

# Self-averaging and Spectral Commutativity of Isotropic Random Matrices

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## Abstract

We show that the eigenvalue distribution for the product of arbitrary number of random matrices from isotropic unitary ensemble (IUE) is independent of the order of multiplication in  $N \rightarrow \infty$  limit. We present the procedure of calculating the average eigenvalue distribution for product of such matrices and employ it to numerous examples involving products of Girko-Ginibre matrices and so called "matrices with predefined eigenvalue distribution". In each case we obtain perfect agreement with Monte-Carlo simulations considering finite size effects.

## Isotropic Random Matrices

- Isotropic random variables:      Isotropic random matrices:
- ▶  $z = r e^{i\phi}$ ,
  - ▶  $r \in \mathbb{R}, r \geq 0$ ,
  - ▶  $\phi$  uniform on  $[0, 2\pi)$ .
- ▶  $x = hu$ ,
  - ▶  $h$  - positive semi-definite Hermitian,
  - ▶  $u$  - Haar measure on  $U(N)$ .

In other words: For  $N \times N$  Isotropic Random Matrix  $x$  one has  $P(x) = P(xv)$  where  $v \in U(N)$

Examples:

- ▶ Girko-Ginibre matrix
- ▶ Matrix of the form  $vhv$  where  $v$  and  $u$  are Unitary Haar measure random matrices and  $h$  is positive semi-definite hermitian random matrix.

Properties:

- ▶ Eigenvalue spectrum is rotationally symmetric on the complex plane.
- ▶ They form Isotropic Unitary Ensemble (IUE).

## Order invariance of average eigenvalue spectrum

Our main result:

- ▶ Average eigenvalue distribution for the product of  $n$  matrices generated from any type of IUE is independent of the order of multiplication in the  $N \rightarrow \infty$  limit.

Or in other words:

- ▶ Consider  $N$  independent IUE matrices  $A_i, i = 1, 2, \dots, N$ . Defining  $x = A_1 A_2 \dots A_N$  and for any permutation  $\pi$ :  
 $x_\pi = A_{\pi(1)} A_{\pi(2)} \dots A_{\pi(N)}$  probability that the eigenvalue of  $x_\pi$  lies within a circle of radius  $r$ :  $\text{Prob}(\lambda_{x_\pi} < r) = \text{Prob}(\lambda_x < r)$

The proof is based on Haagerup Larsen Theorem and multiplicative properties of S-transform in Free Random Variables calculus.

## Haagerup-Larsen Theorem

Lets define radial cumulative density function for eigenvalues of IUE matrix  $x$  by:

$$F_x(r) = \int_0^r 2\pi |z| \rho_x(z, \bar{z}) d|z|, \quad (1)$$

where  $\rho_x(z, \bar{z})$  is the average eigenvalue distribution. Then one has the relation between it and S-transform for  $h^2$  from usual decomposition  $x = hu$ , known as Haagerup-Larsen Theorem:

$$S_{h^2}(F_x(r) - 1) = \frac{1}{r^2}. \quad (2)$$

Additionally  $S_h = S_{h'}$  if eigenvalue distribution for  $h$  are the same as of  $h'$ , so one can rewrite equation (2) in terms of  $x$  only:

$$S_{x^\dagger x}(F_x(r) - 1) = \frac{1}{r^2}. \quad (3)$$

as  $h^2 = xx^\dagger$  and it has the same eigenvalues as  $x^\dagger x$ .

Finally, using properties of S-transforms, one has relation for  $X = \prod_{i=1}^n x_i$ :

$$\prod_{i=1}^n S_{x_i^\dagger x_i}(F_X(r) - 1) = \frac{1}{r^2} \quad (4)$$

## References

- ▶ Z. Burda, M.A. Nowak and A. Swiech, *New spectral relations between products and powers of isotropic random matrices*, arXiv:1205.1625
- ▶ U. Haagerup and F. Larsen, *Journal of Functional Analysis*, 176, 331 (2000).
- ▶ D. V. Voiculescu, *J. Operator Theory* 18, 223 (1987).

## Numerical results

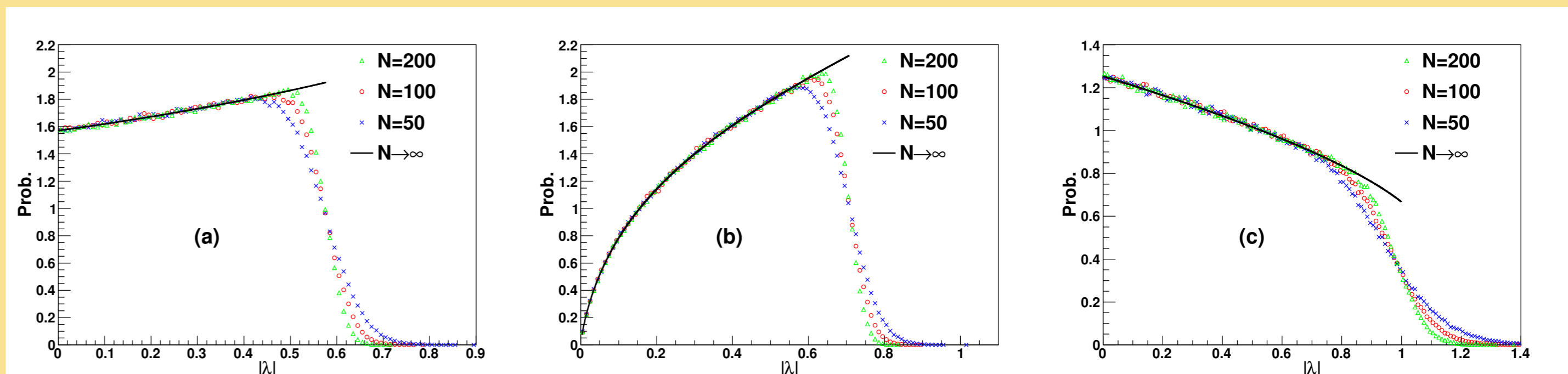


Figure: Numerical verification of theoretical predictions  $p(r) = F'(r)$  (1) for the mean radial spectral density of eigenvalues for the product of independent isotropic matrices. (a) Numerical histograms for the product of independent Girko-Ginibre matrix and matrix with  $\rho_h(r) = 1$  on interval  $(0,1)$  for  $N = 200$  (green triangles),  $N = 100$  (red circles) and  $N = 50$  (blue crosses) compared with theoretical predictions for  $N \rightarrow \infty$ . Each histogram is made for  $10^7$  eigenvalues. The numerical histograms approach theoretical curve as the size of matrices is increased. (b) An analogous plot to (a) for the product of Girko-Ginibre matrix and matrix with  $\rho_h(r) = 2r$  on interval  $(0,1)$ . (c) An analogous plot to (a) and (b) for the product of Girko-Ginibre matrix and matrix with  $\rho_h(r)$  being half-normal distribution.

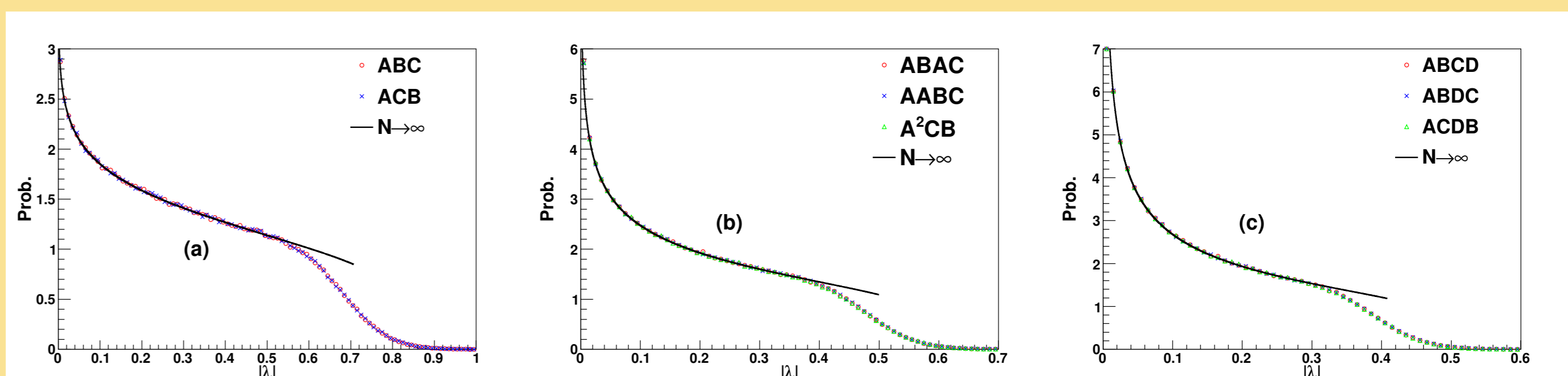


Figure: Numerical verification of theoretical predictions of the order-invariance of the mean radial spectral density,  $\rho(z, \bar{z})$ , of the eigenvalues for the product of independent isotropic matrices. By A we define matrix with  $\rho_h(r) = 2r$  on interval  $(0,1)$ , by B - Girko-Ginibre matrix, by C - matrix with  $\rho_h(r)$  being half-normal distribution and by D - matrix with  $\rho_h(r) = 1$  on interval  $(0,1)$ . (a) Numerical histograms for the product of 3 different IUE matrices for  $N = 100$ , in different order compared with theoretical predictions for  $N \rightarrow \infty$ . ABC (red circles), ACB (blue crosses). Each histogram is made for  $10^7$  eigenvalues. The numerical histograms agrees with each other as well as with theoretical prediction for  $N \rightarrow \infty$  limit taking into account finite size effects. (b) An analogous plot to (a) for the product of 4 IUE matrices, of which two are from same type.  $A_1 B A_2 C$  (red circles),  $A_1 A_2 B C$  (blue crosses) and  $A_1^2 C B$  (green triangles). (c) An analogous plot to (a) for the product of 4 different IUE matrices. ABCD (red circles), ABDC (blue crosses) and ACDB (green triangles).

## Numerical details

Having  $\rho_x(r)$  or  $\rho_h(r)$  ( $x = hu$ ), one can compute the S-transform for matrix  $h^2$ , via Haagerup-Larsen Theorem or series of equations:

$$\rho_{h^2}(r) = \frac{\partial}{\partial r} \int_0^{\sqrt{r}} \rho_h(x) dx, \quad (5)$$

$$G_{h^2}(z) = \int_{-\infty}^{\infty} \frac{\rho_{h^2}(\lambda)}{z - \lambda} d\lambda, \quad (6)$$

$$\phi(z) = \frac{1}{z} G\left(\frac{1}{z}\right) - \frac{1}{z}, \quad (7)$$

$$\chi(\phi(z)) = \phi(\chi(z)) = z, \quad (8)$$

$$S_{h^2}(z) = \frac{z+1}{z} \chi_{h^2}(z). \quad (9)$$

Having those, one can simply multiply S-transforms pointwise and use Haagerup-Larsen Theorem again to get eigenvalue distribution.

Unfortunately, the procedure is almost always impossible to perform fully analytically. There are few hints for numerical evaluation:

- ▶ The cdf takes values only from  $[0, 1]$  interval.
- ▶ The S-transforms have to be calculated and multiplied only on the  $[-1, 0]$  interval.
- ▶ The S-transforms has to be positive on the  $[-1, 0]$  interval.
- ▶ The  $\chi(z)$  is always negative and on  $[-1, 0]$  interval. It approaches 0 for  $z \rightarrow 0^-$ .
- ▶ The S-transform (and  $\chi(z)$ ) has to be continuous and injective. Altogether,  $\chi(z)$  is a continuous, monotonically rising function. This facts allows one to specify good starting parameters to solve relevant equations numerically.