

Optimal transport and the Bergman determinantal point process

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Abstract

We study the Bergman determinantal point process in a theoretical point of view aiming at its simulation. We construct restricted and restricted-truncated variants of the Bergman kernel and show optimal transport inequalities involving the Kantorovitch-Rubinstein Wasserstein distance to show to what extent it is fair to truncate the restriction of this point process to a compact ball of radius $1 - \varepsilon$. We also investigate the deviation of the number of points of the restricted Bergman determinantal point process, indicate which number of points looks like an optimal choice, and provide upper bounds to its deviation, providing an answer to an open question asked in [6].

Contents

I	Introduction	1
II	Determinantal point processes	3
III	The Bergman DPP on a disc	6
IV	The Bergman DPP on an annulus	11
V	General results	13

I Introduction

Introduced for the first time in [13], determinantal point processes (DPPs) are a particular class of point process, the correlations functions of which can be written as determinants. After having been thoroughly studied from a probabilistic point of view, (see [19] and [18] for instance, among many others), they have been used to model fermion particles [20]. They have suprisingly and remarkably appeared in the study of two very elegant mathematical problems : the first one is the point process defined as the set of eigenvalues of random $n \times n$ matrices with independent

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complex standard gaussian entries. It turns out that, as n goes to infinity, this point process converges in law to a point process that can be described [9] as a determinantal point process, the kernel of which was thereafter called the Ginibre kernel. The second one is the point process defined as the set of zeroes of Gaussian Analytical Functions (GAF) in the unit disc of the complex plane - the former are analytical functions (i.e., power series), the coefficients of which are all independent complex standard gaussian random variables : as it turns out, this point process can also be described as a determinantal point process, the kernel of which is this time the Bergman kernel (see [11], [17]). DPPs also appear and are used in the study of a wide variety of different mathematical problems, such as self-avoiding walks, random spanning trees or random integer partitions, see [1], [16], [21]. More recently, some peers attempted to apply them to machine learning [22], and to model phenomena arising in the field of networking [14], [23], [24]. The fact that these processes exhibit repulsion between particles is the core property that makes them very fitting for many applications.

The two aforementioned kernels, respectively Bergman and Ginibre kernels, benefit from several interesting properties. While the former, taking its values in the unit disc of the complex plane, is invariant with respect to rotations, the latter, which is valued in the whole complex plane, is invariant with respect to rotations and translations. Whether it be for quantum mechanics, machine learning or networking modelling, for practical uses it is necessary to benefit from a practical simulation algorithm. This work has been initiated in [10] wherein the authors give a practical algorithm for the simulation of DPPs. The simulation procedure which is hinted in [9] was fully developed in [12].

The main problem of these algorithms is that, strictly speaking, they are unfeasible due to the fact that the point processes involved exhibit an infinite number of points almost surely, and that, in the case of the Ginibre kernel, the points scatter throughout the whole plane. In [8], the authors construct restricted and truncated variants of the Ginibre point process, that allow simulation to become closer to feasible. More precisely, the Ginibre point process is therein restricted onto a ball of arbitrary radius and centered at the origin. The number of points then becomes almost surely finite. However, the stochastic dynamics of DPPs implies that computing the number of points that the point process requires to simulate a countable infinity of Bernoulli random variables, which is still unfeasible. The only option left is then to *truncate* the DPPs to a fixed (and finite) number of points, and brush off the remaining ones. But then, is it possible to quantify the approximation error that comes inevitably with this method ? To what extent is this new probability measure close to the original one ? Theoretical results were shown in [8], providing answers to this question. For instance, the authors have shown that denoting \mathfrak{S}_R and \mathfrak{S}_R^N the Ginibre DPP restricted to a ball of radius R centered at the origin of the complex plane and its truncation to N points respectively, then truncating this DPP to $N_R = (R + c)^2$ yields for all $c > 0$

$$\mathcal{W}_{KR}(\mathfrak{S}_R, \mathfrak{S}_R^{N_R}) \leq \sqrt{\frac{2}{\pi}} R e^{-c^2} \quad (1)$$

whenever $R > c$, where \mathcal{W}_{KR} denotes the Kantorovitch-Rubinstein (or Wassertein-1) distance on the set of probability measures on the set of configurations on \mathbb{C} . This result shows that the (Wasserstein) distance between the two probability measures is exponentially small in c^2 as the number $N_R = (R + c)^2$ deviates from R^2 , and implies that the two DPPs $\mathfrak{S}_R, \mathfrak{S}_R^{N_R}$ actually coincide with high probability. As a result, truncating to R^2 points looks like a very good choice.

This is all very nice and convenient, but only applies to one DPP law, the Ginibre DPP. In this paper, we extend this work, by constructing restricted (onto balls of radius R) and restricted-truncated versions of the Bergman DPP, which will henceforth be at the very core of our study. We find the compromise as to the right number of points depending on the chosen radius R to truncate our restricted DPP and provide the theoretical results that ensure that the use of the simulation algorithm presented in [8] produces a result that is not blatantly far from the (theoretical and non-feasible) simulation of the restricted Bergman DPP. In other words, the truncation error is small when choosing the right number of points, hence providing an answer to the fifth open question asked in [6] for this DPP. Some of our results are far from specific to the Bergman DPP and apply in the general case.

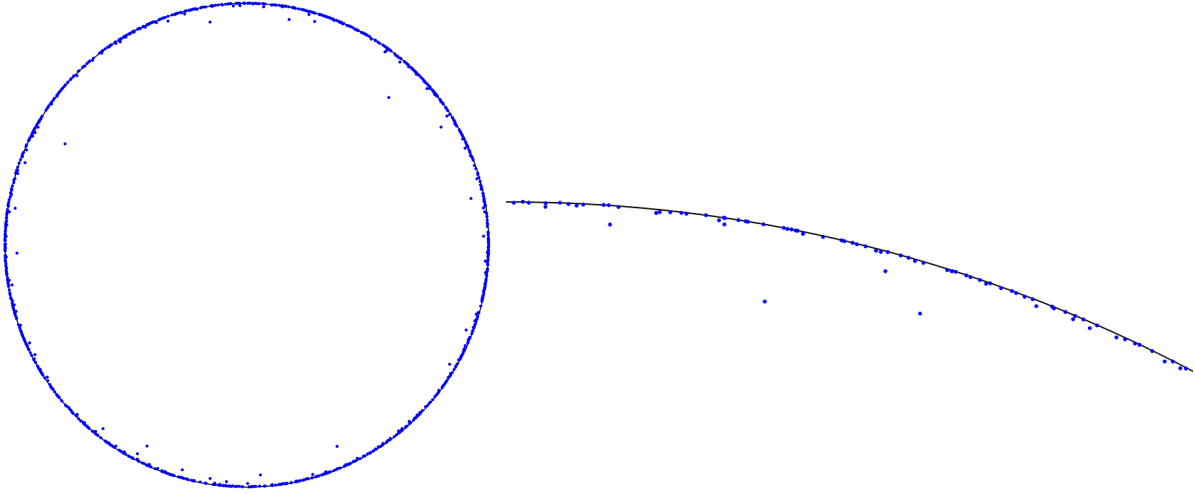


Figure 1: Left : Simulation of the Bergman DPP restricted to a radius of 0.9995, using [15]. Number of points : 985. Right : A zoom on the boundary

II Determinantal point processes

Let E be a Polish space, equipped with the Borel σ -algebra generated by the open subsets of E . In the sequel, λ is a reference Radon measure on this measurable space. Denote \mathfrak{N} the space of locally finite subsets in E , also called the configuration space :

$$\mathfrak{N} = \{\xi \subset E, \forall \Lambda \subset E \text{ compact}, |\Lambda \cap \xi| < \infty\}.$$

It is equipped with the topology of vague convergence, also called weak topology, which is the weakest topology such that for all continuous and compactly supported functions f on E , the mapping $\xi \mapsto \langle f, \xi \rangle := \sum_{x \in \xi} f(x)$ is continuous. Once topologized, it naturally comes with a σ -algebra. Elements of \mathfrak{N} , i.e., locally finite configurations on E , are identified with atomic Radon measures on E .

Next, considering a compact subset $\Lambda \subset E$, we may consider the set $\mathfrak{N}_\Lambda = \{\xi \subset \Lambda, |\xi| < \infty\}$ of finite configurations on Λ , equipped with the trace σ -algebra.

A point process is then any random variable valued in \mathfrak{N} .

Definition 2. Let η be a point process on E . For $n \geq 1$, the n -th factorial moment of η is the point process defined on E^n as the set of n -tuples of points of η :

$$M_n^\eta = \{(x_1, \dots, x_n) \mid x_1, \dots, x_n \in \eta\}.$$

Definition 3. The n -th factorial moment measure of η is the measure defined by

$$\mu_n^\eta(B_1 \times \dots \times B_n) = \mathbb{E}(M_n^\eta(B_1 \times \dots \times B_n)),$$

where B_1, \dots, B_n are measurable subsets of E . This relation defines a unique measure on E^n equipped with the product σ -algebra.

Definition 4. If μ_n^η admits a Radon-Nikodym derivative ρ_n with respect to $\lambda^{\otimes n}$, the former is called n -th correlation function of μ . In other words, we have, assuming their existence

$$\rho_n(x_1, \dots, x_n) = \frac{d(\mu_n^\eta)}{d(\lambda^{\otimes n})}(x_1, \dots, x_n)$$

Equivalently, a point process η is said to have correlation functions $(\rho_n)_{n \geq 1}$ if for all disjoint bounded Borel subsets A_1, \dots, A_n of E , we have

$$\mathbb{E} \left[\prod_{k=1}^n \xi(A_k) \right] = \int_{A_1 \times \dots \times A_n} \rho_n(x_1, \dots, x_n) d\lambda^{\otimes n}.$$

Correlation functions are symmetric and characterized the correlation functions. ρ_1 is interpreted as the average density of particles with respect to λ . More generally,

$$\rho(x_1, \dots, x_n) d\lambda(x_1) \dots d\lambda(x_n),$$

represents the probability of the event « For each $k \in \llbracket 1, n \rrbracket$, there is a (at least one) point at the vicinity dx_k of x_k ». See [4] for an introduction to this topic.

For any compact set $\Lambda \subset E$, we denote by $L^2(\Lambda, \lambda)$ the Hilbert space of complex-valued square integrable functions with respect to the restriction of the Radon measure λ on Λ , equipped with the usual inner-product. Recall that an integral operator $K : L^2(E, \lambda) \rightarrow L^2(E, \lambda)$ is said to have kernel $k : E^2 \rightarrow \mathbb{C}$ if K is written

$$K : f \mapsto \int k(x, y) f(y) d\lambda(y).$$

If k is in $L^2(E^2, \lambda^{\otimes 2})$, the operator K is a linear bounded operator on $L^2(E, \lambda)$.

Definition 5. Let now be a compact subset $\Lambda \subset E$, the *restriction* of the integral operator K to Λ is the integral operator

$$K^\Lambda : f \in L^2(\Lambda, \lambda) \rightarrow \int_\Lambda k(x, y) f(y) d\lambda(y) \in L^2(\Lambda, \lambda),$$

and K^Λ is then a compact operator.

Definition 6. The operator K is said to be Hermitian or self-adjoint if its kernel k verifies $k(x, y) = \overline{k(y, x)}$ for $\lambda^{\otimes 2}$ -almost every $(x, y) \in E^2$.

Equivalently, this means that the integral operators K^Λ are self-adjoint for any compact set $\Lambda \subset E$. If K^Λ is self-adjoint, by the spectral theorem for self-adjoint and compact operators we have that $L^2(\Lambda, \lambda)$ has an orthonormal basis $(\phi_j^\Lambda)_{j \geq 0}$ of eigenfunctions of K^Λ . The corresponding eigenvalues $(\lambda_j^\Lambda)_{j \geq 0}$ have finite multiplicity (except possibly the zero eigenvalue) and the only possible accumulation point of the eigenvalues is zero. In that case, Mercer's theorem indicates that the kernel k^Λ of K^Λ can be written

$$k^\Lambda(x, y) = \sum_{n \geq 0} \lambda_n^\Lambda \phi_n^\Lambda(x) \overline{\phi_n^\Lambda(y)},$$

where the $(\phi_k^\Lambda)_{k \geq 0}$ form a Hilbert basis of $L^2(E, \lambda)$ composed of eigenfunctions of K^Λ .

Recall that K is positive if its spectrum is included in \mathbf{R}^+ , and is of trace-class if

$$\sum_{n=1}^{\infty} |\lambda_n| < \infty,$$

its trace is then $\text{Tr}(K) := \sum_{n=1}^{\infty} \lambda_n$.

Definition 7. If K^Λ is of trace-class for all compacts Λ , K is said to be locally trace-class.

Hypothesis 8. Throughout this paper, our kernels will be self-adjoint, locally trace class, with spectrum contained in $[0, 1]$.

We refer to [2] and [3] for further developments on these notions.

Definition 9. A locally finite and simple point process on E is a determinantal point process if its correlation functions with respect to the reference Radon measure λ on E exist and are of the form

$$\rho_n(x_1, \dots, x_n) = \det(k(x_i, x_j))_{1 \leq i, j \leq n},$$

where k satisfies Hypothesis 8.

The dynamics of DPPs are described by the following fundamental theorem.

Theorem 10. Under the aforementioned assumptions, consider Mercer's decomposition of the kernel k of the determinantal point process η :

$$k(x, y) = \sum_{k=1}^n \lambda_k \phi_k(x) \overline{\phi_k(y)}.$$

Here, the eigenvalues are all in $[0, 1]$ and can all be chosen in $(0, 1]$; n is equal to the rank of K , which can be either finite or infinite. The (ϕ_n) form a Hilbert basis of the space $L^2(E, \lambda)$. Consider then a sequence of independent Bernoulli random variables $(B_k)_{1 \leq k \leq n}$, and consider the random kernel

$$k_B(x, y) = \sum_{k=1}^n B_k \phi_k(x) \overline{\phi_k(y)},$$

Then the point process η_B with (random) kernel k_B has the same law as that of k :

$$k \stackrel{\text{Law}}{=} k_B.$$

Corollary 11. If the n above is finite, then the point process has a finite number of points almost surely. By independence, the Borel-Cantelli lemmas imply that the number of points is almost surely finite if and only if K is of trace class. Since the integral operators that we will consider will all be locally of trace-class, the restrictions of our DPPs to compact subsets will all have a finite number of points almost surely.

For more details, we refer to [10] and [18].

Definition 12. The Bergman determinantal point process is the determinantal point process on the open unit disc centered at the origin of the complex plane with kernel

$$k(x, y) = \frac{1}{\pi} \frac{1}{(1 - x\bar{y})^2}.$$

See [17] and [11] for a thorough study.

The key motivation of this article is the algorithm that was introduced and studied in [7], [8]. It is recalled down below. Its goal is to simulate a DPP restricted to a ball centered at the origin and of radius R . It assumes that the set

$$I = \{n \geq 0, B_n = 1\},$$

of "active" Bernoulli random variables has been computed, and computes the positions of the points $(X_k)_{k \in I}$. Recall that this makes sense because restricting a DPP to a compact reduces the number of points to almost surely finite.

Since simulating a countable infinity of Bernoulli random variables is unfeasible, we introduce the truncation to N points of the restricted DPP, which is defined by its kernel

$$k_N^\Lambda = \sum_{k=0}^{N-1} \lambda_n^\Lambda \phi_n^\Lambda(x) \overline{\phi_n^\Lambda(y)},$$

which is the only one that's numerically feasible.

III The Bergman DPP on a disc

The big question is now to know to what extent this truncation greatly affects the simulation of the DPP law (or if the error is only a minor issue). We provide the theoretical results that answer the last question in the negative for the Bergman DPP (and partially for general DPPs). This is the question that was asked in [6].

Notation. Throughout this paper and unless expressly stated otherwise, R and r denote real numbers in the interval $(0, 1)$ such that $r < R$.

Let us first construct the restriction of the Bergman DPP to a compact ball centered at the origin with radius R .

Proposition 13. Denote $\mathcal{B}(0, R)$ the compact ball centered at 0 with radius R . Mercer's decomposition of the kernel $k^R(x, y)$ of the Bergman determinantal point process restricted to $\mathcal{B}(0, R)$ is

$$k^R(x, y) = \sum_{n \geq 0} \lambda_n^R \phi_n^R(x) \overline{\phi_n^R(y)},$$

Algorithm 1 Sampling of the locations of the points given the set I of active Bernoulli random variables

Input: R, I

Output: $(X_i)_{i \in I}$

Let $\varphi_I^R(x) = (\varphi_i^R(x), i \in I)$

Draw X_1 from the distribution with density $\|\varphi_I(x)\|_{\mathbf{C}^{|I|}}^2/|I|$

$e_1 \leftarrow \frac{\varphi_I^R(X_1)}{\|\varphi_I^R(X_1)\|_{\mathbf{C}^{|I|}}}$

for $i \leftarrow 2$ **to** $|I|$ **do**

Draw X_i from the distribution with density

$$p_i(x) = \frac{1}{|I| - i + 1} \left(\|\varphi_I^R(x)\|_{\mathbf{C}^{|I|}}^2 - \sum_{k=1}^{i-1} |\langle e_k, \varphi_I^R(x) \rangle|^2 \right)$$

$$u_i \leftarrow \varphi_I^R(X_i) - \sum_{k=1}^{i-1} \langle e_k, \varphi_I^R(X_i) \rangle e_k$$

$$e_i \leftarrow \frac{u_i}{\|u_i\|_{\mathbf{C}^{|I|}}}$$

end for

where the eigenvalues are

$$\lambda_k^R = R^{2k+2},$$

and the eigenfunctions

$$\phi_k^R : x \mapsto \sqrt{\frac{k+1}{\pi}} \frac{1}{R^{k+1}} x^k.$$

Proof.

We have, for the (original) Bergman kernel

$$k(x, y) = \frac{1}{\pi} \frac{1}{(1 - x\bar{y})^2},$$

using $\frac{1}{(1-u)^2} = \sum_{k \geq 0} (k+1)u^k$ for all $u \in \mathcal{B}(0, R)$, we have

$$\begin{aligned} \sum_{k \geq 0} \frac{k+1}{\pi} x^k \bar{y}^k &= \sum_{k \geq 0} R^{2k+2} \frac{k+1}{\pi} \frac{x^k}{R^{k+1}} \frac{\bar{y}^k}{R^{k+1}} \\ &= \sum_{k \geq 0} \lambda_n^R \phi_n^R(x) \overline{\phi_n^R(y)}, \end{aligned}$$

taking ϕ_n^R and λ_n^R as in the theorem statement. Because we are restricting onto $L^2(\mathcal{B}(0, R), \lambda)$, we ought to consider the inner product on this space and not $L^2(\mathcal{B}(0, 1), \lambda)$. We have

$$\langle \phi_n^R, \phi_m^R \rangle_{L^2(\mathcal{B}(0, R))} = \frac{\sqrt{(n+1)(m+1)}}{\pi} \frac{1}{R^{n+m+2}} \int_{\mathcal{B}(0, R)} z^n \bar{z}^m dz.$$

However,

$$\int_{\mathcal{B}(0, R)} z^n \bar{z}^m dz = \int_0^R \int_0^{2\pi} r^{n+m} e^{i\theta(n-m)} d\theta dr = 2\pi \delta_{n,m} \int_0^R r^{n+m+1} dr = 2\pi \delta_{n,m} \frac{R^{2n+2}}{2n+2}.$$

Consequently,

$$\langle \phi_n^R, \phi_m^R \rangle_{L^2(\mathcal{B}(0,R))} = \delta_{n,m} \frac{\sqrt{(n+1)(m+1)}}{\pi} \frac{1}{R^{n+m+2}} 2\pi \frac{R^{2n+2}}{2n+2} = \delta_{n,m}.$$

This proves that this family is an orthonormal family of eigenfunctions of the integral operator K^R associated to k^R , and we have hence found its Mercer decomposition. \square

Remark 14. The fact that the eigenvalues and eigenfunctions of the Bergman DPP are explicitly computable makes it very manipulatable. Let us mention that restricting any DPP on any compact rarely ever gives birth to such well-behaving results. The only DPP that was ever discovered to enjoy such properties is the Ginibre DPP as shown in [8].

We will now make use of Optimal transport tools and show that, when choosing the right number of points (which we will exhibit), the law induced by the truncated version of this kernel is close (in the Wasserstein sense) to the non-truncated restriction.

Let X and Y be two Polish spaces. Let μ and ν be probability measures on X and Y respectively. Denote $\Pi(\mu, \nu)$ the set of probability measures on $X \times Y$, the first marginal of which is μ and the second ν . If c is a lower semi-continuous function from $X \times Y$ to \mathbf{R}^+ , the Monge-Kantorovitch problem asks to find

$$\inf_{\gamma \in \Pi(\mu, \nu)} \int_{X \times Y} c(x, y) d\gamma(x, y).$$

We refer to [25] and [26] for proper introductions to the topic of Optimal transport.

Definition 15. Observing that the cardinality of the symmetric difference $d(\xi, \zeta) = |\xi \Delta \zeta|$ induces a distance on the set of configurations on a subset of \mathbf{C} , the Kantorovitch-Rubinstein distance is the Wasserstein-1 distance induced by d , that is

$$\mathcal{W}_{KR}(\mu, \nu) = \inf_{\substack{\text{law}(\xi) = \mu \\ \text{law}(\zeta) = \nu}} \mathbf{E}(|\xi \Delta \zeta|) = \inf_{\substack{\text{law}(\xi) = \mu \\ \text{law}(\zeta) = \nu}} \mathbf{E}(d(\xi, \zeta)).$$

It is a distance on the set of point process laws. See [5] for a thorough study of Wasserstein distances on configuration spaces. The following result exhibits a good compromise as to the number of points the Bergman ought to be truncated to.

Theorem 16. Let

$$N_R := \sum_{n=0}^{\infty} R^{2n+2} = \frac{R^2}{(1-R)(1+R)}$$

Denote \mathfrak{S}^R the law of the restricted Bergman to a compact ball of radius R centered at 0 and \mathfrak{S}_α^R the law of its truncation to α points.

If we truncate it to βN_R points, we have

$$\mathcal{W}_{KR}(\mathfrak{S}^R, \mathfrak{S}_{\beta N_R}^R) \leq N_R e^{-2\beta g(R)} \quad (2)$$

where

$$g(R) = \frac{R^2}{1+R}$$

Proof.

Consider the coupling of $(\xi_{\beta N_R}^R, \xi^R)$ such that the first random variable is a subset of the second one consisting in the point whose indexes are $n \leq \beta N_R$ in Mercer's decomposition. We then have

$$\mathcal{W}_{KR}(\mathfrak{S}^R, \mathfrak{S}_{\beta N_R}^R) \leq \sum_{k=\beta N_R+1}^{\infty} R^{2k+2}.$$

Introducing $\varepsilon = 1 - R$,

$$\begin{aligned} \mathcal{W}_{KR}(\mathfrak{S}^R, \mathfrak{S}_{\beta N_R}^R) &\leq \frac{(1-\varepsilon)^2}{\varepsilon(2-\varepsilon)} (1-\varepsilon)^{2\beta N_R} \\ &\leq \frac{(1-\varepsilon)^2}{\varepsilon(2-\varepsilon)} e^{2\beta N_R \log(1-\varepsilon)} \\ &\leq \frac{(1-\varepsilon)^2}{\varepsilon(2-\varepsilon)} \exp\left(2\beta \frac{(1-\varepsilon)^2}{\varepsilon(2-\varepsilon)} \log(1-\varepsilon)\right) \\ &\leq \frac{(1-\varepsilon)^2}{\varepsilon(2-\varepsilon)} \exp\left(-2\beta \frac{(1-\varepsilon)^2}{2-\varepsilon}\right). \end{aligned}$$

The proof is complete. \square

Remark 17. As a corollary of the previous proof, the two point processes coincide with high probability. The previous proof shows that :

Proposition 18. We have

$$\mathbf{P}(\mathfrak{S}^R \neq \mathfrak{S}_{\beta N_R}^R) \leq N_R e^{-2\beta \frac{R^2}{1+R}}.$$

Proof. Introducing the Bernoulli random variables from Theorem 7, we have

$$\mathbf{P}(\mathfrak{S}^R \neq \mathfrak{S}_{\beta N_R}^R) \leq \mathbf{P}(\exists k > \beta N_R, B_k = 1) \leq \sum_{k > \beta N_R} \mathbf{P}(B_k = 1) = \sum_{k > \beta N_R} \lambda_k^R \leq N_R e^{-2\beta \frac{R^2}{1+R}}$$

as wanted. \square

Remark 19. In [8], the authors have chosen to truncate the restricted Ginibre DPP to $N_R = R^2$ points as seen in result (1) that shows that the truncation error is exponentially small in the deviation c from N_R . Though the authors of [8] have described this as a "well-known *observation*", it is interesting to observe that this R^2 is by no means random and can be theoretically forecasted. This is because turns out to correspond exactly to the expectation of the number of points of the Ginibre DPP. See for yourself : though very interesting to be pointed out, the following proposition is to be found nowhere in the litterature. It motivates to truncate to the expectation and to study the deviation that we have written in Theorem 13.

Proposition 20. The expected number of points $\mathbf{E}[|\mathfrak{S}_R^G|]$ that will come out from the restricted Ginibre DPP to $\mathcal{B}(0, R)$ is exactly R^2 (here, R can be any positive real number).

Proof. According to the fundamental theorem of DPPs (Theorem 7 in this paper), this expectation is $\sum_{n \geq 0} \lambda_n^R$ where λ_n^R denotes the n -th eigenvalue of the restricted Ginibre integral operator.

We have (see [8]) $\lambda_n^R = \frac{\gamma(n+1, R^2)}{n!}$ where γ stands for the incomplete gamma function $\gamma(n, x) = \int_0^x t^{n-1} e^{-t} dt$.

We have

$$\sum_{n=0}^{\infty} \lambda_n^R = \sum_{n=0}^{\infty} \frac{1}{n!} \gamma(n+1, R^2) = \sum_{n=0}^{\infty} \frac{1}{n!} \int_0^{R^2} t^{n+1-1} e^{-t} dt = \int_0^{R^2} \sum_{n=0}^{\infty} \frac{1}{n!} t^n e^{-t} dt = \int_0^{R^2} dt = R^2$$

as wanted. \square

Remark 21. The behavior of the rational function $\frac{R^2}{1+R}$ inside the exponential function is to be nicely bounded in the neighborhood of $R \rightarrow 1$. In other words, the bound (the deviation) is indeed exponential (exponentially small) in β . For instance, for $R \in [0.9, 1]$, we have $g(R) = \frac{R^2}{1+R} \geq 0.42$, so we get $\leq \dots \exp(-0.98\beta)$. For $R \in [0.99, 1]$, $g(R) \geq 0.493$, so $\leq \dots \exp(-0.986\beta)$, and so on. (Since $g(R)$ is continuous and vanishes at 0, it cannot be bounded from below on $[0, 1]$ by a constant $K_0 > 0$. This is why here have to consider subintervals that do not touch 0 and see the bounds on those).

Remark 22. We have $N_R \xrightarrow{R \rightarrow 1^-} +\infty$. This comforts in thinking that the number of points is well-chosen, since as $R \rightarrow 1^-$, we seem to find a restriction that comes closer and closer to the original Bergman point process.

The following results further confirm this idea.

Proposition 23. Denoting \mathfrak{S}^R the law of the restricted Bergman to a compact ball of radius R centered at the origin and \mathfrak{S}_α^R the law of its truncation to α points We have

$$\mathfrak{S}_N^R \xrightarrow{N \rightarrow \infty} \mathfrak{S}^R$$

in distribution.

Proof. The goal is to reduce this result to Theorem 25. Denote $\|\cdot\|_\infty$ the norm of uniform convergence. Let $\Lambda \subset \mathcal{B}(0, R)$ be compact. Denote k_N^R the truncated kernel to N points and k^R the asymptotic one. We have

$$\|k_N^R - k^R\|_{\infty, \Lambda} = \left\| \frac{1}{\pi} \sum_{k=N}^{\infty} (k+1) x^k \bar{y}^k \right\|_{\infty, \Lambda}$$

Observe that $\Lambda \subset \mathcal{B}(0, R)$ implies that

$$\begin{aligned} \|k_{N,R} - k_R\|_{\infty, \Lambda} &\leq \frac{1}{\pi} \sum_{k=N}^{\infty} (k+1) \|x^k \bar{y}^k\|_{\infty, \Lambda} \\ &\leq \frac{1}{\pi} \sum_{k=N}^{\infty} (k+1) R^{2k} \xrightarrow{N \rightarrow \infty} 0 \end{aligned}$$

The proof is complete \square .

Proposition 24. Denoting \mathfrak{S} the original Bergman DPP, we have

$$\mathfrak{S}_{N_R}^R \xrightarrow{R \rightarrow 1^-} \mathfrak{S}$$

in distribution.

Proof.

Denote k_N^R the restricted-truncated (to N points) kernel and k the asymptotic one. Consider a compact $\Lambda \subset D(0, 1)$. Recall that the Bergman kernel $k(x, y) = \frac{1}{\pi} \frac{1}{(1 - x\bar{y})^2}$ is defined on the *open* disc $D(0, 1)$ (this relation would not make sense on some points of the boundary).

Since $D(0, 1)$ is open, denoting $m := \max_{z \in \Lambda} |z|$, we have $m < 1$. So, we have

$$\begin{aligned} \|k_{N_R}^R - k\|_{\infty, \Lambda} &= \left\| \frac{1}{\pi} \sum_{k=N_R}^{\infty} (k+1) x^k \bar{y}^k \right\|_{\infty, \Lambda} \\ &\leq \frac{1}{\pi} \sum_{k=N}^{\infty} (k+1) \|x^k \bar{y}^k\|_{\infty, \Lambda} \\ &\leq \frac{1}{\pi} \sum_{k=N}^{\infty} (k+1) m^{2k} \xrightarrow{N \rightarrow \infty} 0 \end{aligned}$$

We conclude using Theorem 25. □

Theorem 25. (Proposition 3.10, [18])

Let $(K_n)_{n \geq 0}$ be integral operators with nonnegative continuous kernels $k_n(x, y)$. Assume that K_n are bounded hermitian locally trace class integral operators. Assume that $(k_n)_{n \geq 0}$ converges to a kernel k uniformly on each compact as $n \rightarrow \infty$. Then, the kernel k defines an integral operator K that is also a bounded hermitian locally trace class.

Then the determinantal measures, i.e. the DPP probability laws μ_n associated to the DPP induced by the integral operators K_n weakly converge to the determinantal probability measure μ induced by K .

IV The Bergman DPP on an annulus

We now consider a variant of the previous restriction. A sight at Figure 1 in the introduction shows that the points in the Bergman DPP are very highly repelled from the center, and very intensely concentrated at the border of the disc they're living in. The following was shown in [17] :

Theorem 26. The law of the set of the moduli $\{|X_k|, k \geq 1\}$ of the points that come out from the Bergman determinantal point process is exactly the law of the set

$$\{U_k^{1/(2k)}, k \geq 1\}$$

where $(U_k)_{k \geq 1}$ is a sequence of independent, uniform in $[0, 1]$ random variables.

Remark 27. (Conjecture)

According to observations (see Figure 1), we may conjecture that, conditionnally on the set $I = \{n \geq 0, B_n = 1\}$ of active Bernoulli random variables, the law of the set of moduli $\{|X_k|, k \in I\}$ of the points for the *restricted* Bergman determinantal point process is exactly the law of the set

$$\{U_k^{1/(2k)}, k \in I\}$$

where $(U_k)_{k \geq 1}$ is a sequence of independent, uniform in $[0, R]$ random variables.

We then have the following.

Proposition 28. Let $m := \min_{1 \leq k \leq n} (|X_k|)$ denote the smallest radius among the n first points of the Bergan point process (as always, "first" in the sense of Mercer's decomposition).

We have, for all $x \in [0, 1]$

$$\mathbf{P}(m \leq x) = 1 - \prod_{k=1}^n (1 - x^{2k})$$

Proof.

We have

$$\mathbf{P}(U_k^{1/(2k)} \geq x) = \mathbf{P}(U_k \geq x^{2k}) = 1 - x^{2k}$$

According to Theorem 26, we have

$$\mathbf{P}(m \geq x) = \prod_{k=1}^n \mathbf{P}(U_k^{1/(2k)} \geq x) = \prod_{k=1}^n (1 - x^{2k})$$

The proof is complete. □

Remark 29. As a corollary, we have

$$\mathbf{P}(m \leq x) \underset{x \rightarrow 0}{\sim} x^2$$

because the polynomial involved is even and vanishes at zero. This suggests that restricting to a annulus instead of a ball could be enough. Though it is to be mentionned that this has not yet been implemented in [15], we construct this restriction.

Theorem 30. Denote $T(r, R)$ the compact annulus centered at 0 with inner radius r and outer radius R . Mercer's decomposition of the kernel $k_{r,R}(x, y)$ of the Bergman determinantal point process restricted to $T(r, R)$ is

$$k_{r,R}(x, y) = \sum_{n \geq 0} \lambda_n^{r,R} \phi_n^{r,R}(x) \overline{\phi_n^{r,R}(y)}$$

where the eigenvalues are

$$\lambda_k^R = R^{2k+2} - r^{2k+2}$$

and the eigenfunctions

$$\phi_k^R : x \mapsto \sqrt{\frac{k+1}{\pi(R^{2k+2} - r^{2k+1})}} x^k$$

Proof.

We have, for the (original) Bergman kernel

$$k(x, y) = \frac{1}{\pi} \frac{1}{(1 - x\bar{y})^2}$$

using $\frac{1}{(1-u)^2} = \sum_{k \geq 0} (k+1)u^k$ for all $u \in T(r, R)$, we have

$$\sum_{k \geq 0} \frac{k+1}{\pi} x^k \bar{y}^k = \sum_{k \geq 0} \lambda_n^{r,R} \phi_n^{r,R}(x) \overline{\phi_n^{r,R}(y)}$$

taking $\phi_n^{r,R}$ and $\lambda_n^{r,R}$ as in the theorem statement. We have

$$\langle \phi_n^{r,R}, \phi_m^{r,R} \rangle_{L^2(T(r,R))} = \frac{1}{\pi} \sqrt{\frac{(n+1)(m+1)}{\lambda_n^{r,R} \lambda_m^{r,R}}} \int_{T(r,R)} z^n \bar{z}^m dz$$

However,

$$\int_{T(r,R)} z^n \bar{z}^m dz = \int_r^R \int_0^{2\pi} r^{n+m} e^{i\theta(n-m)} d\theta r dr = 2\pi \delta_{n,m} \int_r^R r^{n+m+1} dr = 2\pi \delta_{n,m} \left[\frac{R^{2n+2}}{2n+2} - \frac{r^{2n+2}}{2n+2} \right]$$

Consequently,

$$\langle \phi_n^{r,R}, \phi_m^{r,R} \rangle_{L^2(T(r,R))} = \delta_{n,m}$$

The proof is thus complete. \square

Remark 31. The previous proof actually shows that it is also feasible to compute the restriction to a domain of the form $\{z \in \mathbf{C}, |z| \in A\}$ where A is a Borel subset of $[0, 1]$, yielding eigenvalues

$$\lambda_n^A = \int_A r^{2n+1} dr,$$

with the same corresponding eigenfunctions up to a normalization factor.

V General results

To finish with, we have proven results for the Bergman DPP, the proofs of which used methods that turned out to be very general and apply to general DPPs. We list these results below and group them in this section.

As stated in Comment 16, our approach indicates that truncating restricted DPPs to a number of points equal to the expectation of their cardinality yields strong results such as Theorem 13 and Proposition 15. We here further study the deviation of this cardinality through the two following results.

Theorem 32. Let \mathfrak{S} be a determinantal point process. Assume that its associated integral operator is trace-class.

Denote m the average number of points of \mathfrak{S} . m is finite, and we have, for $c \in]0, 1[$

$$\mathbf{P}(|\mathfrak{S}| \leq (1 \pm c)N_R) \leq e^{-N_R(-2c+3c^2/2-c^3/2)}$$

Remark 33. This is, to some extent, a generalization to Lemma 15 in [8], to any DPP.

Proof.

Denote B_n the n -th Bernoulli random variable (see Theorem 7), we have $B_n \sim \text{Ber}(\lambda_n)$. Denote $S = \sum_{n=0}^{\infty} B_n$.

Note that m is the expectation of S . But K is trace-class. So, m is finite.

$$\begin{aligned} \mathbf{P}(|\mathfrak{S}| \leq (1-c)N_R) &= \mathbf{P}(S \leq (1-c)N_R) = \mathbf{P}\left[\left(\frac{1}{1-c}\right)^S \leq \left(\frac{1}{1-c}\right)^{(1-c)N_R}\right] \\ &= \mathbf{P}\left[(1-c)^{(1-c)N_R} \leq (1-c)^S\right] \leq (1-c)^{-(1-c)N_R} \mathbf{E}[(1-c)^S] \\ &= e^{-(1-c)N_R \log(1-c)} \mathbf{E}\left[e^{\log(1-c) \sum_{n=0}^{\infty} B_n}\right] = e^{-(1-c)N_R \log(1-c)} \prod_{n=0}^{\infty} \mathbf{E}[e^{\log(1-c)B_n}] \\ &= e^{-(1-c)N_R \log(1-c)} \prod_{n=0}^{\infty} (1 - R^{2n+2} + R^{2n+2}(1-c)) = e^{-(1-c)N_R \log(1-c)} \prod_{n=0}^{\infty} (1 - cR^{2n+2}) \\ &\leq e^{-(1-c)N_R \log(1-c)} \prod_{n=0}^{\infty} e^{-cR^{2n+2}} = e^{-(1-c)N_R \log(1-c)} e^{-c \sum_{n=0}^{\infty} R^{2n+2}} \\ &= e^{-(1-c)N_R \log(1-c) - cN_R} = e^{N_R(-c - (1-c)\log(1-c))} \\ &\leq e^{N_R(-c - (1-c)(c - c^2/2))} = e^{N_R(-2c + 3c^2/2 - c^3/2)} \end{aligned}$$

The proof for $1+c$ instead of $1-c$ is almost identical.

Corollary 34. This inequality applies to the restriction to a compact of a DPP with locally trace-class integral operator, in particular to the Bergman restrictions we have studied earlier.

The next result is another information about the aforementioned deviation.

Theorem 35. Let \mathfrak{S}_n be the truncation to n points of a determinantal point process \mathfrak{S} . Denote σ_n^2 the variance of its number of points.

Then we have

$$\frac{|\mathfrak{S}_n| - \mathbf{E}[|\mathfrak{S}_n|]}{\sqrt{n}} \rightarrow \mathcal{N}(0, \sigma_n^2)$$

in distribution.

Proof.

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