A RICE METHOD PROOF OF THE NULL-SPACE PROPERTY OVER THE GRASSMANNIAN

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ABSTRACT. The Null-Space Property (NSP) is a necessary and sufficient condition for the recovery of the largest coefficients of solutions to an underdetermined system of linear equations. Interestingly, this property governs also the success and the failure of recent developments in high-dimensional statistics, signal processing, error-correcting codes and the theory of polytopes.

Although this property is the keystone of ℓ_1 -minimization techniques, it is an open problem to derive a closed form for the phase transition on NSP. In this article, we provide the first proof of NSP using random processes theory and the Rice method. As a matter of fact, our analysis gives non-asymptotic bounds for NSP with respect to unitarily invariant distributions. Furthermore, we derive a simple sufficient condition for NSP.

1. INTRODUCTION

1.1. Null-Space Property. One of the simplest inverse problem can be described as follows: given a matrix $X \in \mathbb{R}^{n \times p}$ and $y \in \text{Im}(X)$, can we faithfully recover β^* such that the identity $y = X\beta^*$ holds? In the ideal case where $n \ge p$ and the matrix X is one to one (namely, the model is identifiable), this problem is elementary. However, in view of recent applications in genetics, signal processing, or medical imaging, the frame of high-dimensional statistics is governed by the opposite situation where n < p. To bypass the limitations due to the lack of identifiability, one usually assumes that the matrix X is at random and one considers the ℓ_1 -minimization procedure [14]:

$$(P_{\ell_1}) \qquad \qquad \Delta_X(\beta^*) \in \arg\min_{X\beta = X\beta^*} \|\beta\|_1,$$

where $\beta^* \in \mathbb{R}^p$ is a "target" vector we aim to recover. Interestingly, Program (P_{ℓ_1}) can be solved efficiently using linear programming, e.g. [11]. Furthermore, the highdimensional models often assume that the target vector β^* belongs to the space Σ_s of *s*-sparse vectors:

$$\Sigma_s := \{ \beta \in \mathbb{R}^p, \ \|\beta\|_0 \le s \},\$$

where $\|\beta\|_0$ denotes the size of the support of β . Note that this framework is the baseline of the flourishing Compressed Sensing (CS), see [10, 19, 15, 13] and references therein. A breakthrough brought by CS states that if the matrix X is drawn at random (e.g. X has i.i.d. standard Gaussian entries) then, with overwhelming probability, one can faithfully recovers $\beta^* \in \Sigma_s$ using (P_{ℓ_1}) . More precisely, the interplay between randomness and ℓ_1 -minimization shows that with only $n = \mathcal{O}(s \log(p/s))$, one can faithfully reconstruct any s-sparse vector β^* from the knowledge of X and $y := X\beta^*$. Notably, this striking fact is governed by the Null-Space Property (NSP).

Date: August 22, 2018.

²⁰¹⁰ Mathematics Subject Classification. 62J05; 62H12; 62F12; 62E20;

Key words and phrases. Rice Method; High-dimensional statistics; ℓ_1 -minimization; Null-Space Property; Random processes theory;

Definition (Null-Space Property of order s and dilatation C) — Let 0 < s < p be two integers and G be a sub-space of \mathbb{R}^p . One says that the sub-space G satisfies NSP(s, C), the Null-Space Property of order s and dilatation $C \ge 1$, if and only if:

$$\forall \gamma \in G, \forall S \subset \{1, \dots, p\} \text{ s.t. } |S| \le s, \quad C \|\gamma_S\|_1 \le \|\gamma_{S^c}\|_1,$$

where S^c denotes the complement of S, the vector γ_S has entry equal to γ_i if $i \in S$ and 0 otherwise, and |S| is the size of the set S.

As a matter of fact, one can prove [15] that the operator Δ_X is the identity on Σ_s if and only if the kernel of X satisfies NSP(s, C) for some C > 1.

Theorem 1 ([15]) — For all $\beta^* \in \Sigma_s$ there is a unique solution to (P_{ℓ_1}) and $\Delta_X(\beta^*) = \beta^*$ if and only if the nullspace ker(X) of the matrix X enjoys NSP(s, C) for some C > 1. Moreover, if ker(X) enjoys NSP(s, C) for some C > 1 then for all $\beta^* \in \mathbb{R}^p$,

$$\|\beta^{\star} - \Delta_X(\beta^{\star})\|_1 \le \frac{2(C+1)}{C-1} \min_{|S| \le s} \|\beta^{\star} - \beta_S^{\star}\|_1.$$

Additionally, NSP suffices to show that any solution to (P_{ℓ_1}) is comparable to the *s*-best approximation of the target vector β^* . Theorem 1 demonstrates that NSP is a natural property that should be required in CS and High-dimensional statistics. This analysis can be lead a step further considering Lasso [33] or Dantzig selector [12]. Indeed, in the frame of noisy observations, ℓ_1 -minimization procedures are based on sufficient conditions like Restricted Isometry Property (RIP) [12], Restricted Eigenvalue Condition (REC) [9], Compatibility Condition (CC) [34], Universal Distortion Property (UDP) [18], or $H_{s,1}$ condition [25]. Note that all of these properties imply that the kernel of the matrix X satisfies NSP. While there exists pleasingly ingenious and simple proofs of RIP, see [13] for instance, a direct proof of NSP (without the use of RIP) remains a challenging issue.

1.2. Contribution. Given $(\rho, \delta) \in]0, 1[^2$, set $s_n = \lfloor \rho n \rfloor$ and $p_n = \lfloor \frac{n}{\delta} \rfloor$ where $\lfloor . \rfloor$ denotes the integer part. Consider a matrix $X(n, p_n) \in \mathbb{R}^{n \times p_n}$ with i.i.d. centered Gaussian entries. In this paper, we describe a region of parameters (ρ, δ) such that $\mathbb{P}[\ker(X(n, p_n)) \text{ enjoys NSP}(s_n, C)]$ tends to one as n goes to infinity. Our result provides a new and simple description of such region of parameters (ρ, δ) .

Theorem 2 — Let $C \ge 1$. For all $n \ge 1$, set $s_n = \lfloor \rho n \rfloor$ and $p_n = \lfloor \frac{n}{\delta} \rfloor$. Let $G(n, p_n)$ be uniformly distributed on the Grassmannian $\operatorname{Gr}_m(\mathbb{R}^{p_n})$ where $m = p_n - n$. If $\delta \ge (1 + \pi/2)^{-1}$ and:

$$\rho \log \left(\sqrt{\frac{\pi}{2eC^2}} \frac{(1-\rho)^2}{\rho^2} \right) + \log \left(Ce \frac{\sqrt{\rho(1-\delta)(1+(C^2-1)\rho)}}{(1-\rho)(1+(2C^2-1)\rho)\sqrt{\delta}} \right) + \frac{1}{\delta} \log \left(\sqrt{\frac{2}{e\pi}} \frac{1+(2C^2-1)\rho}{(1-\rho)\sqrt{\delta(1-\delta)(1+(C^2-1)\rho)}} \right) \le 0$$

then $\mathbb{P}[G(n, p_n) \text{ enjoys } NSP(s_n, C)]$ tends exponentially to one as n goes to infinity.

Remark. The condition $\{\delta \geq (1 + \pi/2)^{-1}\}$ is a technical restriction. Indeed, in the proof, we need to consider an union bound on spheres of decreasing dimension. However, concentration bounds are less efficient on smaller spheres and lead to a limitation of the argument. This is explained into more details further. Interestingly, for C = 1, the transition of Theorem 2 compares to the phase transition of Donoho and Tanner [21], see Figure 2. However, numerically, our lower bound is less interesting when $n \ll p$ so we cannot extract the classical $n \sim c_1 s \log(c_2 p/s)$ result. To underline this fact, we only exhibit a result for $\delta \geq (1 + \pi/2)^{-1} \simeq 0.389$.



FIGURE 1. The phase transition of Theorem 2. From top to bottom, C = 1, 2, 3, 4.



FIGURE 2. The panel illustrates numerically the border of the region described by Theorem 4 (blue line) for which NSP holds ($\Pi \simeq 0$) and the strong phase transition of Donoho and Tanner (green line). Note that the region (ρ, δ) such that $\Pi \simeq 0$, i.e. NSP(s, 1) holds, is located below the curve. Simulations have been performed with n = 200.000.

Exemple 1 — In the case C = 1, we can compare our result (see Theorem 4) given by the set of $(\rho, \delta) \in]0, 1[^2$ such that $\delta \geq (1 + \pi/2)^{-1}$ and such that:

$$\rho \log\left[\sqrt{\frac{\pi}{2e}} \frac{(1-\rho)^2}{\rho^2}\right] + \log\left[e\frac{\sqrt{\rho(1-\delta)}}{(1-\rho)(1+\rho)\sqrt{\delta}}\right] + \frac{1}{\delta}\log\left[\sqrt{\frac{2}{e\pi}} \frac{1+\rho}{(1-\rho)\sqrt{\delta(1-\delta)}}\right]$$

is non-positive to the work of Donoho and Tanner [21] (see Figure 3). Observe that, up to a constant bounded by 2, we recover the initial result.



FIGURE 3. On the left, comparison between the border of the region described by Theorem 2 (blue line) and the strong phase transition of Donoho and Tanner (green line) for $\delta \geq 0.39$. On the right, ratio between the green and the blue line.

We outline that explicit expressions of lower bounds on the phase transition can be found in Section 2.

1.3. Direct proofs of NSP with dilatation C = 1. To the best of our knowledge, all the direct proofs of NSP with dilatation C = 1 are based either on integral convex geometry theory, Gaussian widths, the approximate kinematic formula, or empirical process theory. This section is devoted to a short review of some stateof-the-art results on direct proofs of NSP.

1.3.1. Grassmann angles. In a captivating series of papers [21, 20, 23, 22], Donoho and Tanner have proved that the kernel of a matrix $X(n, p_n) \in \mathbb{R}^{n \times p_n}$ with i.i.d. centered Gaussian entries enjoys a phase transition, i.e. there exists a function ρ_S : $[0, 1[\rightarrow]0, 1[$ such that for all $(\rho, \delta) \in]0, 1[^2,$

$$\lim_{n \to +\infty} \mathbb{P}[\ker(X(n, p_n)) \text{ enjoys } \operatorname{NSP}(s_n, 1)] = \begin{cases} 0 & \text{if } \rho > \rho_S(\delta), \\ 1 & \text{if } \rho < \rho_S(\delta), \end{cases}$$

where we recall that $s_n = \lfloor \rho n \rfloor$ and $p_n = \lfloor \frac{n}{\delta} \rfloor$. Moreover, they have characterized implicitly and computed numerically the function ρ_S (note that the subscript Sstands for "Strong" since ρ_S is often named the "strong threshold"). Observe their approach is based on computation of Grassmann angles of a polytope due to Affentranger and Schneider [3] and Vershik and Sporyshev [35]. Furthermore, note their phase transition is characterized implicitly using an equation involving inverse Mills ratio of the standard normal density. However, they have derived a nice explicit expression of the phase transition for small values of δ , i.e. when $\delta \to 0$. Hence, they uncover that, in the regime $n \ll p$, NSP(s,1) holds when $n \ge Cs \log(\frac{p}{s})$ for nlarge enough.

1.3.2. Gaussian widths. In recent works [31, 32], Stojnic has shown a simple characterization of the sign of the exponent appearing in the expression of the "weak threshold" given by Donoho and Tanner. Note the weak threshold governs the exact reconstruction by ℓ_1 -minimization of s-sparse vectors with prescribed support and signs, while NSP characterizes the exact reconstruction of all s-sparse vectors. In the paper [31], using "Gordon's escape through a mesh" theorem, Stojnic have derived a simpler implicit characterization of the strong threshold ρ_S . As in Donoho and Tanner's work, observe this implicit characterization involves inverse Mill's ratio of the normal distribution and no explicit formulation of ρ_S can be given. Predating Stojnic's work, Rudelson and Vershynin (Theorem 4.1 in [30]) were the first to use "Gordon's escape through the mesh" theorem to derive a nonasymptotic bound on sparse recovery. A similar result can found in the astonishing book of Foucart and Rauhut, see Theorem 9.29 in [24]. Observe that these results hold with a probability at least $1 - \alpha$ and their bounds depend on $\log(\alpha)$ so one needs one more step to derive a lower bound on the strong phase transition. We did not pursue in this direction.

1.3.3. Approximate kinematic formula. In the papers [27, 4], the authors present appealing and rigorous quantitative estimates of weak thresholds appearing in convex optimization, including the location and the width of the transition region. Recall that NSP is characterized by the strong threshold. Nevertheless, the weak threshold describes a region where NSP cannot be satisfied, i.e.

 $\lim_{n \to +\infty} \mathbb{P}[G(n, p_n) \text{ enjoys } \operatorname{NSP}(s_n, 1)] = 0.$

Based on the approximate kinematic formula, the authors have derived recent fine estimates of the weak threshold. Although their result has not been stated for the strong threshold, their work should provide, invoking a simple union bound argument, a direct proof of NSP with dilatation C = 1.

1.3.4. Empirical process theory. Using empirical process theory, Lecué and Mendelson [26] gives a direct proof of NSP for matrices X with sub-exponential rows. Although the authors do not pursue an expression of the strong threshold, their work shows that NSP with dilatation C = 1 holds, with overwhelming probability, when:

(1)
$$n \ge c_0 s \log(\frac{ep}{s}),$$

with $c_0 > 0$ a universal (unknown) constant.

1.3.5. A previous direct proof of NSP with dilatation $C \ge 1$. Using integral convex geometry theory as in Donoho and Tanner's works [21, 20, 23, 22], Xu and Hassibi have investigated [36, 37] the property NSP(s, C) for values $C \ge 1$. Their result uses an implicit equation involving inverse Mill's ratio of the normal distribution and no explicit formulation of their thresholds can be derived. To the best of our knowledge, this is the only proof of NSP(s, C) for values C > 1 predating this paper.

1.4. Simple bounds on the phase transition. As mentioned in Proposition 2.2.17 of [13], if NSP holds then

(2)
$$n \ge c_1 s \log(\frac{c_2 p}{s}),$$

with $c_1, c_2 > 0$ are universal (unknown) constants. The result of Section 1.3.4 shows that a similar bound is also sufficient to get NSP. What can be understood is that the true phase transition (as presented in [21, 20, 23, 22]) lies between the two bounds described by (1) (lower bound) and (2) (upper bound). Observe that these bounds can be equivalently expressed in terms of $\rho = s/n$ and $\delta = n/p$. Indeed, one has:

(3)
$$\{n \ge c_1 s \log(\frac{c_2 p}{s})\} \Leftrightarrow \{A_\star \rho \delta \log(A_\star \rho \delta)) \ge -B_\star \delta\},\$$

where $A_{\star} = c_2^{-1} > 0$ and $1/e \ge B_{\star} = c_1^{-1}c_2^{-1} > 0$. Denote by W₀ (resp. W₋₁) the first (resp. the second) Lambert W function, see [16] for a definition. We deduce that (3) is equivalent to:

(4)
$$\rho \leq \frac{\exp(\mathbb{W}_{-1}(-B_{\star}\delta))}{A_{\star}\delta} \quad \text{or} \quad \rho \geq \frac{\exp(\mathbb{W}_{0}(-B_{\star}\delta))}{A_{\star}\delta}.$$



FIGURE 4. The strong threshold ρ_S and the mapping $\delta \mapsto \frac{\exp(W_{-1}(-0.3394\delta))}{1.38\delta}$.

Furthermore, the papers [21, 20, 23, 22] show that NSP enjoys a phase transition that can be described as a region $\rho \leq \rho_S(\delta)$, see Section 1.3. In particular, one can check that the region described by the right hand term of (4) cannot be a region of solutions of the phase transition problem. We deduce from [13, 26] that ρ_S , the phase transition of Donoho and Tanner [21, 20, 23, 22], can be bounded by the left hand term of (4). Hence, it holds the following result.

Theorem 3 — The strong threshold ρ_S (phase transition of NSP) of Donoho and Tanner [21, 20, 23, 22] is bounded by:

(5)
$$\forall \delta \in]0,1[, \quad \frac{\exp(\mathbb{W}_{-1}(-B_1\delta))}{A_1\delta} \le \rho_S(\delta) \le \frac{\exp(\mathbb{W}_{-1}(-B_2\delta))}{A_2\delta}$$

where $A_1, A_2 > 0$ and $1/e \ge B_1, B_2 > 0$ are universal (unknown) constants.

Although bounds (1) (lower bound) and (2) (upper bound) are known, their expressions as exponential of second Lambert W functions remain overlooked in the literature. As a matter of fact, Figure 4 depicts a comparison between ρ_S and:

(6)
$$\delta \mapsto \frac{\exp(\mathbb{W}_{-1}(-0.3394\delta))}{1.38\delta},$$

where the strong threshold curve has been taken from [21, 20, 23, 22]. Roughly speaking, the curve (6) shows empirically that NSP holds when:

$$n \ge 4s \log(0.7p/s),$$

for large values of s, n, p. Recall that it is still an open problem to find a closed form for the weak and the strong thresholds. In the regime $\delta \to 0$, Donoho and Tanner [21, 20, 23, 22] have proved that the phase transition enjoys

$$n \ge 2es \log(p/(\sqrt{\pi s})) \simeq 5.4s \log(0.6p/s),$$

in the asymptotic.

1.5. Outline of the paper. The main theorem (Theorem 4) is stated in the next section and Section 3 proves it. Section 4 is devoted to the proof of Theorem 2. All the numerical experiments can be reproduced using the codes available at [28].

2. Rice method bound for NSP with dilatation $C \ge 1$

In this paper, we prove NSP following a newt path based on stochastic processes theory and more precisely on the Rice method [7, 8]. This latter is specially design to study the tail of the maximum of differentiable random processes or random fields. Similarly to the case of a deterministic function, it consists of studying the maximum through the zeros of the derivative. For the tail of a stationary Gaussian process defined on the real line, it is known from the work of Piterbarg [29] that it is super-exponentially sharp.

However, the situation here is more involved than in the aforementioned papers since the considered process X(t) is defined on the sphere (as in the recent work [6] for example), non Gaussian and, last but not least, non differentiable. Note that the paper [17] studies the maximum of locally linear process by a smoothing argument. A contrario to this paper, we will use a partition of the sphere and directly the Rice method. This provides a short and direct proof of NSP(s, C) for any value $C \geq 1$.

2.1. An explicit sufficient condition. Our main result reads as follows.

Theorem 4 (Explicit lower bound) — Let 0 < s < n < p and m = p - n. Let G(n,p) be the Kernel of X(n,p), a $(n \times p)$ random matrix with i.i.d. centered Gaussian entries, then for all $C \ge 1$, it holds:

$$\mathbb{P}[G(n,p) \text{ enjoys } NSP(s,C)] = 1 - \Pi,$$

with Π satisfying (7)

$$\Pi \leq \sqrt{\pi} \Big[\sum_{k=0}^{p-n-1} {p \choose k} \Big(\frac{C^2 s}{\tilde{p}_{C,k}} \Big)^{\frac{p-n-1-k}{2}} \frac{\Gamma(\frac{2p-2k-n-1}{2})}{\Gamma(\frac{p-k}{2})\Gamma(\frac{p-n-k}{2})} \psi_{p-k}(C) \mathbf{Q}(k, \tilde{p}_{C,k}, m) \Big],$$

where Γ denotes the Gamma function, $\psi_{p-k}(C)$ is defined by Lemma 4, $\mathbf{Q}(k, \tilde{p}_{C,k}, m)$ is defined by Lemma 2 and $\tilde{p}_{C,k} := (C^2 - 1)s + p - k$.



FIGURE 5. Numerical computation of the lower bound ($\Pi \simeq 0$) on the phase transition for C = 1, 2, 3, 4 (from bottom to top). An explicit expression can be found in Theorem 4. Simulations have been performed with n = 200.000.

3. Proof of Theorem 4

3.1. Model and notation. Let 0 < s < n < p, let C > 1 and set m = p - n. Let G(n, p) be uniformly distributed on the Grassmannian $\operatorname{Gr}_m(\mathbb{R}^p)$. Observe that it can be generated by m independent standard Gaussian vectors $g_i \in \mathbb{R}^p$ for $i = 1, \ldots m$. Define $\{Z(t); t \in \mathbb{S}^{m-1}\}$ the process with values in \mathbb{R}^p given by:

$$Z(t) := \sum_{i=1}^m t_i g_i \,.$$

Note this process spans G(n, p) and it can be written as

for
$$j = 1, \ldots, p$$
, $Z_j(t) = \langle t, g^j \rangle$,

where $(g^j)_{j=1}^p$ are independent Gaussian random vectors with standard distribution in \mathbb{R}^m . Let O_p and O_m two orthogonal matrices of size, respectively, $(p \times p)$ and $(m \times m)$. Thanks to unitarily invariance of the Gaussian distribution, remark that:

$$\forall t \in \mathbb{S}^{m-1}, \quad O_p Z(t) O_m \sim Z(t).$$

Consider now the ordered statistics of the absolute values of the coordinates of Z(t):

$$|Z_{(1)}(t)| \ge \cdots \ge |Z_{(p)}(t)|,$$

where the ordering $((1), \ldots, (p))$ is always uniquely defined if we adopt the convention of keeping the natural order in case of ties. Given a sparsity s, a degree of freedom m, and a degree of constraint p, consider the real valued process $\{X(t); t \in \mathbb{S}^{m-1}\}$ such that:

(8)
$$X(t) = C|Z_{(1)}(t)| + \dots + C|Z_{(s)}(t)| - [|Z_{(s+1)}(t)| + \dots + |Z_{(p)}(t)|].$$

NSP is equivalent to the fact that this process is always non positive. We will prove that it happens with an overwhelming probability.

3.2. Cutting the sphere out. As we will see later, the process X(.) is locally linear over some subsets and to take benefit of that, we need to consider a particular partition of the sphere.

Let $A \subseteq \{1, \ldots, p\}$, define the random subsets S_A and \dot{S}_A of the unit sphere S^{m-1} by:

$$\begin{split} & \$_A = \{t \in \$^{m-1}; \ Z_A(t) = 0\}, \\ & \mathring{\$}_A = \{t \in \$^{m-1}; \ Z_A(t) = 0 \text{ and } \forall j \notin A, \ Z_j(t) \neq 0\}, \end{split}$$

where we denote $Z_A(t) = (Z_j(t))_{j \in A}$. One can check that S_A is the unit sphere of the orthogonal of $V_A := \text{Span}\{g^j; j \in A\}$. This implies that a.s. S_A is a random sphere of dimension m - 1 - |A| if $|A| \leq m - 1$ and is almost surely empty if |A| > m - 1. It follows that the quantities $|Z_{(1)}(t)|, \ldots, |Z_{(n+1)}(t)|$ are a.s. positive and that a.s.

$$\mathbb{S}^{m-1} = \bigcup_{|A| \le m-1} \dot{\mathbb{S}}_A \,,$$

giving a partition of the sphere. We define also, for later use, the random subset ${\mathcal W}$ by:

$$\mathcal{W} := \{ t \in \mathbb{S}^{m-1} ; |Z_{(s)}(t)| = |Z_{(s+1)}(t)| \}.$$

Observe that, conditionally to g^j , the set \mathcal{W} is closed with empty interior.

3.3. Probability of failure. We consider the probability:

(9)
$$\Pi = \mathbb{P}\left\{\mathcal{M} > 0\right\} \le \sum_{|A| \le m-1} \mathbb{P}\left\{\mathcal{M}_{\dot{\mathbb{S}}_A} > 0\right\},$$

where \mathcal{M} and $\mathcal{M}_{\dot{\mathbf{S}}_A}$ are respectively the number of positive local maximum of X(.)along \mathbb{S}^{m-1} and $\dot{\mathbf{S}}_A$. The baseline of our proof is to upper-bound each right hand side probabilities, using the expected number of positive local maximum above zero and Markov inequality. The first element is Lemma 4 proving that:

$$\forall t \in \mathbb{S}_A, \quad \mathbb{P}\{X(t) > 0\} \le \psi_{p-k}(C),$$

where k := |A| and:

$$\psi_{p-k}(C) = \binom{p-k}{s} \left(\frac{C^2 4s}{\pi}\right)^{\frac{p-k-s}{2}} \frac{\Gamma((p-k)/2)}{\Gamma(s/2)\Gamma(p-k-s+1)},$$

where Γ denotes the Gamma function. The second element is that X(t) admits a density $p_{X(t)}$. To check that, note that $|Z_{(1)}(t)|, \ldots |Z_{(p)}(t)|$ are the order statistics of the absolute values of i.i.d. Gaussian variables and thus they have a known joint density on the simplex $|Z_{(1)}| \geq \ldots \geq |Z_{(p)}|$. Formula (8) implies the existence of a density for X(t). Moreover, this density does not depend on t due to invariance of Gaussian distribution.

3.4. Initialization: local maxima on \dot{S}_{\emptyset} . By considering the symmetry properties of the sphere \dot{S}_{\emptyset} , we have:

$$\mathbb{P}\left\{\mathcal{M}_{\dot{\mathbb{S}}_{\emptyset}} > 0\right\} \leq \frac{1}{2}\mathbb{E}(\mathcal{M}_{\dot{\mathbb{S}}_{\emptyset}})\,.$$

In this part, our aim will be to give bound to the expectation using a Kac-Rice formula. One can check that if t belongs to \dot{S}_{\emptyset} and does not belong to \mathcal{W} , X(.) is locally the sum of the absolute values of some s coordinates multiplied by C minus the sum of the absolute values of the other coordinates. It can be written as:

$$X(u) = C\varepsilon_1 Z_{(1)}(u) + \dots + C\varepsilon_s Z_{(s)}(u) + \varepsilon_{s+1} Z_{(s+1)}(u) + \dots + \varepsilon_p Z_{(p)}(u),$$

where $\varepsilon_1, ..., \varepsilon_p$ are random variables taking values ± 1 .

Lemma 1 — Let $t \in \mathbb{S}^{m-1}$ then, almost surely, it holds $t \in \dot{\mathbb{S}}_{\emptyset}$ and $t \notin \mathcal{W}$. Furthermore, the spherical gradient X'(t) and the spherical Hessian X''(t) of X(.) along \mathbb{S}^{m-1} at t exist and:

- $X''(t) = -X(t)I_{m-1}$.
- (X(t), X''(t)) and X'(t) are independent.
- X'(t) has a Gaussian centered isotropic distribution onto t^{\perp} with variance $\tilde{p}_{C,0} = (sC^2 + (p-s))$.

Proof. The fact that, with probability 1, $t \in \hat{\mathbb{S}}_{\emptyset}$ and $t \notin \mathcal{W}$ implies that the process X(.) is locally given by

$$X(u) = C\varepsilon_1 Z_{(1)}(u) + \dots + C\varepsilon_s Z_{(s)}(u) + \varepsilon_{s+1} Z_{(s+1)}(u) + \dots + \varepsilon_p Z_{(p)}(u),$$

where the signs $(\varepsilon_1, ..., \varepsilon_p)$ and the ordering (1), ..., (p) are those of t. The process X(.) is locally linear and thus differentiable around t and its gradient in \mathbb{R}^m at t, denoted $\dot{X}(t)$, is given by

$$C\varepsilon_1 g^{(1)} + \dots + C\varepsilon_s g^{(s)} + \varepsilon_{s+1} g^{(s+1)} + \dots + \varepsilon_p g^{(p)}$$

Moreover, note that its Hessian on \mathbb{R}^m vanishes.

Let us consider now the spherical gradient X'(t) and the spherical Hessian X''(t). It is well known that $X'(t) = P_{t^{\perp}} \dot{X}(t)$, where $P_{t^{\perp}}$ is the orthogonal projection onto the orthogonal of t. As for the spherical Hessian, it is defined on the tangent space t^{\perp} and is equal to the projection of the Hessian in \mathbb{R}^m , which vanishes, minus the product of the normal derivative by the identity matrix. This is detailed in Lemma 5. In the case of the unit sphere, the vector normal to the sphere at t is t itself and

$$X''(t) = -\langle X(t), t \rangle I_{m-1} = -X(t)I_{m-1}.$$

In the case of X'(t), remark that Z(t) and thus X(t), $(\varepsilon_1, \ldots, \varepsilon_p, (1), \ldots, (p))$ are functions of $(P_t(g^1), \ldots, P_t(g^p)) = (Z_1(t)t, \ldots, Z_p(t)t)$ (with obvious notation). They are therefore independent of X'(t) which is a function of $(P_{t\perp}(g^1), \ldots, P_{t\perp}(g^p))$. Conditionally to $(\varepsilon_1, \ldots, \varepsilon_p, (1), \ldots, (p))$, X'(t) can be written as

$$X'(t) = (C\varepsilon_1 P_{t^{\perp}} g^{(1)} + \dots + C\varepsilon_s P_{t^{\perp}} g^{(s)} + \varepsilon_{s+1} P_{t^{\perp}} g^{(s+1)} + \dots + \varepsilon_p P_{t^{\perp}} g^{(p)}),$$

which implies that the conditional distribution of X'(t) is Gaussian with variancecovariance matrix $(sC^2 + (p - s))\mathrm{Id}_{t^{\perp}}$, where $\mathrm{Id}_{t^{\perp}}$ is the identity operator on t^{\perp} . Since X'(t) is independent of $(\varepsilon_1, \ldots, \varepsilon_p, (1), \ldots, (p))$ this conditional distribution is in fact equal to the unconditional distribution.

The next step is to prove that a.s. there is no local maximum on \mathcal{W} . The case where there are tied among the $|Z_i(t)|$ has to be considered (though it happens with probability 0 for a fixed t). Note that the order statistics and the ordering remain uniquely defined because of our convention.

Suppose that $t \in \mathcal{W}$. Since all the possible ordering $((1), \ldots, (p))$ and signs $(\varepsilon_1, \ldots, \varepsilon_p)$ play the same role by unitarily invariance of the distribution of Z(t) for all t, we make the proof in the particular case where $((1), \ldots, (p))$ is the identity and all the signs $(\varepsilon_1, \ldots, \varepsilon_p)$ are positive:

$$Z_1(t) \ge \dots \ge Z_{s-h-1}(t) > Z_{s-h}(t) = \dots = Z_{s+k}(t) > Z_{s+k+1}(t) \ge \dots \ge Z_p(t) > 0.$$

Then, for w in some neighborhood N of t (not included in \mathcal{W}), we have:

$$X(w) = CZ_1(w) + \dots + CZ_{s-h-1}(w) + (1+C) \operatorname{Max}_h (Z_{s-h}(w) + \dots + Z_{s+k}(w)) - (Z_{s-h}(w) + \dots + Z_p(w)),$$

where Max_h is the sum of the *h* largest element of its (h + k + 1) arguments. As being the maximum of $\binom{h}{(s+k)+1}$ linear forms the function Max_h is convex.

Let us consider in detail the vectors g^{s-h}, \ldots, g^{s+k} . With probability 1, they are pairwise different. The point t is chosen such that their projection on t coincide. As a consequence the derivatives of the linear forms $Z_{\ell}(w) = \langle g^{\ell}, w \rangle$, $\ell = (s - h) \ldots (s + k)$ on the tangent space t^{\perp} are pairwise different. This implies that the function Max_h has some direction in which it is strictly convex and as a consequence t cannot be a local maximum.

Suppose that $t \notin W$, and suppose that we limit our attention to points t such that X(t) > 0, then Lemma 1 implies that X''(t) cannot be singular.

This last condition implies that we can apply Theorem 5.1.1 of [2]. This lemma is a Kac type formula that shows that the zeros of the derivative X'(t) are isolated an thus in finite number. In addition recalling that $\mathcal{M}_{\dot{S}_{\emptyset}}$ is the number of positive local maximum of X(.) and belonging to \dot{S}_{\emptyset} , this number satisfies

$$\mathcal{M}(\dot{\mathbb{S}}_{\emptyset}) = \lim_{\delta \to 0} \frac{1}{V(\delta)} \int_{\mathbb{S}^{m-1}} \mathbb{E}(|\det X''(t)| \mathbb{1}_{|X'(t)-0| < \delta} \mathbb{1}_{t \in \dot{\mathbb{S}}_{\emptyset}} \mathbb{1}_{X(t)>0}) \sigma(\mathrm{d}t),$$

where σ is the surfacic measure on \mathbb{S}^{m-1} and $V(\delta)$ is the volume of the ball $B(\delta)$ with radius δ . Passing to the limit using the Fatou lemma gives:

$$\mathbb{E}(\mathcal{M}(\dot{\mathbf{S}}_{\emptyset})) \leq \liminf_{\delta \to 0} \int_{0}^{\infty} \mathrm{d}x \int_{\dot{\mathbf{S}}_{\emptyset}} \mathrm{d}t \, p_{X(t)}(x)$$

$$\frac{1}{V(\delta)} \int_{B(\delta)} \mathrm{d}x' p_{X'(t)}(x') \mathbb{E}\left(|\det(X''(t))| \left| X(t) = x, X'(t) = x'\right) \right.$$

$$\leq \left(2\pi \tilde{p}_{C,0}\right)^{\frac{1-m}{2}} 2 \frac{\pi^{\frac{m}{2}}}{\Gamma(\frac{m}{2})} \int_{0}^{\infty} x^{m-1} p_{X(t)}(x) dx \,,$$

where $p_{X(t)}(x)$ denotes the density of X(t) at x and Γ denotes the Gamma function. Note that we have used:

- the fact that every point t is equivalent so we can replace the integral on the unit sphere by the volume of the unit sphere $2\pi^{\frac{m}{2}}/\Gamma(\frac{m}{2})$ and the value at a given point,
- $\mathbb{E}(|\det(X''(t))| | X(t) = x, X'(t) = x') = x^{m-1},$
- the Gaussian density $p_{X'(t)}(x')$ is bounded by $(2\pi \tilde{p}_{C,0})^{\frac{1-m}{2}}$.

So it remains to bound $\mathbb{E}[(X(t)^+)^{m-1}]$. For that purpose we write X(t) as the independent product $||Z(t)||_2 Y(t)$, where the process Y(t) is constructed exactly as the process X(t) but starting now from a uniform distribution U on the unit sphere \mathbb{S}^{p-1} instead of the standard Gaussian distribution of Z(t). Using standard results on the moments of the χ^2 distribution we have:

$$\mathbb{E}((X(t)^+)^{m-1}) = 2^{\frac{m-1}{2}} \frac{\Gamma(\frac{m-1+p}{2})}{\Gamma(\frac{p}{2})} \mathbb{E}((Y(t)^+)^{m-1}).$$

We use now the fact that $Y(t) \leq C\sqrt{s}$ to get that:

$$\mathbb{E}((Y(t)^+)^{m-1}) \le (C\sqrt{s})^{m-1} \mathbb{P}\{Y(t) > 0\}.$$

Moreover, Lemma 4 shows that, with probability greater than $1 - \psi_p(C)$, a standard Gaussian vector g in \mathbb{R}^p enjoys:

$$C \|g_S\|_1 \le \|g_{S^c}\|_1$$

This implies that:

(10)
$$\mathbb{P}\{Y(t) > 0\} \le \psi_p(C)$$

and consequently the probability of having a local maximum above 0 on $\dot{\mathbb{S}}_{\emptyset}$ is bounded by:

$$\sigma(\dot{\$}_{\emptyset}) (2\pi\tilde{p}_{0})^{\frac{1-m}{2}} 2^{\frac{m-1}{2}} \frac{\Gamma(\frac{m-1+p}{2})}{\Gamma(\frac{p}{2})} (C\sqrt{s})^{m-1} \psi_{p}(C) \le 2\sqrt{\pi} \left(\frac{C^{2}s}{\tilde{p}_{C,0}}\right)^{\frac{m-1}{2}} \frac{\Gamma(\frac{m-1+p}{2})}{\Gamma(\frac{p}{2})\Gamma(\frac{m}{2})} \psi_{p}(C)$$

Denote the right hand side of this last inequality by $h_C(s, m, p)$.

3.5. Maximum on smaller spheres. Let us now consider the case of a maximum on \dot{S}_A , $A \neq \emptyset$. A point $t \in \dot{S}_A \setminus W$ is a local maximum on S^{m-1} if it satisfies the following conditions:

- it is a local maximum along S_A ,
- its super-gradient along the orthogonal space V_A contains zero,

where the super-gradient is defined as the opposite of the sub-gradient. One can easily check that the two conditions are independent. Indeed, recall that k = |A| and $V_A = \text{Span}\{g^i; i \in A\}$ (see Section 3.2) and consider the process X(.) conditionally to V_A . In that case, S_A becomes a deterministic sphere of dimension m - k - 1. Moreover, note that the behavior of X(.) on \mathbb{S}_A depends only on the $g^j, j \notin A$ and that for such j,

$$Z_j(t) = \langle g^j, t \rangle = \langle \Pi_{V_{\mathbf{A}}^{\perp}} g^j, t \rangle,$$

so, conditionally to V_A , the distribution of X(t) corresponds to the case \mathbb{S}_{\emptyset} in the space of dimension m - k instead of m and with p - k vectors. In conclusion, the first condition leads to the same computations as the case \mathbb{S}_{\emptyset} and is bounded by

$$h_C(s, m-k, p-k) = 2\sqrt{\pi} \left(\frac{C^2 s}{\tilde{p}_{C,k}}\right)^{\frac{p-n-1-k}{2}} \frac{\Gamma(\frac{2p-2k-n-1}{2})}{\Gamma(\frac{p-k}{2})\Gamma(\frac{p-n-k}{2})} \psi_{p-k}(C) \,.$$

Let us look to the second one which depends only on the $g^j, j \in A$. Thus we have to compute the probability of the super-gradient to contain zero. Indeed, locally around t, the behavior of X(w) along V_A is the sum of some linear forms (for $j \notin A$) and of absolute value of linear forms (for $j \in A$) thus it is locally concave and we can define its super-gradient. More precisely, for w in a neighborhood of $t \in \hat{S}_A \setminus W$,

$$X(w) = X_A(w) + X_{A^c}(w) \,,$$

where, because $k \leq p - s$:

$$X_A(w) = -\sum_{i \in A} |Z_i(w)|.$$

Around t, $X_{A^c}(w)$ is differentiable and, with a possible harmless change of sign (see Lemma 1), its gradient is given by:

$$\sum_{i\in A^c} C_i g^i\,,$$

where the coefficient C_i takes the value C for s of them and -1 for the others. This gradient is distributed as an isotropic normal variable $\xi \in V_A$ with variance:

$$\tilde{p}_{C,k} = (C^2 - 1)s + p - k$$

By this we mean that the distribution of ξ , in a convenient basis, is $\mathcal{N}(0, \tilde{p}_{C,k}I_k)$. Let us now consider the case $i \in A$. Observe that the super-gradient along V_A of the concave function $-|Z_i(t)|$ at point t is the segment $[-g^i, g^i]$ and thus the super-gradient of $X_A(t)$ is the zonotope:

(12)
$$Zo = \sum_{i \in A} [-g^i, g^i]$$

where the sum denotes the Minkowsky addition. Recall that the distribution of X(t) does not depend on t.

In conclusion, the probability of the super-gradient to contain zero is equal to $\mathbf{P}(k, \tilde{p}_{C,k}, m)$ the probability of the following event:

- draw k standard Gaussian variables g^1, \ldots, g^k in \mathbb{R}^m and consider the zonotope Zo given by formula (12),
- draw in the space V_A generated by g^1, \ldots, g^k an independent isotropic normal variable ξ of variance $\tilde{p}_{C,k}$,
- define $\mathbf{P}(k, \tilde{p}_{C,k}, m)$ as the probability of ξ to be in Zo.

Lemma 2 — Define the orthonormal basis e^1, \ldots, e^k obtained by Gram-Schmidt orthogonalization of the vectors g^1, \ldots, g^k . Then:

(a) $\mathbf{P}(k, \tilde{p}_{C,k}, m)$ is less than the probability $\mathbf{Q}(k, \tilde{p}_{C,k}, m)$ of ξ to be in the hyper-rectangle:

$$R = \sum_{i \in A} [-\langle e^i, g^i \rangle e^i, \langle e^i, g^i \rangle e^i],$$



FIGURE 6. The standard Gaussian measure of the zonotope (in blue) is smaller than that of the rectangle (in red) with basis $\tilde{Z}o_{k-1}$ (in green).

(b) this last probability satisfies:

$$\left(\mathbf{Q}(k,\tilde{p}_{C,k},m)\right)^2 \le \left(\frac{2}{\pi\tilde{p}_{C,k}}\right)^{H_k+k-m} \frac{H_k!}{(m-k)!}$$

with
$$H_k = \lfloor (\frac{\pi}{2} \tilde{p}_{C,k}) \wedge m \rfloor$$
, where $\lfloor . \rfloor$ is the integer part.

Proof. (a) We prove the result conditionally to the g^{i} 's and by induction on k. When k = 1 the result is trivial since the zonotope and the rectangle are simply the same segment.

Let φ_h be the standard Gaussian distribution on \mathbb{R}^h , $\mathbf{P}(k, \tilde{p}_{C,k}, m)$ is equal to:

$$\varphi_k((\widetilde{p}_{C,k})^{-1/2}.Zo) =: \varphi_k(\widetilde{Zo})$$

Via Gram-Schmidt ortogonalisation at step k, we can compute this probability using the Fubini theorem:

$$\mathbf{P}(k, \tilde{p}_{C,k}, m) = \int_{-\frac{\langle e^k, g^k \rangle}{\sqrt{\tilde{p}_{C,k}}}}^{\frac{\langle e^k, g^k \rangle}{\sqrt{\tilde{p}_{C,k}}}} \varphi_{k-1} \big(\widetilde{Zo}_{k-1} + vz \big) \varphi(z) dz,$$

where φ is the standard Gaussian density on \mathbb{R} , \widetilde{Zo}_{k-1} is the zonotope generated by g^1, \ldots, g^{k-1} and normalized by $(\widetilde{p}_{C,k})^{-1/2}$ and v is some vector in \mathbb{R}^m . By use of the Anderson inequality [5], the non-centered zonotope $(\widetilde{Zo}_{k-1} + vz)$ has a smaller standard Gaussian measure than the centered one so

$$\begin{aligned} \mathbf{P}(k, \tilde{p}_{C,k}, m) &\leq \int_{-\frac{\langle e^k, g^k \rangle}{\sqrt{\tilde{p}_{C,k}}}}^{\frac{\langle e^k, g^k \rangle}{\sqrt{\tilde{p}_{C,k}}}} \varphi_{k-1} \big(\widetilde{Zo}_{k-1} \big) \varphi(z) dz \\ &\leq \int_{R} \varphi(z_1) \dots \varphi(z_k) dz_1 \dots dz_k =: \mathbf{Q}(k, \tilde{p}_{C,k}, m). \end{aligned}$$

The last inequality is due to the induction hypothesis. It achieves the proof.

(b) We use the relation above and deconditioning on the g^i . Note the dimension of the edges of the rectangle R are independent with distribution:

$$2\chi(m), 2\chi(m-1), \ldots, 2\chi(m-k+1),$$

where the law $\chi(d)$ is defined as the square root of a $\chi^2(d)$. As a consequence, using the independence of the components of ξ in the basis e^1, \ldots, e^k and the fact that a

Student density T is uniformly bounded by $(2\pi)^{-1/2}$, we get that:

$$\mathbf{Q}(k, \tilde{p}_{C,k}, m) = \mathbb{P}(\xi \in R) = \prod_{\ell=0}^{k-1} \mathbb{P}\left[|T(m-\ell)| \le \sqrt{\frac{m-\ell}{\tilde{p}_{C,k}}} \right]$$
$$= \prod_{\ell=m-k+1}^{m} \mathbb{P}\left[|T(\ell)| \le \sqrt{\frac{\ell}{\tilde{p}_{C,k}}} \right].$$

Suppose that $\pi \tilde{p}_{C,k} \geq 2m$, then a convenient bound is obtained by using the fact that a Student density is uniformly bounded by $(2\pi)^{-1/2}$:

$$\left(\mathbf{Q}(k,\tilde{p}_{C,k},m)\right)^2 \le \left(\frac{2}{\pi\tilde{p}_{C,k}}\right)^k \frac{m!}{(m-k)!}.$$

In the other case, set $H_k = \lfloor (\pi \tilde{p}_{C,k})/2 \rfloor$, where $\lfloor . \rfloor$ is the integer part. Observe that $H_k > m - k + 1$ for $k \ge 1$ to remove factors that are greater than 1 in the computation and obtain

$$\left(\mathbf{Q}(k,\tilde{p}_{C,k},m)\right)^2 \le \left(\frac{2}{\pi\tilde{p}_{C,k}}\right)^{H_k+k-m} \frac{H_k!}{(m-k)!},$$

which conclude the proof.

Eventually, summing up over the $\binom{p}{k}$ sets of size k, we get Theorem 4.

4. INFLUENCE OF SMALLER SPHERES

4.1. General bound on the sum. In this part, we simplify the general bound of Theorem 4 to derive a simpler one, exponentially decreasing in n, as presented in Theorem 2. Considering (7), we have:

(13)
$$\Pi \le \sqrt{\pi} \sum_{k=1}^{m} B_k(s, n, p)$$

where:

$$B_k(s,n,p) = \binom{p}{n+k} \binom{n+k}{s} \left(\frac{C^2 s}{\tilde{p}_{m-k}}\right)^{\frac{k+1}{2}} \frac{\Gamma(\frac{n}{2}+k-\frac{1}{2})}{\Gamma(\frac{n+k}{2})\Gamma(\frac{k}{2})} \times \left(\frac{C^2 4s}{\pi}\right)^{\frac{n+k-s}{2}} \frac{\Gamma(\frac{n+k}{2})\sqrt{H_{m-k}!}}{\Gamma(\frac{s}{2})(n+k-s)!\sqrt{k!}} \left(\frac{2}{\pi \tilde{p}_{C,m-k}}\right)^{\frac{H_{m-k}-k}{2}}$$

In order to derive a lower bound (the aforementioned bound goes exponentially fast towards zero), we limit our attention to the case described by

- $\begin{array}{l} \bullet \ \ ({\rm H1}) \ \rho \leq 1/2, \\ \bullet \ \ ({\rm H2}) \ \frac{1}{\delta} \leq 1+\pi/2(1+\rho(C^2-1)). \end{array}$

Observe that (H1) is not a restriction since we know that NSP does not hold for $\rho \geq 0.2$. Under (H2), note that $\forall k, H_k = m$, and hence

$$\begin{split} \Pi &\leq \mathbf{R}(s,n,p) \frac{p!}{s!\sqrt{m!}\Gamma(s/2)} \left(\frac{4C^2s}{\pi}\right)^{\frac{n-s}{2}} \sum_{i=1}^m \binom{m}{k} \frac{(2C^2s)^k \Gamma(\frac{n}{2}+k)}{(n-s+k)!^2} \left(\frac{2}{\pi \tilde{p}_{C,m-k}}\right)^{\frac{m}{2}} \\ &\leq \mathbf{R}(s,n,p) \frac{p!}{s!(n-s)!\sqrt{m!}\Gamma(s/2)} \left(\frac{4C^2s}{\pi}\right)^{\frac{n-s}{2}} \\ &\qquad \times \sum_{i=1}^m \binom{m}{k} \frac{(2C^2s)^k \Gamma(\frac{n}{2}+k)}{(n-s)^k(n-s+k)!} \left(\frac{2}{\pi \tilde{p}_{C,m-k}}\right)^{\frac{m}{2}}, \end{split}$$



FIGURE 7. Comparison between the phase transition of Theorem 2 and the numerical approximation given by Theorem 4 ($\Pi \simeq 0$ with n = 200.000, dashed line). From top to bottom, C = 1, 2, 3, 4.

where $\mathbf{R}(s, n, p)$ is a polynomial term in (s, n, p). Consider now the quantity

$$\alpha(k) := \frac{\Gamma(\frac{n}{2} + k)}{(n - s + k)!}$$

which is a decreasing function of k under assumption (H1) and the fact that $\tilde{p}_{C,k}$ is an increasing function of k, then we obtain

$$\Pi \le \mathbf{R}(s,n,p) \frac{p! \mathbf{\Gamma}(n/2)}{s!(n-s)!^2 \sqrt{m!} \mathbf{\Gamma}(s/2)} \left(\frac{4C^2 s}{\pi}\right)^{\frac{n-s}{2}} \sum_{i=1}^m \binom{m}{k} \frac{(2C^2 s)^k}{(n-s)^k} \left(\frac{2}{\pi \tilde{p}_{C,m-1}}\right)^{\frac{m}{2}}.$$

At last, using Stirling Formula (see Lemma 3), it yields

$$\Pi \leq \mathbf{R}(s,n,p) \frac{p! \mathbf{\Gamma}(n/2)}{s!(n-s)!^2 \mathbf{\Gamma}(s/2)} \left(\frac{4C^2 s}{\pi}\right)^{\frac{n-s}{2}} \left(\frac{2e}{\pi(n+(C^2-1)s)m}\right)^{\frac{m}{2}} \left(1+\frac{2C^2 s}{n-s}\right)^m.$$

fathering the piece, one has:

$$\Pi \leq \mathbf{R}(s,n,p) \left(\sqrt{\frac{\pi}{2eC^2}} \frac{(n-s)^2}{s^2}\right)^s \left(Ce \frac{\sqrt{nsm(n+(C^2-1)s)}}{(n-s)(n+(2C^2-1)s)}\right)^n \\ \times \left(\sqrt{\frac{2}{e\pi}} \frac{p(n+(2C^2-1)s)}{(n-s)\sqrt{m(n+(C^2-1)s)}}\right)^p,$$

which gives the result of Theorem 2.

Remark. The upper bound on $\alpha(k)$ and the lower bound on $\tilde{p}_{C,k}$ may seem weak but they do not change the result on the phase transition because first terms of the sum give the right order on ρ and δ . To ensure that, see figure 7, which compare the numerical bound with the one of Theorem 2.

APPENDIX A. STIRLING'S FORMULA

Lemma 3 — Let z > 0 then there exists $\theta \in (0, 1)$ such that:

$$\mathbf{\Gamma}(z+1) = (2\pi z)^{\frac{1}{2}} \left(\frac{z}{e}\right)^z \exp(\frac{\theta}{12z}).$$

In particular, if z > 1/12,

$$\left(\frac{z}{e}\right)^z \le \Gamma(z+1) \le \sqrt{2\pi z} \left(\frac{z}{e}\right)^z$$

Proof. See [1] Eq. 6.1.38.

Appendix B. Concentration $\psi_l(C)$

Lemma 4 — Let $C \ge 1$, then, except with a probability smaller than:

$$\psi_l(C) := \binom{l}{s} \left(\frac{C^2 4s}{\pi}\right)^{\frac{l-s}{2}} \frac{\Gamma(l/2)}{\Gamma(s/2)\Gamma(l-s+1)},$$

a standard Gaussian vector $g \in \mathbb{R}^l$ enjoys for all $S \subset \{1, \ldots, l\}, |S| \leq s$,

$$C \|g_S\|_1 \le \|g_{S^c}\|_1.$$

Proof. Let $\xi_C := \{v \in \mathbb{R}^l \text{ that does not satisfy } NSP(s, C)\}$ and consider the joint law of standard Gaussian ordered statistics $(W_{(1)}, \ldots, W_{(l)})$, then

$$\begin{split} \mathbb{P}(\xi) &= 2^{l} l! \int_{\mathbb{R}^{l}} \mathbbm{1}_{t \in \xi_{C}} \mathbbm{1}_{t_{1} \ge \dots \ge t_{l}} \varphi(t_{1}) \dots \varphi(t_{l}) dt_{1} \dots dt_{l} \\ &= \frac{2^{l} l!}{(l-s)!} \int_{\mathbb{R}^{l-s} \times \mathbb{R}^{s}} \mathbbm{1}_{t \in \xi_{C}} \mathbbm{1}_{t_{1} \ge \dots \ge t_{s}} \varphi(t_{1}) \dots \varphi(t_{s}) \varphi(t_{s+1}) \dots \varphi(t_{l}) dt_{1} \dots dt_{l} \\ &\leq \frac{2^{l} l!}{(l-s)!} \left(\frac{1}{2\pi}\right)^{\frac{l-s}{2}} \int_{\mathbb{R}^{s}} \mathbbm{1}_{t_{1} \ge \dots \ge t_{s}} \frac{\lambda_{l-s} (B_{1}(C(t_{1}+\dots+t_{l})))}{2^{l-s}} \varphi(t_{1}) \dots \varphi(t_{s}) dt_{1} \dots dt_{s} \,, \end{split}$$

where the last inequality relies on $\mathbb{P}(\mathcal{N}(0, I_{l-s}) \in B_1(t_1 + \dots + t_s))$ is bounded by the density function of $\mathcal{N}(0, I_{l-s})$ in 0 times the Lebesgue measure of the l_1 ball of radius $C(t_1 + \dots + t_s)$ in \mathbb{R}^{l-s} . Finally, as

$$\lambda_{l-s}(B_1(R)) = \frac{(2R)^{l-s}}{(l-s)!},$$

it implies,

$$\mathbb{P}(\xi) \leq \left(\frac{2C^2}{\pi}\right)^{\frac{l-s}{2}} \frac{2^{s}l!}{(l-s)!^2} \int_{\mathbb{R}^s} (t_1 + \dots + t_s)^{l-s} \mathbb{1}_{t_1 \geq \dots \geq t_s} \varphi(t_1) \dots \varphi(t_s) dt_1 \dots dt_s
= \left(\frac{2C^2}{\pi}\right)^{\frac{l-s}{2}} \frac{l!}{(l-s)!^2 s!} \mathbb{E}\left((|W_1| + \dots + |W_s|)^{l-s}\right)
= \left(\frac{2C^2}{\pi}\right)^{\frac{l-s}{2}} \binom{l}{s} \frac{1}{(l-s)!} \mathbb{E}(||W||_1^{l-s}),$$

where W is a standard Gaussian vector in \mathbb{R}^s . At last, using bound on l_1 norm, it comes,

,

$$\mathbb{P}(\xi) \le \left(\frac{2C^2}{\pi}\right)^{\frac{l-s}{2}} {l \choose s} \frac{1}{(l-s)!} s^{\frac{l-s}{2}} \mathbb{E}(||W||_2^{l-s}) \\ = \left(\frac{2C^2}{\pi}\right)^{\frac{l-s}{2}} {l \choose s} \frac{1}{(l-s)!} s^{\frac{l-s}{2}} 2^{\frac{l-s}{2}} \frac{\Gamma(l/2)}{\Gamma(s/2)} ,$$

where the last equality follows from classical results on the moment of the χ distribution. $\hfill \Box$

Appendix C. Spherical Hessian

Lemma 5 — Denote X''(.) the Hessian of X(.) along the sphere \mathbb{S}^{m-1} then $X''(t) = -X(t)I_{m-1}.$

Proof. To compute the spherical Hessian, since every point plays the same role, we can compute it at the "east pole" $t = e_1$, the first vector of the canonical basis. Consider a basis (w_2, \ldots, w_n) of the tangent space at $t = e_1$ and use, as a chart of the sphere, the orthogonal projection on this space.

Let $Y(t_2, \ldots, t_m)$ be the process X(.) written in this chart in some neighborhood of e_1 . By the Pythagorean theorem,

$$Y(t_2,...,t_m) = X(\sqrt{1-t_2^2-\dots-t_m^2},t_2,\dots,t_m).$$

Plugging this into the order two Taylor expansion of the process X(.) at $t = e_1$ gives

$$Y(t_2, \dots, t_m) = X(e_1) + t_2 X_2'(e_1) + \dots + t_m X_m'(e_1) + \sum_{2 \le i, j \le m} \frac{t_i t_j}{2} X_{ij}''(e_1) - X_1'(e_1) \frac{t_2^2 + \dots + t_m^2}{2} + o(t_2^2 + \dots + t_m^2),$$

where $X'_k(e_1) = \frac{\partial X}{\partial w_k}(e_1)$ and $X''_{ij}(e_1) = \frac{\partial X}{\partial w_i \partial w_j}(e_1)$. As the process is locally linear, in a small enough neighborhood of e_1, X''_{ij} is equal to zero and, by identification,

$$X''(e_1) = -X'_1(e_1)I_{m-1} = -X(e_1)I_{m-1}$$

giving the desired result.

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