and
2. one of the random matrices are standard unitary (theorem 3.2) or Gaussian (theorem 3.3).

These results expand known results from [7] for asymptotic freeness. The proofs of theorems 3.2 and 3.3 use random matrix approximations with deterministic matrices. Asymptotic freeness of Gaussian/standard unitary random matrices and uniformly norm-bounded deterministic matrices are well-known (lemma 4.3.2 in [7]). Unfortunately, norm-bounded deterministic matrices are not able to approximate the random matrices under consideration. We solve the problem by generalizing to matrices satisfying uniform $\|\cdot\|_p$ -norm bounds instead, where $\|\cdot\|_p$ denotes the Schatten p -norm (with respect to tr_n), defined for $p \geq 1$ by $\|A\|_p = tr_n(|A|^p)^{\frac{1}{p}}$ ($A \in M_n(C)$): We prove that matrices satisfying such bounds can be used to approximate our random matrices, and that they also give asymptotic freeness as in lemma 4.3.2 in [7] (theorem 3.1).

Theorem 1.1 is actually proved by combining theorems 3.3 and 3.4 through another approximation argument (see theorem 3.5). While [3] restricts to the distribution of $\frac{1}{N}(R_n + \sigma X_n)(R_n + \sigma X_n)^*$, we show more in that any mixed moments of $\frac{1}{N}R_n R_n^*$ and $\frac{1}{N}X_n X_n^*$ are obtained through our asymptotic freeness results.

Recent works [17, 18] show that multiplicative free convolution also admits an efficient implementation in terms of the moments of the operand measures. The basic results on free probability we need for this are proved in this paper (theorems 2.1 and 2.2). A consequence is that existing computational frameworks can be used in obtaining μ_Γ and μ_W . In [18], μ_Γ and μ_W are illustrated in terms of signal processing applications, and simulations are run using a computational framework building on theorems 2.1 and 2.2. A useful consequence of the link with free probability is that the "inverse problem" (i.e. that of finding μ_Γ from μ_W) can be solved within the same framework, since the framework embraces convolution as well as deconvolution.

The eigenvalue distribution of Γ_n provides us with possibilities for estimat-

ing the covariance matrices of the system through the so-called G -estimators [5]. These will be reviewed, and it will be shown how multiplicative free convolution can be used to rewrite such estimators to a very simple form. It will be apparent from this that the G^2 -estimator actually can be viewed as a step in expressing μ_W from μ_Γ .

While the results mentioned here are hard to prove, some of them should come as no surprise. For instance, [14] has already made the connection between Information-Plus-Noise type matrices and multiplicative free convolution. This paper also indicates that some of the mentioned results are already known, by saying that random matrices with Haar-distributed eigenvectors are asymptotically free from any random matrices independent from them. However, the generality in which this should hold is not indicated. Also, [14] considers only Gaussian matrices, and the connection with already existing estimators of covariance matrices was not made.

This paper is organized as follows. Section 2 contains notation and preliminaries for various free probability tools, like free transforms and combinatorial aspects. The mentioned implementation of free convolution builds on the combinatorial expression of freeness, and the results needed on this are explained in section 2.1. The proof for theorem 1.1 is presented in section 3. A sketch of the proof is first given, followed by the proofs for theorems 3.1, 3.2 3.3 and 3.4. Section 4 first states the results we need from [3], and sketch the proof for the equivalence of these and theorem 1.1. This sketch is then followed by the rest of the details. The various transforms used in free probability (section 2) are used in this direction. In section 5 we state the principles of G -analysis and the expression for the G^2 -estimator. We also prove the theorem which expresses the G^2 -estimator in terms of free probability.

2. Notation and preliminaries. In the following, uppercase symbols will be used for matrices, and $(\cdot)^*$ will denote hermitian transpose. I_n will represent the identity matrix of order n . We will focus here on certain noncommutative probability spaces. A noncommutative probability space is a pair (A, ϕ) where A is a unital $*$ -algebra and ϕ is a normalized (i.e. $\phi(I) = 1$) linear functional on A . The elements of A are called random variables. The probability spaces we will encounter are mostly $(M_n(C), tr_n)$, i.e. $n \times n$ -matrices equipped with the normalized trace. Any matrix can be associated with a probability measure through it's eigenvalue distribution. We will mostly be concerned with probability measures with compact support.

DEFINITION 2.1. *A family of unital $*$ -subalgebras $(A_i)_{i \in I}$ will be called*

a free family if

$$(2.1) \quad \left\{ \begin{array}{l} a_j \in A_{i_j} \\ i_1 \neq i_2, i_2 \neq i_3, \dots, i_{n-1} \neq i_n \\ \phi(a_1) = \phi(a_2) = \dots = \phi(a_n) = 0 \end{array} \right\} \Rightarrow \phi(a_1 \cdots a_n) = 0.$$

(2.1) enables us to calculate the mixed moments of a_1 and a_2 when they are free. In particular, the moments of $a_1 + a_2$ and $a_1 a_2$ can be calculated. This gives us two new probability measures, which depend on the probability measures of a_1, a_2 only (i.e. not on their realizations). Therefore we can define two operations on the set of probability measures: Additive free convolution

$$(2.2) \quad \mu_1 \boxplus \mu_2$$

for the sum of free random variables, and multiplicative free convolution

$$(2.3) \quad \mu_1 \boxtimes \mu_2$$

for the product of free random variables.

Let F^{μ_A} denote the empirical distribution function (e.d.f.) of the eigenvalues of A (so that $F^{\mu_A}(x)$ is the proportion of eigenvalues of A which are $\leq x$). When we have a series of e.d.f.'s $F^{\mu_{A_n}}$, we will use the notation

$$F^{\mu_{A_n}} \xrightarrow{\mathcal{D}} F^\mu$$

for weak convergence, where F^μ is the cumulative distribution function of the measure μ . We will also write a.s. as shorthand notation for almost sure convergence.

Some random matrices and limit distributions occur naturally in many contexts. If the entries of the $n \times N$ (with $\lim_{n \rightarrow \infty} \frac{n}{N} = c$) random matrices W_n have zero mean and unit variance, the empirical eigenvalue distribution of $\frac{1}{N} W_n W_n^*$ converges almost surely to the so-called Marčenko Pastur law μ_c ([21] page 9). These are also called the free Poisson distributions, and are characterized by the density

$$(2.4) \quad f^{\mu_c}(x) = \left(1 - \frac{1}{c}\right)^+ \delta(x) + \frac{\sqrt{(x-a)^+(b-x)^+}}{2\pi c x},$$

where $(z)^+ = \max(0, z)$, $a = (1 - \sqrt{c})^+$ and $b = (1 + \sqrt{c})^+$. Similar notation to the Marčenko Pastur law is used for the distribution μ_a of a random variable a . We avoid confusion by never using c to denote random variables. μ_1 will always mean the Marčenko Pastur law with parameter one. Marčenko

Pastur laws are some of the most basic random matrix building blocks, as they appear as limits for large random matrices in many contexts. This paper will demonstrate that this is indeed the case for the type of systems we consider also.

We will not use the characterization of the Marčenko Pastur law as in (2.4) directly. Rather we will work with equivalent expressions of it through the transforms defined in this section. The transforms we define will only be applied for probability measures with support contained on the positive real line.

The *Stieltjes transform* ([21] page 38) of a probability measure μ is the analytic function on $C^+ = \{z \in C : \Im z > 0\}$ defined by

$$(2.5) \quad m_\mu(z) = \int_{-\infty}^{\infty} \frac{1}{\lambda - z} dF^\mu(\lambda).$$

A convenient inversion formula for the Stieltjes transform also exists, so that m_μ uniquely identifies μ . If μ is assumed to have nonnegative support, m_μ can be analytically continued to the negative part of the real line. If $\mu = \mu_X$ for a non-negative random variable X , m_μ is strictly monotone on the negative real line, taking values in the interval $[0, E\left(\frac{1}{X}\right)]$. We will use the fact that if we know $m_\mu(z)$ in an interval $(-z, 0)$ for $z < 0$, we also know m_μ for all other values of z , and hence we also know μ (use the Stieltjes inversion formula).

The η -transform ([21] page 40) is defined for measures μ with support on the positive real line, and for nonnegative real numbers by

$$(2.6) \quad \eta_\mu(z) = \int_{-\infty}^{\infty} \frac{1}{1 + z\lambda} dF^\mu(\lambda).$$

$\eta(z)$ is a strictly monotonically decreasing function. As such it simplifies many derivations and statements of results. The inverse is tightly connected to the S -transform (see below). It's connection with the Stieltjes transform is

$$(2.7) \quad \eta_\mu(z) = \frac{m_\mu(-\frac{1}{z})}{z}, \quad m_\mu(z) = -\frac{\eta_\mu(-\frac{1}{z})}{z}.$$

Therefore $\eta_\mu(z)$ uniquely identifies $m_\mu(z)$, since $m_\mu(z)$ for real, negative z can be continued analytically to C^+ . We will use the fact that if we know $\eta_\mu(z)$ in an interval $(0, z)$ for $z > 0$, we also know μ .

The R -transform ([21] page 48) has domain of definition C^+ and can be defined in terms of the Stieltjes transform as

$$(2.8) \quad \mathcal{R}_\mu(z) = m_\mu^{-1}(-z) - \frac{1}{z}.$$

The importance of the R -transform comes from its additive property for the distribution of the sum of free random variables A_1 and A_2 ,

$$(2.9) \quad \mathcal{R}_{\mu_{a_1+a_2}}(z) = \mathcal{R}_{\mu_{a_1}}(z) + \mathcal{R}_{\mu_{a_2}}(z).$$

Slightly different versions of the R -transform are encountered in the literature. The one above is from [21]. In connection with free combinatorics, another definition is used, namely $R_\mu(z) = z\mathcal{R}_\mu(z)$. Of course, $R_\mu(z)$ also satisfies (2.9).

The S -transform ([21] page 50) is defined on $(-1, 0)$. It can be defined in terms of the η -transform by

$$(2.10) \quad S_\mu(z) = -\frac{z+1}{z}\eta_\mu^{-1}(z+1).$$

The Marčenko Pastur law (2.4) can be shown to have S -transform $S_{\mu_c}(z) = \frac{1}{1+cz}$ ([21] page 51). The importance of the S -transform comes from its multiplicative property for the distribution of the product of free random variables a_1 and a_2 :

$$(2.11) \quad S_{\mu_{a_1a_2}}(z) = S_{\mu_{a_1}}(z)S_{\mu_{a_2}}(z).$$

If the values of $\eta_\mu(z)$ or $S_\mu(z)$ are known in an interval, one also knows μ .

Freeness, additive and multiplicative free convolution have a combinatorial description involving these transforms which we will use for in some of our proofs. These combinatorial descriptions build on the concept of *non-crossing partitions*:

DEFINITION 2.2. *A partition π is called noncrossing if whenever we have $i < j < k < l$ with $i \sim k$, $j \sim l$ (\sim meaning belonging to the same block), we also have $i \sim j \sim k \sim l$ (i.e. i, j, k, l are all in the same block). The set of noncrossing partitions of $\{1, \dots, n\}$ is denoted $NC(n)$.*

$NC(n)$ becomes a lattice under the refinement order of partitions. An ingredient we need in making the connections between freeness and the noncrossing partitions is the complementation map of Kreweras, which is a lattice anti-isomorphism of $NC(n)$. To define this we need the circular representation of a partition: We mark n equidistant points $1, \dots, n$ (numbered clockwise) on the circle, and form the convex hull of points lying in the same block of the partition. This gives us a number of convex sets H_i , equally many as there are blocks in the partition, which do not intersect if and only if the partition is noncrossing. Put names $\bar{1}, \dots, \bar{n}$ on the midpoints of the $1, \dots, n$ (so that \bar{i} is the midpoint of the segment from i to $i+1$). The

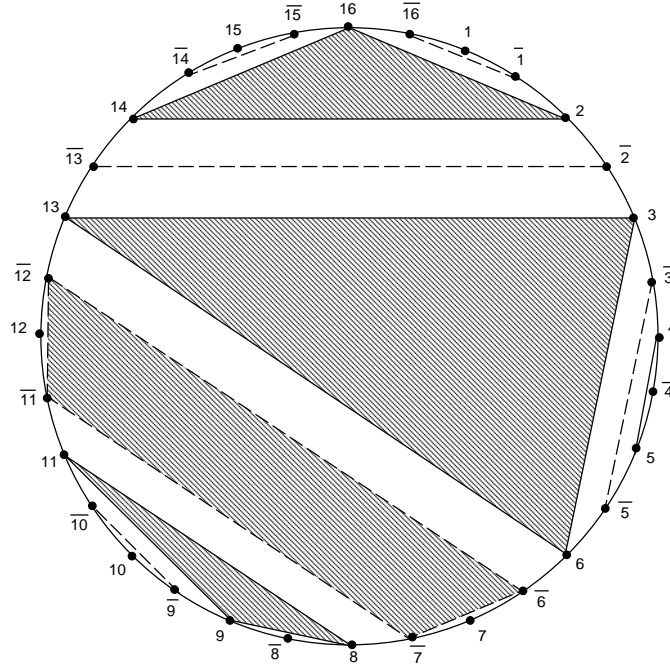


FIG 1. *The circular representation of a partition of $\{1, \dots, 16\}$.*

complement of the set $\cup_i H_i$ is again a union of disjoint convex sets \tilde{H}_i . All this is demonstrated in figure 1, where the \tilde{H}_i are the scrambled areas with dashed borders, the H_i are the scrambled areas with non-dashed borders. We will refer to this figure heavily during the proof of theorem 3.4. We can now define the Kreweras complementation map:

DEFINITION 2.3. *The Kreweras complement of π , denoted $K(\pi)$, is the partition on $\{\bar{1}, \dots, \bar{n}\}$ determined by*

$$i \sim j \text{ in } K(\pi) \iff \bar{i}, \bar{j} \text{ belong to the same convex set } \tilde{H}_k.$$

The connection between the R -transform and noncrossing partitions comes through the *moment-cumulant formula*, which relates the moments and the R -transform coefficients (also called cumulants) for the distribution of a random variable.

LEMMA 2.1. *Write the R -transform as a power series, $R_{\mu_a}(z) = \sum_n \alpha_n z^n$.*

Then

$$(2.12) \quad \phi(a^n) = \sum_{\pi=\{B_1, \dots, B_k\} \in NC(n)} \prod_{i=1}^k \alpha_{|B_i|},$$

This can be used as an alternative definition of the R -transform. We also need to define the multidimensional R -transform for the joint distribution of a sequence of random variables. Denote by $C\langle z_1, \dots, z_m \rangle$ the space of complex power series in m noncommuting variables z_i with vanishing constant term. These can be written in the form

$$\sum_{k \geq 1} \sum_{i_1, \dots, i_k} a_{i_1, \dots, i_k} z_{i_1} \cdots z_{i_k}.$$

In referring to the coefficients of a power series f on this form we will write

$$[\text{coef}(i_1, \dots, i_k)](f) = a_{i_1, \dots, i_k},$$

and if $\pi = \{B_1, \dots, B_m\}$,

$$\begin{aligned} [\text{coef}(i_1, \dots, i_k)|B_i](f) &= a_{(i_j)_{j \in B_i}} \\ [\text{coef}(i_1, \dots, i_m); \pi](f) &= \prod_i [\text{coef}(i_1, \dots, i_k)|B_i](f). \end{aligned}$$

For power series in one variable, the coefficients will also be written in the form $[\text{coef}_k](f)$. For n random variables a_1, \dots, a_n we define their joint moment series as the power series $M_{\mu_{a_1, \dots, a_n}} \in C\langle z_1, \dots, z_n \rangle$ such that

$$M_{\mu_{a_1, \dots, a_n}}(z_1, \dots, z_k) = \sum_{m \geq 1} \sum_{i_1, \dots, i_m} \phi(a_{i_1} \cdots a_{i_m}) z_{i_1} \cdots z_{i_m},$$

and we define their joint R -series as the unique power series $R_{\mu_{a_1, \dots, a_n}} \in C\langle z_1, \dots, z_n \rangle$ such that

$$(2.13) \quad \phi(a_{i_1} \cdots a_{i_m}) = \sum_{\pi \in NC(m)} [\text{coef}(i_1, \dots, i_m); \pi](R_{\mu_{a_1, \dots, a_n}}).$$

The result we will use connecting the joint R -series and freeness is the following:

LEMMA 2.2. $(\{a_1, \dots, a_n\}, \{b_1, \dots, b_m\})$ is a free family if and only if

$$\begin{aligned} &R_{\mu_{a_1, \dots, a_n, b_1, \dots, b_m}}(z_1, \dots, z_{n+m}) \\ &= R_{\mu_{a_1, \dots, a_n}}(z_1, \dots, z_n) + R_{\mu_{b_1, \dots, b_m}}(z_{n+1}, \dots, z_{n+m}). \end{aligned}$$

This lemma is often summarized by saying that the joint R -series of free random variables has no *mixed* terms. A special form of (2.13) and lemma 2.2 we will use is the following: If (a_1, a_2) is a free family, and a mixed term $a_{i_1} \cdots a_{i_m}$ is given, form the partition with two blocks $\sigma = \{\sigma_1, \sigma_2\}$, where $\sigma_k = \{j | a_{i_j} = k\}$. Then

$$(2.14) \quad \phi(a_{i_1} \cdots a_{i_m}) = \sum_{\pi \leq \sigma \in NC(m)} [\text{coef}(i_1, \dots, i_m); \pi](R_{\mu_{a_1, a_2}}).$$

Our combinatorial connection with multiplicative free convolution can be made complete with the help of the following definition [10], [11]:

DEFINITION 2.4. *Given two power series f and g , their boxed convolution $f \boxtimes g$ is defined by*

$$(2.15) \quad \begin{aligned} & [\text{coef}(i_1, \dots, i_m)](f \boxtimes g) \\ &= \sum_{\pi \in NC(m)} [\text{coef}(i_1, \dots, i_m); \pi](f) [\text{coef}(i_1, \dots, i_m); K(\pi)](g). \end{aligned}$$

Boxed convolution is commutative only on power series in one variable [13]. It satisfies the associative law, but not the distributive law. It does not satisfy linearity properties w.r.t. scalar multiplication. However, the following holds and will be useful to us:

$$[\text{coef}_n]((cf) \boxtimes (cg)) = [\text{coef}_n](c^{n+1}(f \boxtimes g))$$

and

$$[\text{coef}_n](f \boxtimes (cId)) = [\text{coef}_n](c^n f).$$

Here we used the shorthand notation $c^n f$ for the power series defined by $[\text{coef}_n](c^n f) = c^n [\text{coef}_n](f)$. The first statement is easily proved using the fact that $|\pi| + |K(\pi)| = n + 1$ for any $\pi \in NC(n)$ [13]. The second statement is trivial. The following result holds for multiplicative free convolution [13]:

LEMMA 2.3. *If $(\{a_1, \dots, a_n\}, \{b_1, \dots, b_n\})$ is a free family, then*

$$(2.16) \quad R_{\mu_{a_1 b_1, \dots, a_n b_n}} = R_{\mu_{a_1, \dots, a_n}} \boxtimes R_{\mu_{b_1, \dots, b_n}}$$

One can also define additive and multiplicative free deconvolution in most cases, i.e. finding μ_2 in (2.3) when $\mu_1 \boxtimes \mu_2$ are known.

DEFINITION 2.5. *Given probability measures μ and μ_2 . When there is a unique probability measure μ_1 such that $\mu = \mu_1 \boxtimes \mu_2$, we will denote $\mu_1 = \mu \boxtimes \mu_2$. We say that μ_1 is the multiplicative free deconvolution of μ with μ_2 .*

We can define additive free deconvolution similarly. Note that free deconvolution is defined only for a subset of all probability measures, since measures exist which can't be expressed on the forms $\mu_1 \boxplus \mu_2$ or $\mu_1 \boxtimes \mu_2$. Deconvolution can, however, also be viewed as a formal operation on a sequence of moments. Viewed as such, multiplicative free deconvolution is well-defined when we have non-vanishing first moments. This can be seen from the combinatorial description of multiplicative free convolution (2.15). In light of (2.16), it is obvious from (2.15) that the cumulants in R_{μ_2} can be calculated recursively from those of R_{μ_1} and $R_{\mu_1 \boxtimes \mu_2}$, when the first coefficient of R_{μ_1} (which equals the first moment) is known. Since the main theorem relates to the moments of the involved measures (it is a statement on convergence in distribution), we will in the following view deconvolution in terms of the moments only.

A form of (2.16) which will be useful to us is for the case $n = 1$. If we write $\mu_a = (\mu_a \boxtimes \mu_c) \boxtimes \mu_c$, we get

$$(2.17) \quad R_{\mu_a \boxtimes \mu_c} = R_{\mu_a} \boxtimes R_{\mu_c}^{-1}.$$

The facts we will use concerning boxed convolution are the following, relating moment series, R -series of general random variables (in particular projections and free Poisson random variables), the *Zeta* series, which is defined as

$$Zeta(z_1, \dots, z_n) = \sum_k \sum_{i_1, \dots, i_k} z_{i_1} \cdots z_{i_k},$$

the *Moeb* series (which is the inverse of *Zeta* under composition with \boxtimes) and *Id* (which is the unit under composition with \boxtimes):

$$(2.18) \quad \begin{aligned} M_\mu &= R_\mu \boxtimes Zeta & \text{and} & & R_\mu &= M_\mu \boxtimes Moeb \\ M_{\mu_p} &= cZeta & \text{and} & & R_{\mu_c} &= c^{n-1}Zeta. \end{aligned}$$

Here p is a projection with $\phi(p) = c$. Our definition of μ_c differs from that of [7], for purposes of compatibility with [3] [14]. Consequently, the expressions for the R -transforms are different. In the terminology of [7], the R -series would be $cZeta$. The following definition [12] will also be in use:

DEFINITION 2.6. *A pair (a, b) of noncommutative random variables is called an R -diagonal pair if its R -series is of the form*

$$(2.19) \quad R_{\mu_{a,b}}(z_1, z_2) = \sum_{n=1}^{\infty} \alpha_n ((z_1 z_2)^n + (z_2 z_1)^n).$$

An element a will be said to be an R -diagonal element if (a, a^) is an R -diagonal pair. The one-variable series $\sum_{n=1}^{\infty} \alpha_n z^n$ will be called the determining series of the R -diagonal pair (a, b) .*

We will use the fact that if a is an R -diagonal element, its determining series can be written as $R_{\mu_{aa^*}} \boxtimes Moeb$ [7]. Two important R -diagonal elements are

1. the *Haar unitary*, which can be defined as a unitary u satisfying $\phi(u^n) = 0$ for all $n \in \mathbb{Z} \neq 0$, and
2. the *circular element*, which can be defined as an element s whose $*$ -distribution μ_{s,s^*} satisfies $R_{\mu_{s,s^*}}(z_1, z_2) = z_1 z_2 + z_2 z_1$.

The concept of R -diagonality was in fact invented in search of a common approach for Haar unitaries and circular elements [12]. Haar unitaries are very important in asymptotic random matrix results. In W^* -probability spaces, when the isometric part of an R -diagonal element has kernel equal to zero, the isometric part is actually a Haar unitary.

2.1. Implementation of free convolution. While free convolution has an abstract definition, the combinatorial description given in this section can actually be used to obtain an efficient implementation. In many practical cases, free convolution with μ_c is what we are interested in. Such free convolution is simplified through the following result.

THEOREM 2.1.

$$(2.20) \quad (cM_\mu) \boxtimes Zeta = c(M_{\mu \boxtimes \mu_c}).$$

PROOF. To see this, start by combining (2.16) with (2.18) to get

$$R_{\mu \boxtimes \mu_c} = R_\mu \boxtimes R_{\mu_c} = R_\mu \boxtimes (c^{m-1} Zeta).$$

After convolving both sides with $Zeta$, we get

$$(2.21) \quad M_{\mu \boxtimes \mu_c} = M_\mu \boxtimes (c^{m-1} Zeta).$$

To prove (2.20), rewrite the left hand side as

$$\sum_{\pi \in NC(m)} c^{|\pi|} [coef_m; \pi] M_\mu.$$

Since $|\pi| + |K(\pi)| = m + 1$, this equals

$$\begin{aligned} & c \sum_{\pi \in NC(m)} [coef_m; \pi] (M_\mu) c^{m-|K(\pi)|} \\ &= c \sum_{\pi \in NC(m)} [coef_m; \pi] (M_\mu) c^{m-|K(\pi)|} [coef_m; K(\pi)] (Zeta) \\ &= c \sum_{\pi \in NC(m)} [coef_m; \pi] (M_\mu) c^m [coef_m; K(\pi)] (c^{-1} Zeta) \\ &= c \sum_{\pi \in NC(m)} [coef_m; \pi] (M_\mu) [coef_m; K(\pi)] (c^{m-1} Zeta) \\ &= c (M_\mu \boxtimes (c^{m-1} Zeta)), \end{aligned}$$

substituting (2.21) proves the claim. \square

In summary, if we need to compute the moments of $\mu \boxtimes \mu_c$, one can first compute the moment series cM_μ , then use this to compute the left hand side of (2.20). According to (2.21) and (2.20), the moment series of $\mu \boxtimes \mu_c$ can then be computed from this with an additional scaling with $\frac{1}{c}$.

In other words, convolving with μ_c is equivalent to convolving with μ_1 (with additional scalings of power series taken into account), since $R_{\mu_1} = Zeta$. It turns out that boxed convolution with *Zeta* is easy to compute, as the following result shows. The result is stated in terms of the moment-cumulant formula, since the relation between cumulants and moments are given by boxed convolution with *Zeta*.

THEOREM 2.2.

$$(2.22) \quad [coef_m](M_\mu) = \sum_{k=1}^m [coef_k](R_\mu)[coef_{m-k}](1 + M_\mu)^k.$$

PROOF. For each $\pi \in NC(m)$, fix the block $B_1 = \{b_{11}, \dots, b_{1k}\}$ in π containing 1, and let $NC(m, B_1)$ be the set of all noncrossing partitions which contain B_1 as a block. Rewrite the definition of boxed convolution (2.15) to

$$(2.23) \quad \begin{aligned} [coef_m](M_\mu) &= \sum_{B_1} \sum_{\pi \in NC(m, B_1)} [coef_m; \pi](R_\mu)[coef_m; K(\pi)](Zeta) \\ &= \sum_{B_1} \sum_{\pi \in NC(m, B_1)} [coef_m; \pi](R_\mu). \end{aligned}$$

Blocks in $\pi \in NC(m, B_1)$ other than B_1 must be entirely contained in one of $\{b_{11} + 1, \dots, b_{12} - 1\}, \dots, \{b_{1k} + 1, \dots, b_{11} - 1\}$. This means that the inner summand in (2.23) can be rewritten to

$$(2.24) \quad [coef_k](R_\mu) \prod_{i=1}^k \left(\sum_{\pi \in NC(b_{1(i+1)} - b_{1i} - 1)} [coef_m; \pi](R_\mu) \right).$$

From the moment-cumulant formula it is seen that each sum here is simply a moment, so we can rewrite to

$$[coef_k](R_\mu) \prod_{i=1}^k [coef_{b_{1(i+1)} - b_{1i} - 1}](1 + M_\mu),$$

where the summand 1 in $1 + M_\mu$ accounts for elements i in (2.24) with $b_{1(i+1)} = b_{1i} + 1$ (i.e. consecutive elements in a block). All in all, (2.23) can be rewritten to

$$(2.25) \quad \sum_k \sum_{\substack{B_1 \\ |B_1|=k}} [coef_k](R_\mu) \prod_{i=1}^k \left([coef_{b_{1(i+1)} - b_{1i} - 1}](1 + M_\mu) \right)$$

Write $a_i = b_{1(i+1)} - b_{1i} - 1$, and note that

$$\sum_{i=1}^k a_i = \sum_{i=1}^k (b_{1(i+1)} - b_{1i} - 1) = m - k.$$

The a_i are in one-to-one correspondence with all candidates for B_1 , so that we can rewrite (2.25) to

$$\sum_k [\text{coef}_k](R_\mu) \sum_{\substack{a_1, \dots, a_k \\ \sum a_i = m - k}} \prod_{i=1}^k ([\text{coef}_{a_i}](1 + M_\mu)).$$

The inner sum here is easily recognized as coefficient $m - k$ in the power series $(1 + M_\mu)^k$ (one factor for each a_i). Putting things together we get (2.22). \square

In (2.22) we see that there is no reference to noncrossing partitions. (2.22) can be used easily in calculating moments recursively from cumulants. The coefficients in the power series $(1 + M_\mu)^k$ can be computed in terms of k -fold (classical) convolution. This is done in [18], where many multiplicative free convolutions are computed based on (2.22). The actual implementation of (2.22) used in [18] is contained in [16].

Free convolution as introduced here is just defined for compactly supported probability measures.

3. Proof of theorem 1.1. In what follows we first sketch the proof of theorem 1.1. After this follows proofs for theorems needed in the proof.

First we prove the following variant of lemma 4.3.2 in [7], which can be used together with the Borel-Cantelli lemma to prove almost sure convergence. It is slightly more general in the sense that boundedness in the operator norm $\|\cdot\|$ is not assumed, only boundedness in $\|\cdot\|_p$, for $p \geq 1$. This weaker boundedness assumption is needed since uniformly norm-bounded matrices are not sufficient to approximate all compactly supported probability measures almost surely. Recall that an $n \times n$ unitary random matrix is called *standard unitary* if its distribution equals the Haar probability measure on $\mathcal{U}(n)$.

THEOREM 3.1. *Let $U(s, n)_{s \in S}$ be an independent family of $n \times n$ standard unitary random matrices. Let $s_1, \dots, s_l \in S$, $m_1, \dots, m_l \in \mathbb{Z} \setminus \{0\}$, and let $R_p \geq 0$, $p \geq 1$ be constants. Then*

$$(3.1) \quad E \left(\left| \text{tr}_n (U(s_1, n)^{m_1} D_1(n) U(s_2, n)^{m_2} D_2(n) \cdots U(s_l, n)^{m_l} D_l(n)) \right|^2 \right)$$

is $O(n^{-2})$ as $n \rightarrow \infty$ uniformly for the choice of any $D_r(n) \in M_n(C)$ ($1 \leq r \leq l$) such that for $1 \leq r \leq l$ either

$$\operatorname{tr}_n(D_r(n)) = 0 \text{ and } \|D_r(n)\|_p \leq R_p \text{ (} n \in N \text{)}$$

or

$$D_r(n) = I_n \text{ (} n \in N \text{) and } s_r \neq s_{r+1} \text{ (with } s_{l+1} = s_1 \text{)}.$$

Also, for a given l , there exists a p_l such that the same statement holds as long as the $\|\cdot\|_p$ -norm bounds are satisfied for $p \leq p_l$ only.

The proof is in section 3.1. It somewhat simplifies the proof of lemma 4.3.2 in [7], and can also be used to simplify the proof of theorem 4.3.5 in [7]. As in [7], theorem 3.1 is sufficient to prove asymptotic freeness almost everywhere for the family

$$(\{U(s, n), U(s, n)^*\})_{s \in S}, \{D(t, n), D(t, n)^* : t \in T\}$$

when the $D_r(n)$ is known to have a limit distribution. It will also be useful to us that theorem 3.1 gives us bounds also in cases where the $D_r(n)$ do *not* converge to a limit. The $D_r(n)$ model in our case concerns $\frac{1}{\sqrt{n}}R_n$ random matrices, for which it is not known whether an almost sure limit exists (only that $\Gamma_n = \frac{1}{N}R_nR_n^*$ has an almost sure limit). Also, theorem 3.1 gives us grounds for proving that only the lower mixed moments converge to zero. It can be applied to cases where only the lower $\|\cdot\|_p$ -norms are known to be bounded, in which only lower mixed moments can be bounded.

What we really want is to use random matrices R_n independent from the U_n instead of the deterministic matrices $D_r(n)$. This is addressed by the following theorem. We restrict to the case of one standard unitary random matrix.

THEOREM 3.2. *Let U_n be $n \times n$ standard unitary random matrices, and let R_n be random matrices independent from U_n , such that $R_nR_n^*$ converges in distribution almost surely to a compactly supported probability measure ρ . Then*

$$(3.2) \quad |\operatorname{tr}_n(U_n^{m_1}P_1(R_n)U_n^{m_2}P_2(R_n)\cdots U_n^{m_l}P_l(R_n))| \rightarrow 0 \text{ a.s.}$$

uniformly for any choice of polynomials P_1, \dots, P_l such that $\operatorname{tr}_n(P_i(R_n)) = 0$ for all $1 \leq i \leq l$.

The proof is in section 3.2. As for theorem 3.1, theorem 3.2 is sufficient to prove asymptotic freeness almost everywhere for the family (U_n, R_n) when the R_n are additionally known to have a limit distribution.

The proof is split in two: First (3.2) is shown for random matrices satisfying bounds of the form $\|R_n\|_p \leq R_p$ ($p \geq 1$). The proof in this case uses theorem 3.1 and is quite short. The more general case of compactly supported probability measures is proved with an approximation argument.

The next step is to pass from standard unitary random matrices U_n to standard Gaussian random matrices X_n . Note that it could be possible to skip starting with standard unitary random matrices altogether, by building directly on results for almost sure convergence of Gaussian random matrices like those in [15]. We have chosen the approach with standard unitary random matrices for compatibility with [7]. We will prove the following:

THEOREM 3.3. *Let X_n be $n \times n$ standard Gaussian random matrices, and let R_n be random matrices independent from X_n , such that $R_n R_n^*$ converges in distribution almost surely to a compactly supported probability measure ρ . Then*

$$|tr_n(Q_1(X_n)P_1(R_n)Q_2(X_n)P_2(R_n) \cdots Q_l(X_n)P_l(R_n))| \rightarrow 0 \text{ a.s.}$$

uniformly for any choice of polynomials $Q_1, \dots, Q_l, P_1, \dots, P_l$ such that

$$tr_n(Q_i(R_n)) = 0 \text{ and } tr_n(P_i(R_n)) = 0$$

for all $1 \leq i \leq l$.

The proof is quite short, and also presented in section 3.2. Note that the approximation argument used in the proof of theorem 4.3.5 in [7] does not work in this case. As for theorem 3.2, theorem 3.3 is enough to prove asymptotic freeness almost everywhere when the R_n are additionally known to have a limit distribution. Just as theorem 3.1 gives bounds for mixed moments also in cases where the deterministic matrices do not converge in distribution, theorem 3.2 and its counterpart for Gaussian random matrices can be used to bound mixed moments in cases where it is only known that the R_n matrices satisfy $\|\cdot\|_p$ -norm bounds.

To finish the proof we will model our situation through the following theorem, which is stated independently of a random matrix setting.

THEOREM 3.4. *Suppose that a and $\{p, b\}$ are $*$ -free, with a R -diagonal and p a projection with $\phi(p) = c$. In the reduced probability space $(pAp, \phi(p)^{-1}\phi)$, $\mu_{p(a+b)(a+b)^*p}$ is uniquely identified by μ_{paa^*p} and μ_{pbb^*p} through the equation*

$$(3.3) \quad \mu_{p(a+b)(a+b)^*p} \boxtimes \mu_c = (\mu_{paa^*p} \boxtimes \mu_c) \boxplus (\mu_{pbb^*p} \boxtimes \mu_c)$$

*In particular, $\mu_{p(a+b)(a+b)^*p}$ has no dependence on mixed moments of a and b .*

This will be proved in section 3.3. Note that there is no assumption on freeness between p and b . The case $c = 1$ is particularly interesting, and corresponds to $p = I$. In this case, $\mu_{(a+b)(a+b)^*}$ is uniquely identified by μ_{aa^*} and μ_{bb^*} , and (3.3) is simply

$$(3.4) \quad \mu_{(a+b)(a+b)^*} \boxtimes \mu_1 = (\mu_{aa^*} \boxtimes \mu_1) \boxplus (\mu_{bb^*} \boxtimes \mu_1).$$

This equation has an interpretation in terms of square random matrices.

Due to (3.3), R -diagonality relieves us from dependencies of many mixed moments, so that some cancellation phenomenon must occur. This also happens in other cases. If a and b are free, [13] expresses the distribution of the *free commutator*, i.e.

$$(3.5) \quad R_{\mu_{i(ab-ba)}}(z) = 2 \left(R_{\mu_a}^{\text{even}} \boxtimes R_{\mu_b}^{\text{even}} \boxtimes \text{Zeta} \right) (z^2),$$

where $R^{\text{even}}(z) = \sum_{n=1}^{\infty} \alpha_{2n} z^n$ whenever $R(z) = \sum_{n=1}^{\infty} \alpha_n z^n$. (3.5) holds also when a and b are not R -diagonal. (3.5) also expresses a connection with multiplicative free convolution with μ_1 , since boxed convolution with the *Zeta*-series is involved.

Theorem 3.4 has a more general flavour than theorem 4.1, since the limits $\frac{1}{N} X_n X_n^*$ from theorem 4.1 do not include all R -diagonal pairs. The following limiting version of theorem 3.4 will be useful in finishing the proof of theorem 4.1:

THEOREM 3.5. *Let the random variables $\{a_n, b_n, p_n\} \in (\mathcal{A}_n, \phi_n)$, $\{a, p\} \in (\mathcal{A}, \phi)$ be given, where a is R -diagonal and p, p_n are projections with $\phi(p) = \phi_n(p_n) = c$. Form the random variable paa^*p in $(p\mathcal{A}p, \phi(p)^{-1}\phi)$, and the random variables $p_n b_n b_n^* p_n$ and $p_n(a_n + b_n)(a_n + b_n)^* p_n$ in $(p_n \mathcal{A}_n p_n, \phi(p_n)^{-1}\phi_n)$. If*

$$\mu_{a_n, a_n^*} \rightarrow \mu_{a, a^*}, \quad \mu_{p_n} \rightarrow \mu_p,$$

and

$$\mu_{p_n b_n b_n^* p_n} \rightarrow \mu,$$

in distribution, moments are uniformly bounded in n , and mixed moments of $(a_n, \{p_n, b_n\})$ go to 0, then $\mu_{p_n(a_n + b_n)(a_n + b_n)^* p_n}$ converges in distribution and the limit is uniquely identified by the equation

$$(3.6) \quad \lim_{n \rightarrow \infty} \mu_{p_n(a_n + b_n)(a_n + b_n)^* p_n} \boxtimes \mu_c = (\mu_{paa^*p} \boxtimes \mu_c) \boxplus (\mu \boxtimes \mu_c)$$

PROOF. The limiting moments of $\mu_{p_n(a_n + b_n)(a_n + b_n)^* p_n} \boxtimes \mu_c$ do not change if we "zero out" the mentioned mixed moments (i.e. that we assume freeness

of $(a_n, \{p_n, b_n\})$, due to the assumption on their vanishing and of uniform boundedness on moments. It is also easily seen that the limiting moments do not change if we change the distribution of a_n to $\mu_{a_n} = \mu_a$ for all n . But then

$$\begin{aligned} \lim_{n \rightarrow \infty} \mu_{p_n(a_n+b_n)(a_n+b_n)^*p_n} \boxtimes \mu_c &= \lim_{n \rightarrow \infty} \mu_{p_n(a+b_n)(a+b_n)^*p_n} \boxtimes \mu_c \\ &= \lim_{n \rightarrow \infty} (\mu_{p_n a a^* p_n} \boxtimes \mu_c) \boxplus (\mu_{p_n b_n b_n^* p_n} \boxtimes \mu_c) = (\mu_{p a a^* p} \boxtimes \mu_c) \boxplus (\mu \boxtimes \mu_c) \end{aligned}$$

where we used theorem 3.4, so that (3.6) holds. \square

The rest of the proof of theorem 4.1 now goes as follows: The rectangular random matrices R_n can be viewed as the $N \times N$ random matrices $p_n S_n$, where the projection p_n is a diagonal constant matrix, with the fraction of 1's on the diagonal equal to c , and S_n is an extension of the $n \times N$ matrix R_n to an $N \times N$ -matrix, obtained by adding zeros. Similarly, the random matrices X_n can be viewed as the $N \times N$ -matrices $p_n Y_n$, where Y_n is an extension of the $n \times N$ matrix X_n to an $N \times N$ -matrix, obtained by adding more independent standard Gaussian entries.

Since $\frac{1}{N} S_n S_n^*$ almost surely converges to a compactly supported probability measure, $(\frac{1}{\sqrt{n}} Y_n, \frac{1}{\sqrt{n}} S_n)$ satisfies the requirements of theorem 3.3. Thus, mixed moments of $\frac{1}{\sqrt{n}} X_n$ and $\frac{1}{\sqrt{n}} S_n$ go to zero almost surely. It is also seen that $\frac{1}{\sqrt{n}} S_n$ has its moments bounded as $n \rightarrow \infty$ almost surely. It is well known [7] that $\frac{1}{\sqrt{N}} Y_n$ converges in distribution almost surely to the circular law, which is R -diagonal.

Thus, all assumptions of theorem 3.5 are satisfied for $a_n = \frac{1}{\sqrt{N}} Y_n$, $b_n = \frac{1}{\sqrt{N}} S_n$ and p_n , almost surely. Thus, almost surely,

$$\begin{aligned} \lim_{n \rightarrow \infty} \mu_{p_n \frac{1}{N} (S_n + \sigma Y_n) (S_n + \sigma Y_n)^* p_n} \boxtimes \mu_c &= \left(\mu_{p \sigma^2 \frac{1}{N} Y Y^* p} \boxtimes \mu_c \right) \boxplus (\mu_\Gamma \boxtimes \mu_c) \\ &= (\mu_{\sigma^2 I} \boxtimes \mu_c \boxtimes \mu_c) \boxplus (\mu_\Gamma \boxtimes \mu_c) = \mu_{\sigma^2 I} \boxplus (\mu_\Gamma \boxtimes \mu_c), \end{aligned}$$

or

$$\lim_{n \rightarrow \infty} \mu_{\frac{1}{N} (R_n + \sigma X_n) (R_n + \sigma X_n)^*} \boxtimes \mu_c = \mu_{\sigma^2 I} \boxplus (\mu_\Gamma \boxtimes \mu_c),$$

which is the statement of theorem 4.1.

The proof as sketched here assumes $c \leq 1$. An explanation for how the proof goes for $c > 1$ is given succeeding the proof of theorem 3.4.

3.1. *The proof of theorem 3.1.* The proof will use the (generalized) Hölder inequality:

LEMMA 3.1. *For matrices A_1, \dots, A_k , the following holds:*

$$\|A_1 \cdots A_k\|_p \leq \|A_1\|_{p_1} \cdots \|A_k\|_{p_k} \text{ when } \sum_{i=1}^k \frac{1}{p_i} = \frac{1}{p}.$$

In the proof of lemma 4.3.2 in [7], (3.1) is written as

$$(3.7) \quad \left(\frac{1}{n}\right)^2 \sum_{i_1, \dots, i_{2k}=1}^n \sum_{j_{k(1)}, \dots, j_{k(l)}, j_{k(l+1)+1}, \dots, j_{k(2l)+1}}^n \left(\prod_{r=1}^l d_{j_{k(r)} i_{k(r)+1}}(t_r, n) \right) \\ \times \left(\prod_{r=l+1}^{2l} \bar{d}_{i_{k(r)} j_{k(r)+1}}(t_r, n) \right) E \left(\prod_{h=1}^{2k} u_{i_h j_h}(s(h), \varepsilon(h), n) \right),$$

where for $1 \leq r \leq l$,

$$k(r) = |m_1| + \cdots + |m_r|,$$

$$k = k(l), \quad k(l+r) = k + k(r), \quad t_{l+r} = t_r.$$

Moreover, for h such that $k(r-1) + 1 \leq h \leq k(r)$,

$$s(h) = s_r, \quad \varepsilon(h) = \begin{cases} 1 & \text{if } m_r > 0 \\ -1 & \text{if } m_r < 0 \end{cases}$$

Here $s(h+k) = s(h)$ and $\varepsilon(k+h) = -\varepsilon(h)$ for $1 \leq h \leq k$, and

$$u_{ij}(s, \varepsilon, n) = \begin{cases} U_{ij}(s, n) & \text{if } \varepsilon = 1 \\ \bar{U}_{ji}(s, n) & \text{if } \varepsilon = -1 \end{cases}$$

Since (3.7) is a matrix product written out, the following must hold:

$$(3.8) \quad \begin{cases} j_h = i_{h+1} & \text{for } h \in \{1, \dots, k\} \setminus \{k(1), \dots, k(l)\}, \\ i_h = j_{h+1} & \text{for } h \in \{k+1, \dots, 2k\} \setminus \{k(l+1), \dots, k(2l)\} \end{cases}$$

Also, due to the vanishing of many mixed moments of entries in standard unitary random matrices (lemma 4.2.2 in [7]), two pair partitions \mathcal{U} and \mathcal{V} can be chosen so that if $\{h, h'\} \in \mathcal{U}$ then

$$(3.9) \quad s(h) = s(h'), \quad \varepsilon(h) = 1, \quad \varepsilon(h') = -1, \quad i_h = j_{h'} (= i_{h'} + 1),$$

and if $\{h, h'\} \in \mathcal{V}$ then

$$(3.10) \quad s(h) = s(h'), \varepsilon(h) = -1, \varepsilon(h') = 1, i_h = j_{h'} (= i_{h'} + 1)$$

These two pair partitions and (3.8) cause many equalities among the i_1, \dots, i_{2k} , and define the equivalence relation $\mathcal{R}(\mathcal{U}, \mathcal{V})$ on $\{1, \dots, 2k\}$ so that $i_h = i_{h'}$ whenever h and h' are in the same equivalence class of $\mathcal{R}(\mathcal{U}, \mathcal{V})$. We let k_0 denote the number of equivalence classes of $\mathcal{R}(\mathcal{U}, \mathcal{V})$, and let $h(1), \dots, h(k_0)$ be representatives from the equivalence classes.

Recall the expressions

$$(3.11) \quad C_n(\iota_1, \dots, \iota_{k_0}) = \left(\prod_{r=1}^l d_{j_{k(r)} i_{k(r)+1}}(t_r, n) \right) \left(\prod_{r=l+1}^{2l} \bar{d}_{i_{k(r)} j_{k(r)+1}}(t_r, n) \right),$$

$$(3.12) \quad Q_n(\iota_1, \dots, \iota_{k_0}) = E \left(\prod_{h=1}^{2k} u_{i_h j_h}(s(h), \varepsilon(h), n) \right)$$

from [7], where $(\iota_1, \dots, \iota_{k_0})$ in (3.11) are defined as $(i_{h(1)}, \dots, i_{h(k_0)})$ (i.e. representatives of the equivalence classes), and j_1, \dots, j_{2k} are determined subject to \mathcal{U}, \mathcal{V} . (4.3.6) of [7] says that it is enough to prove that for any partition \mathcal{W} of $\{1, \dots, k_0\}$ we have that

$$(3.13) \quad \sum_{(\iota_1, \dots, \iota_{k_0}) : \mathcal{W}} C_n(\iota_1, \dots, \iota_{k_0}) Q_n(\iota_1, \dots, \iota_{k_0}) = O(1) \text{ as } n \rightarrow \infty,$$

where the summation is over $(\iota_1, \dots, \iota_{k_0})$ such that $\iota_p = \iota_q$ if and only if p and q are in the same block of \mathcal{W} . For a given \mathcal{W} , it is known that $|Q_n(\iota_1, \dots, \iota_{k_0})| = O(n^{-k})$ uniformly for i_k, j_k as $n \rightarrow \infty$, and that it has the same value for all $(\iota_1, \dots, \iota_{k_0})$ taking part in the sum (3.13). So, from (3.14) we deduce that it is enough to show that for any choice of $\mathcal{U}, \mathcal{V}, \mathcal{W}$ (there is a finite number of such choices),

$$(3.14) \quad \sum_{(\iota_1, \dots, \iota_{k_0}) : \mathcal{W}} C_n(\iota_1, \dots, \iota_{k_0})$$

is $O(n^k)$. In [7] this is proved using the fact that (3.11) are bounded uniformly. This is not true in our case since only uniform boundedness in $\|\cdot\|_p$ is assumed. Instead, we will group sums of terms into matrix multiplication units, and use the Hölder inequality together with the $\|\cdot\|_p$ -norm bounds. Instead of the terms in (3.14), where the sum is over

$$(\iota_1, \dots, \iota_{k_0}) : \iota_p = \iota_q \text{ if and only if } p \text{ and } q \text{ are in the same block of } \mathcal{W},$$

it will be better for us to sum over

$$(\iota_1, \dots, \iota_{k_0}): \iota_p = \iota_q \text{ if } p \text{ and } q \text{ are in the same block of } \mathcal{W}.$$

The latter set is more compatible with indices in multiplications of many matrices. This second set is larger than the first, and can be written as

$$(3.15) \quad \sum_{\mathcal{W}' \geq \mathcal{W}} \sum_{(\iota_1, \dots, \iota_{k_0}): \mathcal{W}' } C_n(\iota_1, \dots, \iota_{k_0})$$

It is obvious that (3.14) can be written

$$(3.16) \quad \sum_{\mathcal{W}' \geq \mathcal{W}} a_{\mathcal{W}'} \sum_{\mathcal{W}'' \geq \mathcal{W}'} \sum_{(\iota_1, \dots, \iota_{k_0}): \mathcal{W}''} C_n(\iota_1, \dots, \iota_{k_0}),$$

where $a_{\mathcal{W}'}$ are integer constants which can easily be calculated (proving (3.16) boils down to splitting all values of ι_1, ι_2 into those where $\iota_1 = \iota_2$, and those where $\iota_1 \neq \iota_2$. This is done recursively and for all ι_i to yield (3.16)). Since there is a finite number of elements in the two outer sums in (3.16), to prove that (3.14) is $O(n^k)$ it is enough to show that (3.15) is $O(n^k)$ for any choice of \mathcal{W} .

Let l_0 denote the number of equivalence classes in $\mathcal{R}(\mathcal{U}, \mathcal{V})$ with only one entry, and let $h(1), \dots, h(l_0)$ be the corresponding representatives. According to [7], equivalence classes with only one entry give rise to factors of the form $\tilde{d}_{\iota_i \iota_i}$ in (3.11), where \tilde{d} is either $d_{\iota_i \iota_i}(t_r, n)$ or $\bar{d}_{\iota_i \iota_i}(t_r, n)$ for some r . Equivalence classes with only one entry thus leads (through summation over one ι_i appearing in just one factor) to factors in (3.15) which are (non-normalized) traces of the $D(t_r, n)$. These are zero, so we can assume that $l_0 = 0$ when we attempt to bound (3.15). Had we used the sum (3.14) instead of (3.15), we would not obtain zero.

So we assume that there are no singleton equivalence classes, i.e. $k_0 \leq k$. Let K_0 be the number of equivalence classes actually appearing in (3.11) (this is a function of \mathcal{U} and \mathcal{V}). we have that $K_0 \leq k_0$, but equality does not necessarily hold. We will use matrix units E_{ij} (i.e. $E_{ij}(i, j) = \delta_{ij}$). By placing matrix units F_i with indices from $\iota_1, \dots, \iota_{k_0}$ in between the terms in (3.11), (3.15) can be written as

$$(3.17) \quad \sum_{\mathcal{W}' \geq \mathcal{W}} \sum_{(\iota_1, \dots, \iota_{k_0}): \mathcal{W}' } ntr_n \left(\prod_{i=1}^{2l} (F_i D_i) \right),$$

where D_i are matrices from $D(t_r, n)$ or one of their transposes/conjugates. Since $\|E_{ij}\|_p = n^{-\frac{1}{p}}$ and the number of possible choices of matrix units is

n^{K_0} , lemma 3.1 implies that (3.17) is bounded by

$$(3.18) \quad n^{K_0+1} n^{-\frac{1}{2}} \prod_{i=1}^{2l} \|D_i\|_{4l} = n^{K_0+\frac{1}{2}} \prod_{i=1}^{2l} \|D_i\|_{4l}.$$

Since $K_0 \leq k$, this is $O(n^k)$ except possibly in the case when $K_0 = k$, i.e. when all equivalence classes have exactly two elements.

So, for the rest of the proof, we assume that all equivalence classes have exactly two elements. Note that the number of times an equivalence class appears as an i is equal to the number of the times the same class appears as a j in (3.11). This is obvious from the way the equivalence relation is defined (3.8), (3.9), (3.10) in order to avoid a zero value in (3.12). This means that we can take the first of the K_0 equivalence classes appearing in (3.11), and rearrange the terms in (3.11) so that the equivalence class appear in alternating order as an i and as a j . (3.17) can thus be rewritten to

$$(3.19) \quad \sum_{\mathcal{W}' \geq \mathcal{W}(\iota_1, \dots, \iota_{k_0})} \sum_{\mathcal{W}'} n \operatorname{tr}_n (F_1 D_1 G_1 D_2 F_2 D_3 G_2 D_4 F_3),$$

where D_1, D_2, D_3, D_4 are the matrices where the first equivalence class appear as an i or a j , and in alternating order. Also, $F_1 = F_2 = F_3 = E_{rr}$ are matrix units, r is a given number, and the G_i are products man of the matrices D_i in (3.17). (3.19) can also be written

$$(3.20) \quad \sum_{\mathcal{W}' \geq \mathcal{W}(\iota_1, \dots, \iota_{k_0})} \sum_{\mathcal{W}'} n \operatorname{tr}_n (\operatorname{diag}(D_1 G_1 D_2) \operatorname{diag}(D_3 G_2 D_4)),$$

where $\operatorname{diag}(A)$ stands for the diagonal of the matrix A . Similarly to the calculation of the bound (3.18), (3.20) is seen to be bounded by

$$(3.21) \quad n^{K_0} \|\operatorname{diag}(D_1 G_1 D_2)\|_2 \|\operatorname{diag}(D_3 G_2 D_4)\|_2.$$

Note that $\|\operatorname{diag}(A)\|_2 \leq \|A\|_2$, since

$$\|\operatorname{diag}(A)\|_2^2 = \frac{1}{n} \sum_i |a_{ii}|^2 \leq \frac{1}{n} \sum_{i,j} |a_{ij}|^2 = \operatorname{tr}_n(A^* A) = \|A\|_2^2,$$

where $A = (a_{ij})_{i,j}$. This means that (3.21) is bounded by

$$n^{K_0} \|D_1 G_1 D_2\|_2 \|D_3 G_2 D_4\|_2,$$

which is $O(n^{K_0})$ and hence $O(n^k)$ since all D_i are bounded in p -norm, and the only other factors are matrix units, which have p -norm $n^{-\frac{1}{p}}$.

That there exists a p_l for a given l as in the last statement of the theorem is obvious from the proof and the way the Hölder inequality was used. This completes the proof.

3.2. *The proofs of theorem 3.2 and 3.3.* First assume that R_n satisfies $\|R_n\|_p \leq R_p$ ($p \geq 1$) almost surely for some constants R_p . $P_i(R_n)$ satisfies similar $\|\cdot\|_p$ -norm bounds due to lemma 3.1. Call the underlying probability space Ω . Denote by $f_{U_n, R_n}(U, R)$ the joint density of R_n and U_n , and by $f_{U_n}(U)$ and $f_{R_n}(R)$ the marginal densities. Due to independence, $f_{U_n, R_n}(U, R) = f_{U_n}(U)f_{R_n}(R)$, and therefore

$$\begin{aligned}
 & E \left(|tr_n (U_n^{m_1} P_1(R_n) U_n^{m_2} P_2(R_n) \cdots U_n^{m_l} P_l(R_n))|^2 \right) \\
 (3.22) \quad &= \int_{\Omega} |tr_n (U_n^{m_1} P_1(R_n) \cdots U_n^{m_l} P_l(R_n))|^2 ds \\
 &= \int_{M_n(C)} \int_{M_n(C)} |tr_n (U^{m_1} P_1(R) \cdots U^{m_l} P_l(R))|^2 f_{U_n, R_n}(U, R) dU dR \\
 &= \int_{M_n(C)} \int_{M_n(C)} |tr_n (U^{m_1} P_1(R) \cdots U^{m_l} P_l(R))|^2 f_{U_n}(U) dU f_{R_n}(R) dR \\
 &\leq \int_{M_n(C)} C n^{-2} f_{R_n}(R) dR = C n^{-2}
 \end{aligned}$$

where we have used the bounds for deterministic matrices from theorem 3.1. Therefore

$$|tr_n (U_n^{m_1} P_1(R_n) U_n^{m_2} P_2(R_n) \cdots U_n^{m_l} P_l(R_n))| \rightarrow 0 \text{ a.s.}$$

for such random matrices R_n . If $R_n R_n^*$ is just known to converge in distribution almost surely to a compactly supported probability measure, observe that almost surely there exists a value R so that $\|R_n\|_p \leq R$ for n large enough [7]. For each l , choose p_l as in the statement of theorem 3.1. Denote by $\Omega_{p_l, N}$, $p \geq 1$ $N \in \mathbb{N}$ the subset of Ω determined by values s such that

$$(3.23) \quad \|R_n(s)\|_{p_l} \leq N$$

for all n . Define $R_{n, p_l, N} = \chi_{\Omega_{p_l, N}} R_n$ with χ denoting the characteristic function. The $R_{n, p_l, N}$ satisfy the estimates (3.22) for mixed moments of length $\leq l$, so that these mixed moments go to zero almost surely in $\Omega_{p_l, N}$. $\cup_{p_l, N} \Omega_{p_l, N}$ has probability 1: Almost surely, the $\|\cdot\|_{p_l}$ -norm of R_n stays bounded by some finite value for large enough n . Thus, for every s in a set with probability one, we can find a value N_s such that $\|R_n(s)\|_{p_l} \leq N_s$ for ALL n . But then $s \in \Omega_{p_l, N_s}$, so that $\cup_{p_l, N} \Omega_{p_l, N}$ has probability 1 as claimed. Since $\cup_N \Omega_{p_l, N}$ has probability 1, theorem 3.2 follows from the fact that $R_n = R_{n, p_l, N}$ on $\Omega_{p_l, N}$. By increasing l we get almost sure convergence of higher mixed moments to zero also.

Now for theorem 3.3. Write

$$X_n = U_n \Lambda_n U_n^*$$

for a unitary random matrix U_n , and diagonal random matrix Λ_n . We may assume that U_n is a standard unitary random matrix, as in the proof of theorem 4.3.5 of [7], since Gaussian random matrices are unitarily invariant. We can also assume that Λ_n is independent from U_n , so that $(U_n, \{\Lambda_n, R_n\})$ is an independent family. R_n converges to a limit which is compactly supported, and Λ_n does the same. Since (3.22) can be easily generalized to the case where the R_n are replaced with many different R_n (with the $\{R_n\}$ all independent from U_n) we conclude also for theorem 3.3 that we get almost sure convergence to zero of mixed moments as in definition 2.1.

3.3. *The proof of theorem 3.4.* First write $\phi((p(a+b)(a+b)^*p)^m)$ as a sum of mixed moments of length $3m$ by multiplying out $(p(a+b)(a+b)^*p)^m$:

$$(3.24) \quad \phi((p(a+b)(a+b)^*p)^m) = \sum_{\sigma_1 \leq \sigma} \phi(x_1 x_2^* p \cdots x_{2m-1} x_{2m}^* p),$$

where $\sigma = \{1, 2, 4, 5, \dots, 3m-2, 3m-1\}$ ($\{1, 2, 4, 5, \dots, 3m-2, 3m-1\}$ correspond to the indices of the locations of the x_i, x_i^* in the moments $x_1 x_2^* p \cdots x_{2m-1} x_{2m}^* p$), σ_1 runs over all subsets of σ , and $x_i = a$ if $i \in \sigma_1$, $x_i = b$ if $i \in \sigma \setminus \sigma_1$. Denote by $|\sigma_1|$ the cardinality of σ_1 . We denote by α the cumulants of μ_{a,a^*} and β the cumulants of $\mu_{b,b^*,p}$, so that the moment-cumulant formula for (a, a^*) is

$$(3.25) \quad \phi(x_{i_1} \cdots x_{i_n}) = \sum_{\pi = \{B_1, \dots, B_k\} \in NC(n)} \prod_{i=1}^k [\text{coef}(i_1, \dots, i_n) | B_i](R_{\mu_{a,a^*}})$$

with $x_1 = a, x_2 = a^*$ and $i = 1$ or 2 , and the moment-cumulant formula for (b, b^*, p) is

$$(3.26) \quad \phi(x_{i_1} \cdots x_{i_n}) = \sum_{\pi = \{B_1, \dots, B_k\} \in NC(n)} \prod_{i=1}^k [\text{coef}(i_1, \dots, i_n) | B_i](R_{\mu_{b,b^*,p}}),$$

with $x_1 = b, x_2 = b^*, x_3 = p$ and $i = 1, 2$ or 3 . We will use the shorthand notation

$$\begin{aligned} \alpha_{B_i} &= [\text{coef}(i_1, \dots, i_n) | B_i](R_{\mu_{a,a^*}}) \\ \beta_{B_i} &= [\text{coef}(i_1, \dots, i_n) | B_i](R_{\mu_{b,b^*,p}}) \end{aligned}$$

Due to the freeness of a and $\{p, b\}$, the moment-cumulant formula applied to all moments in (3.24) and (2.14) yields

$$(3.27) \quad \sum_{\sigma_1 \leq \sigma} \sum_{\substack{\pi_1 \in NC(|\sigma_1|) \\ \pi_1 \leq \sigma_1}} \sum_{\substack{\pi_2 \in NC(|\sigma_1^c|) \\ \pi_2 \leq \sigma_1^c, \text{ no crossings between } \pi_1 \text{ and } \pi_2}} \alpha_{\pi_1} \beta_{\pi_2},$$

where $\sigma_1^c = \{1, \dots, 3m\} \setminus \sigma_1$ and $|\sigma_1^c|$ is the cardinality of σ_1^c . π_1 divides $\{1, \dots, |\sigma_1|\}$ into $|K(\pi_1)|$ sets (see definition 2.3 and figure 1) when π_1 is viewed as an element in $NC(|\sigma_1|)$. π_1 also divides $\{1, \dots, 3m\}$ into the same number of sets, according to the circular representation of $\{1, \dots, 3m\}$. Let us denote these blocks by B_1, \dots, B_k , so that $\sigma_1^c = \{B_1, \dots, B_k\}$ as a subpartition of $1_{|\sigma_1^c|}$. Since π_1 and π_2 have no crossings if and only if $\pi_2 \leq \{B_1, \dots, B_k\}$, (3.27) can be written as

$$(3.28) \quad \sum_{\sigma_1 \leq \sigma} \sum_{\substack{\pi_1 \in NC(|\sigma_1|) \\ \pi_1 \leq \sigma_1}} \sum_{\substack{\pi_2 \in NC(|\sigma_1^c|) \\ \pi_2 \leq \{B_1, \dots, B_k\}}} \alpha_{\pi_1} \beta_{\pi_2}$$

When $\pi_2 \leq \{B_1, \dots, B_k\}$ we can write $B_i = \pi_{2i1} \cup \pi_{2i2}, \dots$, where the π_{2ij} are the reindexed blocks of π_2 which are contained in B_i , and where $\pi_{2i} = \{\pi_{2i1}, \pi_{2i2}, \dots\}$. This is in $NC(|B_i|)$ since π_2 is noncrossing. First rewrite (3.28) to

$$(3.29) \quad \sum_{\sigma_1 \leq \sigma} \sum_{\substack{\pi_1 \in NC(|\sigma_1|) \\ \pi_1 \leq \sigma_1}} \alpha_{\pi_1} \left(\sum_{\pi_{2i} \in NC(|B_i|)} \prod_{i=1}^k \beta_{\pi_{2i}} \right).$$

Then note that the $\pi_{2i} \in NC(|B_i|)$ can be summed independently of one another, so that we can rewrite to

$$(3.30) \quad \sum_{\sigma_1 \leq \sigma} \sum_{\substack{\pi_1 \in NC(|\sigma_1|) \\ \pi_1 \leq \sigma_1}} \alpha_{\pi_1} \prod_{i=1}^k \left(\sum_{\pi_{2i} \in NC(|B_i|)} \beta_{\pi_{2i}} \right).$$

Note also that only π_1 with blocks

$$C_k = \{c_{k1}, \dots, c_{kr}\}$$

where $x_{i_{c_{k1}}}, \dots, x_{i_{c_{kr}}}$ are alternating values of a and a^* , give contribution in (3.30), due to R -diagonality of a . Hold such a π_1 fixed in (3.30), and take a look at the inner sum in (3.30) for a given i . This is simply the moment-cumulant formula (3.26) for a moment of length $|B_i|$, where the mixed moment is on the form

$$pbb^*pb^*p \cdots bb^*p,$$

or on the form

$$b^*pbb^*pb^*p \cdots bb^*pb$$

due to the alternating structure in (3.24). In both cases the moment-cumulant formula yields $\phi\left((pbb^*p)^{\frac{|B_i|}{2}}\right)$ for the inner sum in (3.30). Therefore, we get that (3.30) equals

$$(3.31) \quad \sum_{\sigma_1 \leq \sigma} \sum_{\substack{\pi_1 \in NC(|\sigma_1|) \\ \pi_1 \leq \sigma_1}} \alpha_{\pi_1} \prod_{i=1}^k \phi\left((pbb^*p)^{\frac{|B_i|}{2}}\right).$$

Since a is R -diagonal, the α_{π_1} which give contribution in (3.31) are uniquely identified by the moments $\phi((aa^*)^m)$. Therefore, the moments

$$\phi((p(a+b)(a+b)^*p)^m)$$

are entirely identified by the moments $\phi((aa^*)^m)$ and $\phi((pbb^*p)^m)$, so that $\mu_{p(a+b)(a+b)^*p}$ only depends on μ_{aa^*} and μ_{pbb^*p} . All distributions are here in the noncommutative probability space (\mathcal{A}, ϕ) , not yet in the reduced space $(p\mathcal{A}p, \phi(p)^{-1}\phi)$. If we can prove the theorem when p and b are free, it will also hold when p and b are not free since $\mu_{p(a+b)(a+b)^*p}$ only depends on μ_{aa^*} and μ_{pbb^*p} .

We can replace μ_{b,b^*} with the (unique) R -diagonal pair b_0 so that $\mu_{bb^*} = \mu_{b_0b_0^*}$. So, we assume that both a and b give rise to free R -diagonal pairs. Their determining series are $R_{\mu_{aa^*}} \boxtimes Moeb$ and $R_{\mu_{bb^*}} \boxtimes Moeb$, respectively. $(a+b, (a+b)^*)$ is also an R -diagonal pair, with determining series $R_{\mu_{aa^*}} \boxtimes Moeb + R_{\mu_{bb^*}} \boxtimes Moeb$. This means that

$$(3.32) \quad R_{\mu_{(a+b)(a+b)^*}} \boxtimes Moeb = R_{\mu_{aa^*}} \boxtimes Moeb + R_{\mu_{bb^*}} \boxtimes Moeb.$$

If x is free from p , we next calculate $R_{\mu_{pxx^*p}}$ in the reduced space $(p\mathcal{A}p, \phi(p)^{-1}\phi)$. We call this $R_{\mu_{pxx^*p}}^{p\mathcal{A}p}$ in the rest of the proof, with similar notation for the moment series (In the rest of the paper, this notation is dropped since pxx^*p is assumed to be in $(p\mathcal{A}p, \phi(p)^{-1}\phi)$). Note that $M_{\mu_{pxx^*p}}^{p\mathcal{A}p} = \frac{1}{c} M_{\mu_{pxx^*p}}$. We have

$$\begin{aligned} R_{\mu_{pxx^*p}}^{p\mathcal{A}p} &= M_{\mu_{pxx^*p}}^{p\mathcal{A}p} \boxtimes Moeb = \left(\frac{1}{c} M_{\mu_{pxx^*p}}\right) \boxtimes Moeb \\ &= \left(\frac{1}{c} (R_{\mu_{xx^*}} \boxtimes M_{\mu_p})\right) \boxtimes Moeb = \left(\frac{1}{c} (R_{\mu_{xx^*}} \boxtimes (cZeta))\right) \boxtimes Moeb \\ &= \left(\frac{1}{c} (R_{\mu_{xx^*}} \boxtimes (cZeta))\right) \boxtimes \left(\frac{1}{c} (cMoeb)\right) \\ &= c^{-n-1} (R_{\mu_{xx^*}} \boxtimes (cZeta) \boxtimes (cMoeb)) \\ &= c^{-n-1} (R_{\mu_{xx^*}} \boxtimes (c^{n+1}Id)) = c^{-n-1} (R_{\mu_{xx^*}} \boxtimes (c^2Id)) \\ &= c^{-n-1} (c^{2n} R_{\mu_{xx^*}}) = c^{-n-1} R_{\mu_{xx^*}}. \end{aligned}$$

For a general Marčenko Pastur law μ_d , $R_{\mu_d} = d^{n-1}Zeta$, and it is easily verified that $R_{\mu_d}^{-1} = d^{n-1}Moeb$. We have that

$$R_{\mu_{pxx^*p}}^{p\mathcal{A}p} \boxtimes R_{\mu_d}^{-1} = \left(c^{-n-1} R_{\mu_{xx^*}}\right) \boxtimes (d^{n-1}Moeb) = c^n d^n \left(\left(c^{-1} R_{\mu_{xx^*}}\right) \boxtimes (d^{-1}Moeb)\right)$$

if $c = d$, this can be simplified to

$$(3.33) \quad c^n c^n c^{-n-1} (R_{\mu_{xx}^*} \boxtimes Moeb) = c^{n-1} (R_{\mu_{xx}^*} \boxtimes Moeb).$$

Using this for $x = a$, $x = b$ and $x = a + b$, and also using (3.32), we get

$$\begin{aligned} R_{\mu_{p(a+b)(a+b)^*p}^{pAp}} \boxtimes R_{\mu_c}^{-1} &= c^{n-1} (R_{\mu_{(a+b)(a+b)^*}^*} \boxtimes Moeb) \\ &= c^{n-1} (R_{\mu_{aa}^*} \boxtimes Moeb + R_{\mu_{bb}^*} \boxtimes Moeb) \\ &= c^{n-1} (R_{\mu_{aa}^*} \boxtimes Moeb) + c^{n-1} (R_{\mu_{bb}^*} \boxtimes Moeb) \\ &= R_{\mu_{paa^*p}^{pAp}} \boxtimes R_{\mu_c}^{-1} + R_{\mu_{pbb^*p}^{pAp}} \boxtimes R_{\mu_c}^{-1} \end{aligned}$$

Using (2.17), this can be written

$$R_{\mu_{p(a+b)(a+b)^*p}^*} \boxtimes \mu_c = R_{\mu_{paa^*p}^*} \boxtimes \mu_c + R_{\mu_{pbb^*p}^*} \boxtimes \mu_c,$$

which can equivalently be stated as

$$\left(\mu_{p(a+b)(a+b)^*p} \boxtimes \mu_c \right) = (\mu_{paa^*p} \boxtimes \mu_c) \boxplus (\mu_{pbb^*p} \boxtimes \mu_c),$$

which is what we had to prove.

Note that if $c \neq d$, $R_{\mu_{xx}^*} \boxtimes Moeb$ does not appear as a factor in (3.33). Therefore, there is no reason why the result should hold for other Marčenko Pastur laws than μ_c , since (3.32) can not be used in such cases.

Although the proof of theorem 1.1 is described for $c \leq 1$ only, the methods used in the proof of theorem 3.4 here can help us prove the case for $c > 1$ also. If R_n and X_n are the random matrices from theorem 1.1 and $c > 1$, note that

$$M_{\mu_{\frac{1}{N}R_n R_n^*}} = \frac{1}{c} M_{\mu_{\frac{1}{N}R_n^* R_n}} = \frac{1}{c} M_{\mu_{\frac{1}{c} \frac{1}{N} R_n^* R_n}} = c^{-m-1} M_{\mu_{\frac{1}{N} R_n^* R_n}},$$

where $c^m f$ denoted the power series defined by $[coe f_k](c^m f) = c^k [coe f_k](f)$. From this one can show that

$$(3.34) \quad \begin{aligned} R_{\mu_{\frac{1}{N}R_n R_n^*}} \boxtimes R_{\mu_c}^{-1} &= M_{\mu_{\frac{1}{N}R_n R_n^*}} \boxtimes Moeb \boxtimes (c^{m-1} Moeb) \\ &= \left(c^{-m-1} M_{\mu_{\frac{1}{N}R_n^* R_n}} \right) \boxtimes Moeb \boxtimes (c^{m-1} Moeb) \\ &= \left(c^{-m-1} \left((c^{-m}) M_{\mu_{\frac{1}{N}R_n^* R_n}} \right) \boxtimes (c^m Moeb) \right) \boxtimes Moeb \\ &= \left(c^{-m-1} (M_{\mu_{\frac{1}{N}R_n^* R_n}} \boxtimes Moeb) \right) \boxtimes (c^{-1} (c^{-m+1} Moeb)) \\ &= c^{-m} c^{-m-1} \left(M_{\mu_{\frac{1}{N}R_n^* R_n}} \boxtimes Moeb \boxtimes (c^{-m+1} Moeb) \right) \\ &= c^{-2m-1} \left(R_{\mu_{\frac{1}{N}R_n^* R_n}} \boxtimes R_{\mu_{\frac{1}{c}}}^{-1} \right). \end{aligned}$$

Since theorem 1.1 has been proved for $c \leq 1$, $R_{\mu_{\frac{1}{n}(R_n+X_n)^*(R_n+X_n)}} \boxtimes R_{\mu_{\frac{1}{c}}}$ will converge to

$$R_{\mu_{\frac{1}{n}R_n^*R_n}} \boxtimes R_{\mu_{\frac{1}{c}}}^{-1} + R_{\mu_{\frac{1}{n}X_n^*X_n}} \boxtimes R_{\mu_{\frac{1}{c}}}^{-1}$$

as $n \rightarrow \infty$ ($c > 1$), and the result for $c > 1$ follows from (3.34).

4. Equivalence with known expressions for limit distributions of Information-Plus-Noise Type Matrices. [14] studies systems where the sample covariance matrix is formed by taking independent samples of a system of the form

$$y_n = A_n x_n + \sigma w_n$$

(the noise factor σ does not appear in [14]) where x_n and w_n are independent standard (zero mean, unit variance) Gaussian random vectors, and A_n is an $n \times L$ matrix. The covariance matrix of the system is $\Theta_n = A_n A_n^* + \sigma^2 I$. In particular, when there is no noise (i.e. $\sigma = 0$), the covariance is $\Theta_n = A_n A_n^*$. Denote by μ_{Θ} the limiting eigenvalue distribution of $A_n A_n^*$. [14] states that the limiting eigenvalue distribution of the sample covariance matrix of the system is $(\mu_{\Theta} \boxplus \mu_{\sigma^2 I}) \boxtimes \mu_c$. When there is no noise, the limit is $\mu_{\Theta} \boxtimes \mu_c$. This way of passing from

$$\mu_{\Theta} \boxtimes \mu_c \text{ to } (\mu_{\Theta} \boxplus \mu_{\sigma^2 I}) \boxtimes \mu_c$$

is of course compatible with theorem 1.1. We will also show that it is equivalent with the results in [3]. The following restrictions taken from [3] will be used:

1. For $n = 1, 2, \dots$, $X_n = (X_{ij}^n, n \times N, \text{i.d. for all } i, j, n, \text{ independent across } i, j \text{ for each } n, \text{ and } E|X_{11}^1 - EX_{11}^1|^2 = 1$
2. R_n is $n \times N$ and independent of X_n , with $F^{\mu_{\Gamma_n}} \xrightarrow{\mathcal{D}} F^{\mu_{\Gamma}}$

Theorem 1.1 in [3] expresses a relationship for finding the limiting eigenvalue distribution μ_W of W_n from that of $\Gamma_n = \frac{1}{N} R_n R_n^*$ (denoted μ_{Γ}). More precisely, under the conditions 1) and 2), we have that (in a slightly rewritten form) $F^{\mu_{W_n}} \xrightarrow{\mathcal{D}} F^{\mu_W}$ almost surely, where F^{μ_W} is a nonrandom p.d.f. characterized by

$$(4.1) \quad m_{\mu_W}(z) = \int \frac{dF^{\mu_{\Gamma}}(t)}{\frac{t}{1+\sigma^2 c m_{\mu_W}(z)} - (1 + \sigma^2 c m_{\mu_W}(z))z + \sigma^2(1-c)}$$

for any $z \in C^+$. The connection with multiplicative free convolution is hard to see from this formula. To obtain this connection, the following lemma is needed, which will be proved in section 4.1:

LEMMA 4.1. (4.1) is equivalent to

$$(4.2) \quad m_{\mu_W}^{-1} \left(\frac{z}{1 - \sigma^2 cz} \right) = (1 - \sigma^2 cz)^2 m_{\mu_\Gamma}^{-1}(z) + \sigma^2(1 - c)(1 - \sigma^2 cz),$$

for z in some interval $(0, z_1)$.

In (4.2) and all other places where the inverse of the Stieltjes transform is taken in this paper, we will mean the unique inverse on the negative real line. The inverse will only be calculated for positive values close to 0. It will turn out that (4.2) can be more conveniently expressed in terms of distributions obtained from multiplicative free deconvolution with the Marčenko Pastur law using the following lemma, which will be proved in section 4.2:

LEMMA 4.2. If

$$(4.3) \quad \mu_\Gamma = \mu_\Theta \boxtimes \mu_c,$$

then, for z in some interval $(0, z_1)$,

$$(4.4) \quad \eta_{\mu_\Gamma}^{-1}(z) = \frac{\eta_{\mu_\Theta}^{-1}(z)}{1 - c + cz}$$

and also

$$(4.5) \quad m_{\mu_\Gamma}^{-1} \left(\frac{z}{1 - c - cz m_{\mu_\Theta}^{-1}(z)} \right) = m_{\mu_\Theta}^{-1}(z)(1 - c - cz m_{\mu_\Theta}^{-1}(z)).$$

Using (4.5), the following relationship with multiplicative free convolution will be shown:

THEOREM 4.1. Under the conditions 1) and 2), assume that

$$(4.6) \quad F^{\mu_{\Gamma_n}} \xrightarrow{\mathcal{D}} F^{\mu_\Theta \boxtimes \mu_c} \text{ a.s.}$$

Then

$$(4.7) \quad F^{\mu_{W_n}} \xrightarrow{\mathcal{D}} F^{(\mu_\Theta \boxplus \mu_{\sigma^2 I}) \boxtimes \mu_c} \text{ a.s.}$$

Equivalently, assume that

$$(4.8) \quad F^{\mu_{\Gamma_n}} \xrightarrow{\mathcal{D}} F^{\mu_\Gamma} \text{ a.s.}$$

Then

$$(4.9) \quad F^{\mu_{W_n}} \xrightarrow{\mathcal{D}} F^{\mu_W} \text{ a.s.}$$

where μ_W is uniquely identified by the equation

$$(4.10) \quad \mu_W \boxtimes \mu_c = (\mu_\Gamma \boxtimes \mu_c) \boxplus \mu_{\sigma^2 I},$$

Theorem 4.1 will be proved in section 4.3.

4.1. *The proof of lemma 4.1.* Rewritten in terms of the Stieltjes transform, (4.1) says that (with terms somewhat regrouped)

$$\frac{m_{\mu_W}}{1 + \sigma^2 c m_{\mu_W}} = m_{\mu_\Gamma} \left((1 + \sigma^2 c m_{\mu_W})^2 z - \sigma^2 (1 - c)(1 + \sigma^2 c m_{\mu_W}) \right),$$

where m_{μ_W} , m_{μ_Γ} are evaluated in z when the parameter is omitted. We restrict ourselves to z on the negative real line. The relation holds also for such z , since we can analytically continue to the negative real line. Evaluating in $m_{\mu_W}^{-1}(z)$ we get

$$\frac{z}{1 + \sigma^2 c z} = m_{\mu_\Gamma} \left((1 + \sigma^2 c z)^2 m_{\mu_W}^{-1}(z) - \sigma^2 (1 - c)(1 + \sigma^2 c z) \right)$$

for z in some interval $(0, z_1)$. We will find it convenient to work with the inverse of the Stieltjes transform, so we rewrite the expression to

$$m_{\mu_\Gamma}^{-1} \left(\frac{z}{1 + \sigma^2 c z} \right) = (1 + \sigma^2 c z)^2 m_{\mu_W}^{-1}(z) - \sigma^2 (1 - c)(1 + \sigma^2 c z).$$

Substituting $u = \frac{z}{1 + \sigma^2 c z}$ (or equivalently $z = \frac{u}{1 - \sigma^2 c u}$) (this is an isomorphism of the positive real axis which sends 0 to 0), we get

$$m_{\mu_\Gamma}^{-1}(z) = \frac{m_{\mu_W}^{-1} \left(\frac{z}{1 + \sigma^2 c z} \right)}{(1 - \sigma^2 c z)^2} - \frac{\sigma^2 (1 - c)}{1 - \sigma^2 c z}$$

for z in some interval $(0, z_1)$, so that

$$m_{\mu_W}^{-1} \left(\frac{z}{1 - \sigma^2 c z} \right) = (1 - \sigma^2 c z)^2 m_{\mu_\Gamma}^{-1}(z) + \sigma^2 (1 - c)(1 - \sigma^2 c z),$$

which is (4.2).

4.2. *The proof of lemma 4.2.* By the multiplicative property of the S-transform we have

$$S_{\mu_\Gamma}(z) = \frac{S_{\mu_\Theta}(z)}{1 + cz}.$$

Expressed in terms of the η -transform this can be written

$$\eta_{\mu_\Gamma}^{-1}(z) = \frac{\eta_{\mu_\Theta}^{-1}(z)}{1 - c + cz},$$

which is (4.4). Evaluating in $\eta_{\mu_\Theta}(z)$ and applying η_{μ_Γ} on both sides gives

$$\eta_{\mu_\Gamma} \left(\frac{z}{1 - c + c \eta_{\mu_\Theta}(z)} \right) = \eta_{\mu_\Theta}(z)$$

for $z \geq 0$. This can also be expressed in terms of Stieltjes transforms as

$$\frac{m_{\mu_\Gamma} \left(-\frac{1-c+c\eta_{\mu_\Theta}(z)}{z} \right)}{\frac{z}{1-c+c\eta_{\mu_\Theta}(z)}} = \frac{m_{\mu_\Theta} \left(-\frac{1}{z} \right)}{z}$$

Regrouping terms and substituting $-\frac{1}{z}$ for z we get

$$m_{\mu_\Gamma} (z(1-c-czm_{\mu_\Theta}(z))) = \frac{m_{\mu_\Theta}(z)}{1-c-czm_{\mu_\Theta}(z)}$$

for $z < 0$. Substituting $m_{\mu_\Theta}^{-1}(z)$ for z and taking the inverse Stieltjes transform $m_{\mu_\Gamma}^{-1}$ we get

$$m_{\mu_\Gamma}^{-1} \left(\frac{z}{1-c-czm_{\mu_\Theta}^{-1}(z)} \right) = m_{\mu_\Theta}^{-1}(z)(1-c-czm_{\mu_\Theta}^{-1}(z))$$

for z in some interval $(0, z_1)$, which is (4.5).

4.3. *The proof of theorem 4.1.* Note that if

$$z = \frac{z_1}{1-c-cz_1m_{\mu_\Theta}^{-1}(z_1)}$$

for z_1 positive and close to 0, then we have

$$(4.11) \quad \frac{z}{1-\sigma^2cz} = \frac{z_1}{1-c-cz_1(m_{\mu_\Theta}^{-1}(z_1) + \sigma^2)}.$$

Note also that $\frac{z}{1-\sigma^2cz} \geq 0$ as long as $z < \frac{1}{\sigma^2c}$. Substituting (4.11) and (4.5)

in (4.2) we get $m_{\mu_W}^{-1} \left(\frac{z}{1-\sigma^2cz} \right) =$

$$(4.12) \quad \begin{aligned} & \left(1 - \frac{\sigma^2cz_1}{1-c-cz_1m_{\mu_\Theta}^{-1}(z_1)} \right)^2 m_{\mu_\Theta}^{-1}(z_1)(1-c-cz_1m_{\mu_\Theta}^{-1}(z_1)) \\ & + \sigma^2(1-c) \left(1 - \frac{\sigma^2cz_1}{1-c-cz_1m_{\mu_\Theta}^{-1}(z_1)} \right) \\ & = \frac{(1-c-cz_1m_{\mu_\Theta}^{-1}(z_1)-\sigma^2cz_1)^2 m_{\mu_\Theta}^{-1}(z_1) + \sigma^2(1-c)(1-c-cz_1m_{\mu_\Theta}^{-1}(z_1)-\sigma^2cz_1)}{1-c-cz_1m_{\mu_\Theta}^{-1}(z_1)} \\ & = \frac{(1-c-cz_1m_{\mu_\Theta}^{-1}(z_1)-\sigma^2cz_1)((1-c-cz_1m_{\mu_\Theta}^{-1}(z_1)-\sigma^2cz_1)m_{\mu_\Theta}^{-1}(z_1) + \sigma^2(1-c))}{1-c-cz_1m_{\mu_\Theta}^{-1}(z_1)} \\ & = \frac{(1-c-cz_1(m_{\mu_\Theta}^{-1}(z_1) + \sigma^2))(1-c-cz_1m_{\mu_\Theta}^{-1}(z_1))(m_{\mu_\Theta}^{-1}(z_1) + \sigma^2)}{1-c-cz_1m_{\mu_\Theta}^{-1}(z_1)} \\ & = \left(m_{\mu_\Theta}^{-1}(z_1) + \sigma^2 \right) \left(1 - c - cz_1(m_{\mu_\Theta}^{-1}(z_1) + \sigma^2) \right) \\ & = m_{\mu_{\Theta \boxplus \mu_{\sigma^2 I}}}^{-1}(z_1) \left(1 - c - cz_1m_{\mu_{\Theta \boxplus \mu_{\sigma^2 I}}}^{-1}(z_1) \right) \end{aligned}$$

Here we have used that $m_{\mu_{\Theta} \boxplus \mu_{\sigma^2 I}}^{-1}(z_1) = m_{\mu_{\Theta}}^{-1}(z_1) + \sigma^2$, which follows from the additivity property of the R -transform and the fact that the inverse of the Stieltjes transform is used to define the R -transform (perform additive free convolution with $\mu_{\sigma^2 I}$). (4.12) is thus nothing else than (4.5) (with $m_{\mu_{\Theta}}$ replaced by $m_{\mu_{\Theta} \boxplus \mu_{\sigma^2 I}}$). Since (4.5) is just an equivalent expression for multiplicative free convolution, we therefore have

$$\mu_W = (\mu_{\Theta} \boxplus \mu_{\sigma^2 I}) \boxtimes \mu_c,$$

or equivalently

$$\mu_W \boxtimes \mu_c = \mu_{\Theta} \boxplus \mu_{\sigma^2 I} = (\mu_{\Gamma} \boxtimes \mu_c) \boxplus \mu_{\sigma^2 I}.$$

This completes the proof.

5. Using G -analysis to estimate the spectral function of covariance matrices. It turns out that multiplicative free deconvolution can also be used to estimate covariance matrices. The general statistical analysis of observations, also called G -analysis [5] is a mathematical theory for complex systems where the number of parameters of the underlying mathematical model increase together with the growth of the number of observations of the system. The mathematical models which approach the system in some sense are called G -estimators. The main difficulty in G -analysis is to find good G -estimators. G -estimators have already shown their usefulness in many applications [9]. We denote by N the number of observations of the system, and by n the number of parameters of the mathematical model. The condition used in G -analysis expressing the growth of the number of observations vs. the number of parameters in the mathematical model, is called the G -condition. The G -condition used throughout this paper is (1.2).

Girko restricts to systems where a number of independent random vector observations are taken, and where the random vectors have identical distributions. If a random vector r_n has length n , we will let Θ_n denote its covariance, while Γ_n will still denote sample covariance matrices. The Γ_n we analyze in this section are more restrictive than in previous sections, since independence across samples is assumed. Girko calls estimators for the Stieltjes transform of covariance matrices G^2 -estimators. In chapter 2.1 of [4] he introduces the following expression as candidate for a G^2 -estimator:

$$(5.1) \quad G_n^2(z) = \frac{\hat{\theta}(z)}{z} m_{\mu_{\Gamma_n}}(\hat{\theta}(z)),$$

where the function $\hat{\theta}(z)$ is the solution to the equation

$$(5.2) \quad \hat{\theta}(z) c m_{\mu_{\Gamma_n}}(\hat{\theta}(z)) - (1 - c) + \frac{\hat{\theta}(z)}{z} = 0.$$

Girko claims that a function $G_n^2(z)$ satisfying (5.2) and (5.1) is a good approximation for the Stieltjes transform of the covariance matrices $m_{\Theta_n}(z) = \text{tr}_n \{\Theta_n - zI_n\}^{-1}$. More precisely, he shows that when (5.1), (5.2) and the G -condition (1.2) are fulfilled, under certain conditions there exists a $c > 0$ such that

$$(5.3) \quad \lim_{n \rightarrow \infty} \sup_{\substack{0 < c \leq \Im(z) \leq S \\ |\Re(z)| \leq T}} |G_n^2(z) - m_{\Theta_n}(z)| = 0,$$

with probability one for every $S > 0$ and $T > 0$. According to Girko, analytical continuation of $G_n^2(z)$ can be performed to obtain limits for other z than the ones in (5.3).

As it turns out, the G^2 -estimator can equivalently be expressed in terms of multiplicative free convolution:

THEOREM 5.1. *For the G^2 -estimator given by (5.1), (5.2), the following holds for real $z < 0$:*

$$(5.4) \quad G_n^2(z) = m_{\mu_{\Gamma_n} \boxtimes \mu_c}$$

PROOF. (5.2) can be rewritten to

$$\begin{aligned} -c\eta_{\mu_{R_n}} \left(-\frac{1}{\hat{\theta}(z)} \right) - (1-c) + \frac{\hat{\theta}(z)}{z} &= 0 \\ \eta_{\mu_{R_n}}^{-1} \left(\frac{1}{c} \left(\frac{\hat{\theta}(z)}{z} - (1-c) \right) \right) &= -\frac{1}{\hat{\theta}(z)}, \end{aligned}$$

which we will write

$$(5.5) \quad -\frac{1}{\eta_{\mu_{R_n}}^{-1} \left(\frac{1}{c} \left(\frac{\hat{\theta}(z)}{z} - (1-c) \right) \right)} = \hat{\theta}(z).$$

Denote by μ the measure with Stieltjes transform $G_n^2(z)$. (5.1) can be rewritten using the η -transform as

$$\eta_{\mu} \left(-\frac{1}{z} \right) = \eta_{\mu_{R_n}} \left(-\frac{1}{\hat{\theta}(z)} \right).$$

Since $\eta_{\mu_{R_n}}$ and η_{μ} are monotone, it is easily seen from this that $\hat{\theta}$ is monotone since it is a combination of monotone functions. Forming the inverse functions on both sides, and also applying $\hat{\theta}$, yields

$$(5.6) \quad \hat{\theta} \left(-\frac{1}{\eta_{\mu}^{-1}(z)} \right) = -\frac{1}{\eta_{\mu_{R_n}}^{-1}(z)}.$$

Showing $\mu_{R_n} = \mu \boxtimes \mu_c$ is equivalent to (after rearranging (4.4))

$$-\frac{1}{\eta_{\mu}^{-1}(z)} = -\frac{1}{(1-c+cz)\eta_{\mu_{R_n}}^{-1}(z)}$$

Applying $\hat{\theta}$ on both sides and using (5.6) yields that this is equivalent to

$$(5.7) \quad -\frac{1}{\eta_{\mu_{R_n}}^{-1}(z)} = \hat{\theta} \left(-\frac{1}{(1-c+cz)\eta_{\mu}^{-1}(z)} \right)$$

Observe now that (5.7) and (5.5) are related in the following way: If we substitute $z = \frac{1}{c} \left(\frac{\hat{\theta}(w)}{w} - (1-c) \right)$ into (5.7), the argument on the right hand side can be rewritten using (5.5) to

$$-\frac{1}{\left(1-c+\frac{\hat{\theta}(z)}{z}-(1-c)\right) \left(-\frac{1}{\hat{\theta}(z)}\right)} = z,$$

so that (5.7) is nothing but a restatement of (5.5), at least on values of the form $z = \frac{1}{c} \left(\frac{\hat{\theta}(w)}{w} - (1-c) \right)$. If these values take on an open set of real values, equality in (5.7) follows for all z by analytic continuation. This happens when $\hat{\theta}(z) \neq kz$ for some constant k . If $\hat{\theta}(z) = kz$, then $\eta_{\mu_{R_n}}$ is seen to be constant, which only happens in trivial cases. Thus we have that $\mu_{R_n} = \mu \boxtimes \mu_c$, and we are done \square

Several remarks concerning theorem 5.1 are in place. First of all, the G^2 -estimator has a much shorter expression in terms of multiplicative free deconvolution, which also places it as an ingredient in theorem 1.1. The theorem is nice to combine with continuity results for free convolution. Voiculescu has proved such results when convergence is in the weak- $*$ topology [1]. This enables us in many cases to conclude that

$$\lim_{n \rightarrow \infty} G_n^2(z) = \lim_{n \rightarrow \infty} m_{\mu_{\Gamma_n} \boxtimes \mu_c} = m_{\mu_{\Gamma} \boxtimes \mu_c}$$

for some probability measure μ_{Γ} . Secondly, [14] expresses the exact same estimator, i.e.

$$\lim_{n \rightarrow \infty} \mu_{\Gamma_n} \boxtimes \mu_c = \lim_{n \rightarrow \infty} \mu_{\Theta_n}$$

in the case of Gaussian systems. Theorem 5.1 can be seen as a way of generalizing from the Gaussian case.

6. Further work. The concept of freeness and free convolution can be extended to unbounded random variables and general probability measures. In [1] it is shown how this can be done in the context of unbounded operator spaces, and certain regularity properties are proved. For instance, if $\mu_n \rightarrow \mu$ and $\nu_n \rightarrow \nu$ in the weak-* topology with both $\mu \neq \delta_0$ and $\nu \neq \delta_0$, then $\mu_n \boxtimes \nu_n \rightarrow \mu \boxtimes \nu$ in the weak-* topology also. It is possible that applying such extensions together with the methods applied here can extend the results to the same generality as those in [3]. This may be addressed in a future paper.

The G^2 -estimator is just one of many estimators introduced by Girko. He has estimators for many other quantities also [4], like for the square root and the moments of covariance matrices. Certain of these estimators may also have alternative expressions in terms of free probability constructs.

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