

# Some basic properties and characterizations of Gaussian measures and variables

The note (to be completed!) collects, in a random order, some basic properties and classical characterizations of Gaussian measures and variables, which may be found in standard references in probability theory and mathematical statistics. In the text,  $\gamma_n$ , or  $\mathcal{N}(0, \text{Id})$ , denote the standard Gaussian distribution on the Borel sets of  $\mathbb{R}^n$  with density  $\frac{1}{(2\pi)^{\frac{n}{2}}} e^{-\frac{1}{2}|x|^2}$ ,  $x \in \mathbb{R}^n$ , with respect to the Lebesgue measure  $\lambda_n$ .

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# 1 Rotational invariance

The rotational invariance of Gaussian measures may be described by two main features.

1) If  $X$  is a random variable on a probability space  $(\Omega, \mathcal{A}, \mathbb{P})$  with law  $\mathcal{N}(0, \text{Id})$ , and if  $O$  is an orthogonal matrix (transformation) on  $\mathbb{R}^n$ , then  $OX$  is also distributed according to  $\mathcal{N}(0, \text{Id})$ . This is checked on the covariance matrix.

Maxwell's observation is that the standard Gaussian measure in  $\mathbb{R}^n$  is the only probability measure which is both invariant under orthogonal transformations and is a product measure. More precisely, if  $\mu$  is a probability measure on the Borel sets of  $\mathbb{R}^n$  with those two properties, necessarily  $\mu = \mathcal{N}(0, \sigma^2 \text{Id})$  for some  $\sigma > 0$ . For a quick argument, after convolution with  $\mathcal{N}(0, \varepsilon^2 \text{Id})$  for some  $\varepsilon > 0$ , it may be assumed that  $\mu$  has a smooth, strictly positive, density  $f$  with respect to the Lebesgue measure. But this density is both a function of  $|x|^2$  by rotational invariance and a product (necessarily of the same one-dimensional density), which forces the Gaussian density (take the partial derivatives of  $\log f$ ).

2) If  $X$  and  $Y$  are centered Gaussian vectors on a probability space  $(\Omega, \mathcal{A}, \mathbb{P})$ , independent and identically distributed, for every real  $\theta$ ,  $X(\theta) = X \sin(\theta) + Y \cos(\theta)$  and  $X'(\theta) = X \cos(\theta) - Y \sin(\theta)$  are Gaussian vectors, independent with the same law as  $X$ . In others words, the couples  $(X(\theta), X'(\theta))$  have the same law as  $(X, Y)$ . This may easily be checked on the covariances.

The Kac-Bernstein theorem [6, 4] ensures conversely that if  $X$  and  $Y$  are independent real random variables such that  $X + Y$  and  $X - Y$  are also independent, then both  $X$  and  $Y$  must have normal distributions. More generally, the statement holds true as soon as  $X(\theta)$  and  $X'(\theta)$  are independent for some  $\theta$  which is not an integer multiple of  $\frac{\pi}{2}$ .

Cramér's theorem [5] (initially announced by P. Lévy) expresses that if  $X$  and  $Y$  are two independent (non-constant) real random variables such that  $X + Y$  is normally distributed, then both  $X$  and  $Y$  follow a normal distribution.

## 2 Wick's formula

It is clear that, for a real random variable  $X$  on a probability space  $(\Omega, \mathcal{A}, \mathbb{P})$ , the collection of moments

$$\mathbb{E}(X^{2k+1}) = 0, \quad \mathbb{E}(X^{2k}) = \frac{(2k)!}{2^k k!}, \quad k \geq 0,$$

characterize the law of  $X$  as the standard normal distribution  $\mathcal{N}(0, 1)$ .

Wick's formula is a kind of multilinear extension which expresses any product of the coordinates of a Gaussian random vector by the covariance of this vector. Namely, if  $X$  is a centered Gaussian vector in  $\mathbb{R}^n$ , for any even collection  $(\xi_1, \dots, \xi_{2k})$  of linear functions on  $\mathbb{R}^n$ ,

$$\mathbb{E}(\xi_1(X) \cdots \xi_{2k}(X)) = \sum \prod_{\ell=1}^k \mathbb{E}(\xi_{i_\ell}(X) \xi_{j_\ell}(X))$$

where the sum runs over all unordered sequences of unordered pairs  $\{i_1, j_1\}, \dots, \{i_k, j_k\}$  where each of the integers  $1, \dots, 2k$  appears only once.

For example, if  $X = (X_1, X_2, X_3, X_4)$  is a centered Gaussian vector in  $\mathbb{R}^4$ ,

$$\begin{aligned} \mathbb{E}(X_1 X_2 X_3 X_4) &= \mathbb{E}(X_1 X_2) \mathbb{E}(X_3 X_4) + \mathbb{E}(X_1 X_3) \mathbb{E}(X_2 X_4) \\ &\quad + \mathbb{E}(X_1 X_4) \mathbb{E}(X_2 X_3). \end{aligned}$$

Among other proofs, this type of identity may be obtained from the form of the Fourier transform

$$\mathbb{E}(e^{i\langle u, X \rangle}) = e^{-\frac{1}{2} \mathbb{E}(\langle u, X \rangle^2)}, \quad u = (u_1, u_2, u_3, u_4) \in \mathbb{R}^4.$$

Identification of the terms with all the  $u_\ell$ 's distincts in Taylor expansions of both sides at the order 4 yields the conclusion.

## 3 Fourier transform

The Fourier transform, of characteristic function in the probabilistic language, of the standard Gaussian measure  $\gamma_n$  is given by

$$\varphi_{\gamma_n}(u) = e^{-\frac{1}{2}|u|^2}, \quad u \in \mathbb{R}^n.$$

In other words, the density  $f(x) = \frac{1}{(2\pi)^{\frac{n}{2}}} e^{-\frac{1}{2}|x|^2}$ ,  $x \in \mathbb{R}^n$ , of  $\gamma_n$ , is its own Fourier transform  $\hat{f}$  (up to the standard Fourier normalizations).

## 4 Integration by parts and Stein's characterization

For a function  $f : \mathbb{R} \rightarrow \mathbb{R}$ , locally Lipschitz and such that  $xf$  and  $f'$  are integrable with respect to  $\gamma_1$ ,

$$\int_{\mathbb{R}} xf \, d\gamma_1 = \int_{\mathbb{R}} f' \, d\gamma_1. \quad (1)$$

This immediately follows by integration by parts on the form of the Gaussian density

$$\int_{\mathbb{R}} xf \, d\gamma_1 = \frac{1}{\sqrt{2\pi}} \int_{\mathbb{R}} f(x)x e^{-\frac{1}{2}x^2} d\lambda_1(x).$$

On  $\mathbb{R}^n$ , if  $f : \mathbb{R}^n \rightarrow \mathbb{R}$  is smooth enough,

$$\int_{\mathbb{R}^n} xf \, d\gamma_n = \int_{\mathbb{R}^n} \nabla f \, d\gamma_n$$

as vector integrals.

Stein's observation, leading to various approximation results cf. [2], expresses that (1) ranging over a rich enough family of functions  $f$  characterizes  $\gamma_1$  (over all probability measures with a first moment). Let indeed  $\mu$  be a probability on  $\mathbb{R}$  with a first moment such that  $\int_{\mathbb{R}} xf \, d\mu = \int_{\mathbb{R}} f' \, d\mu$  for every, say, bounded continuous function  $f : \mathbb{R} \rightarrow \mathbb{R}$ . Applied to the real and imaginary parts of the family of functions  $f(x) = e^{iux}$ ,  $x \in \mathbb{R}$ ,  $u \in \mathbb{R}$ , this integration by parts formula yields that the Fourier transform  $\varphi(u)$ ,  $u \in \mathbb{R}$ , of  $\mu$  satisfies the differential equation  $\varphi'(u) = -u\varphi(u)$ ,  $u \in \mathbb{R}$ , so that  $\varphi(u) = e^{-\frac{1}{2}u^2}$ ,  $u \in \mathbb{R}$ , the Fourier transform of  $\gamma_1$ .

## 5 Heat and Mehler kernels

Let

$$h_t(x) = \frac{1}{(4\pi t)^{\frac{n}{2}}} e^{-\frac{1}{4t}|x|^2}, \quad t > 0, \quad x \in \mathbb{R}^n, \quad (2)$$

be the standard heat kernel on  $\mathbb{R}^n$ , fundamental solution of the heat equation  $\partial_t h_t = \Delta h_t$ . In other words,  $h_t$  is the density of the normal law  $\mathcal{N}(0, 2t)$ .

The convolution semigroup  $H_t f(x) = f * h_t(x)$ ,  $t > 0$ , solves

$$\partial_t H_t f = \Delta H_t f = H_t \Delta f$$

with initial data  $f$ . By the definition of  $h_t$ , the semigroup  $H_t$ ,  $t > 0$ , admits the integral representation

$$H_t f(x) = \int_{\mathbb{R}^n} f(x + \sqrt{2t}y) d\gamma_n(x)$$

for all  $t > 0$ ,  $x \in \mathbb{R}^n$ , and any suitable measurable function  $f : \mathbb{R}^n \rightarrow \mathbb{R}$ . At  $t = \frac{1}{2}$ ,  $h_t$  is just the standard Gaussian density so that  $H_{\frac{1}{2}}f(0) = \int_{\mathbb{R}^n} f d\gamma_n$  (while  $H_0f = f$ ).

There is however a related Gaussian kernel which has the advantage to be invariant with respect to  $\gamma_n$  (as the classical heat kernel is invariant under the Lebesgue measure  $\lambda_n$ ).

Define the Mehler kernel, for  $t > 0$ ,  $x, y \in \mathbb{R}^n$ , by

$$p_t(x, y) = p_t(y, x) = \frac{1}{(1 - e^{-2t})^{\frac{n}{2}}} \exp \left( - \frac{e^{-2t}}{2(1 - e^{-2t})} [|x|^2 + |y|^2 - 2e^t \langle x, y \rangle] \right). \quad (3)$$

It holds true that  $\int_{\mathbb{R}^n} p_t(x, y) d\gamma_n(y) = 1$  for all  $t > 0$  and  $x \in \mathbb{R}^n$ . The Mehler kernel satisfies besides the basic semigroup property with respect to  $\gamma_n$ ,

$$\int_{\mathbb{R}^n} p_s(x, z) p_t(z, y) d\gamma_n(z) = p_{s+t}(x, y)$$

for all  $s, t > 0$  and  $x, y \in \mathbb{R}^n$ .

The Mehler kernel generates the Ornstein-Uhlenbeck semigroup

$$P_t f(x) = \int_{\mathbb{R}^n} f(y) p_t(x, y) d\gamma_n(y)$$

for all  $t > 0$ ,  $x \in \mathbb{R}^n$ , and any suitable measurable function  $f : \mathbb{R}^n \rightarrow \mathbb{R}$ , which after a suitable change of variable admits the integral representation

$$P_t f(x) = \int_{\mathbb{R}^n} f(e^{-t}x + \sqrt{1 - e^{-2t}}y) d\gamma_n(y). \quad (4)$$

With the natural extension  $P_0 = \text{Id}$ , the family  $(P_t)_{t \geq 0}$  defines a Markov semigroup, symmetric in  $L^2(\gamma_n)$  and invariant with respect to  $\gamma_n$ , that is  $\int_{\mathbb{R}^n} f P_t g d\gamma_n = \int_{\mathbb{R}^n} g P_t f d\gamma_n$  and  $\int_{\mathbb{R}^n} P_t f d\gamma_n = \int_{\mathbb{R}^n} f d\gamma_n$ . These properties are actually a reformulation of the rotational invariance of Gaussian measures, expressing that under  $\gamma_n \otimes \gamma_n$ , the couples

$$(x \sin(\theta) + y \cos(\theta), x \cos(\theta) - y \sin(\theta)),$$

with  $e^{-t} = \sin(\theta)$ , are distributed as  $(x, y)$ .

The infinitesimal generator  $L = \Delta - x \cdot \nabla$  of the Markov semigroup  $(P_t)_{t \geq 0}$  fulfills the integration by parts formula

$$\int_{\mathbb{R}^n} f(-Lg) d\gamma_n = \int_{\mathbb{R}^n} \nabla f \cdot \nabla g d\gamma_n \quad (5)$$

for every smooth functions  $f, g : \mathbb{R}^n \rightarrow \mathbb{R}$ . The spectrum of the operator  $-L$  is  $\mathbb{N}$ , and the eigenvectors are the Hermite polynomials (cf. [3]). For example, in dimension 1, if  $(h_k)_{k \in \mathbb{N}}$  denotes the sequence of Hermite polynomials (normalized in  $L^2(\gamma_1)$ ),

$$-Lh_k = k h_k, \quad k \in \mathbb{N}.$$

In particular, by the integration by parts formula (5),

$$k \int_{\mathbb{R}} h_k f d\gamma_1 = \int_{\mathbb{R}} (-Lh_k) f d\gamma_1 = \sqrt{k} \int_{\mathbb{R}} h_{k-1} f' d\gamma_1$$

for every smooth function  $f : \mathbb{R} \rightarrow \mathbb{R}$ , which extends (1) since  $h_1(x) = x$ ,  $h_0(x) = 1$ ,  $x \in \mathbb{R}$ .

## 6 Maximum of entropy

If  $P$  is a probability measure on the Borel sets of  $\mathbb{R}^n$ , with density  $f$  with respect to the Lebesgue measure, its *entropy* is

$$H(P) = - \int_{\mathbb{R}^n} f \log(f) d\lambda_n$$

whenever the integral is well-defined,  $H(P) = +\infty$  if not. It is easily seen that  $H(\gamma_n) = \frac{n}{2} \log(2\pi e)$ .

If  $f$  and  $g$  are probability densities, by Jensen's inequality with respect to the convex function  $-\log$  and to the probability measure  $f d\lambda_n$ ,

$$\int_{\mathbb{R}^n} f \log\left(\frac{f}{g}\right) d\lambda_n = \int_{\mathbb{R}^n} (-\log)\left(\frac{g}{f}\right) f d\lambda_n \geq -\log\left(\int_{\mathbb{R}^n} g d\lambda_n\right) = 0.$$

In other words,  $\int_{\mathbb{R}^n} f \log(f) d\lambda_n \geq \int_{\mathbb{R}^n} f \log(g) d\lambda_n$ .

Let now  $P$  be a probability measure with density  $f$  and finite entropy  $H(P)$ , satisfying  $\int_{\mathbb{R}^n} |x|^2 f d\lambda_n \leq n$  ( $= \int_{\mathbb{R}^n} |x|^2 d\gamma_n$ ). If  $g$  is the density of the Gaussian distribution  $\gamma_n$ , by the preceding,

$$\begin{aligned} H(P) &= - \int_{\mathbb{R}^n} f \log(f) d\lambda_n \\ &\leq - \int_{\mathbb{R}^n} f \log(g) d\lambda_n \\ &= \frac{1}{2} \int_{\mathbb{R}^n} (|x|^2 + n \log(2\pi)) f d\lambda_n \\ &\leq \frac{n}{2} (1 + \log(2\pi)) = H(\gamma_n) \end{aligned}$$

since  $\int_{\mathbb{R}^n} |x|^2 f d\lambda_n \leq n$ .

As a conclusion, the standard Gaussian measure  $\gamma_n$  maximizes the entropy over all probability measures with density  $f$  of finite entropy satisfying  $\int_{\mathbb{R}^n} |x|^2 f d\lambda_n \leq n$ . By the case of equality in Jensen's inequality, the proof shows at the same time that the Gaussian measure  $\gamma_n$  is characterized in this way, a result commonly attributed to L. Boltzmann.

## 7 Chi-squared distribution

If  $X_1, \dots, X_n$  are independent variables with common law  $\mathcal{N}(0, 1)$ , then  $X_1^2 + \dots + X_n^2$  follows the classical  $\chi^2$  law with  $n$  degree of freedom, expressed by the gamma distribution with density  $\frac{1}{2^{\frac{n}{2}} \Gamma(\frac{n}{2})} x^{\frac{n}{2}-1} e^{-\frac{x}{2}}$ ,  $x \in (0, \infty)$ , with respect to the Lebesgue measure.

## 8 Independence of empirical mean and variance

If  $X = (X_1, \dots, X_n)$  is random vector with law  $\mathcal{N}(0, \text{Id})$ , the empirical mean and variance of the sample  $(X_1, \dots, X_n)$  defined by

$$\bar{X} = \frac{1}{n} \sum_{k=1}^n X_k \quad \text{and} \quad S^2 = \frac{1}{n-1} \sum_{k=1}^n (X_k - \bar{X})^2$$

are independent. Moreover,  $(n-1)S^2$  has the same distribution as  $\sum_{k=1}^{n-1} X_k^2$  (that is, follows a  $\chi^2$  law with  $n-1$  degree of freedom).

To verify these properties, let  $O$  be a  $n \times n$  orthogonal matrix, and  $Y = OX$ . Then  $\sum_{k=1}^n X_k^2 = \sum_{k=1}^n Y_k^2$ , and  $Y$  has the same distribution as  $X$ . Choosing  $O$  such that the coefficients of the last line are all equal to  $\frac{1}{\sqrt{n}}$ , so that  $Y_n = \sqrt{n} \bar{X}$ , it may be checked that

$$(n-1)S^2 = \sum_{k=1}^n X_k^2 - n\bar{X}^2 = \sum_{k=1}^n Y_k^2 - Y_n^2 = \sum_{k=1}^{n-1} Y_k^2.$$

Since  $Y = (Y_1, \dots, Y_n)$  has law  $\mathcal{N}(0, \text{Id})$ , the coordinates  $Y_1, \dots, Y_n$  are independent standard normal real random variables. As a consequence,  $\bar{X} = \frac{1}{\sqrt{n}} Y_n$  is independent from  $(n-1)S^2 = \sum_{k=1}^{n-1} Y_k^2$ , which proves the various claims.

## 9 Maximum of iid Gaussians

It is a classical exercise to check that if  $X_1, \dots, X_n$ ,  $n \geq 1$ , are independent standard normal random variables on a probability space  $(\Omega, \mathcal{A}, \mathbb{P})$ , then

$$c \sqrt{\log n} \leq \mathbb{E}(\max_{1 \leq k \leq n} X_k) \leq C \sqrt{\log n} \quad (6)$$

for numerical constants  $0 < c < C < \infty$ .

The upper-bound actually holds in a wider generality. Let  $X_1, \dots, X_n$ ,  $n \geq 2$ , be centered Gaussian variables on  $(\Omega, \mathcal{A}, \mathbb{P})$  with respective variances  $\sigma_1^2, \dots, \sigma_n^2$ . Then, for any  $0 < p < \infty$ , there is a constant  $C_p > 0$  only depending on  $p$  such that

$$\left[ \mathbb{E}(\max_{1 \leq k \leq n} |X_k|^p) \right]^{1/p} \leq C_p \max_{1 \leq k \leq n} \sigma_k \sqrt{\log n}. \quad (7)$$

As a quick proof, by homogeneity, it may be assumed that  $\max_{1 \leq k \leq n} \sigma_k \leq 1$ . For  $p \geq 2$ , let  $\Psi : \mathbb{R}_+ \rightarrow \mathbb{R}_+$  be the convex function equal to  $e^{\frac{1}{4}x^{2/p}}$  on the complement of the interval  $[0, (2p-4)^{p/2}]$ , and equal to  $e^{\frac{1}{4}(2p-4)}$  on this interval. By Jensen's inequality,

$$\exp\left(\frac{1}{4} \left[ \mathbb{E}(\max_{1 \leq k \leq n} |X_k|^p) \right]^{2/p}\right) \leq \Psi\left(\mathbb{E}(\max_{1 \leq k \leq n} |X_k|^p)\right) \leq \mathbb{E}\left(\Psi(\max_{1 \leq k \leq n} |X_k|^p)\right).$$

Now

$$\begin{aligned} \mathbb{E}\left(\Psi(\max_{1 \leq k \leq n} |X_k|^p)\right) &\leq e^{\frac{1}{4}(2p-4)} + \mathbb{E}(e^{\frac{1}{4} \max_{1 \leq k \leq n} X_k^2}) \\ &\leq e^{\frac{1}{4}(2p-4)} + \sum_{k=1}^n \mathbb{E}(e^{\frac{1}{4} X_k^2}) \\ &\leq e^{\frac{1}{4}(2p-4)} + n \sqrt{2}, \end{aligned}$$

from which the claim follows.

Towards the lower-bound in (6), since the  $X_1, \dots, X_n$  are independent with common law  $\gamma_1$ , the distribution function of the random variable  $\max_{1 \leq k \leq n} X_k$  is  $\Phi(t)^n$ ,  $t \in \mathbb{R}$  (where  $\Phi$  is the distribution function of  $\gamma_1$ ), so its law has density  $n\Phi(x)^{n-1} \frac{1}{\sqrt{2\pi}} e^{-\frac{1}{2}x^2}$ ,  $x \in \mathbb{R}$ , with respect to the Lebesgue measure  $\lambda_1$  on  $\mathbb{R}$ . Hence

$$\mathbb{E}(\max_{1 \leq k \leq n} X_k) = \int_{\mathbb{R}} x n \Phi(x)^{n-1} \frac{1}{\sqrt{2\pi}} e^{-\frac{1}{2}x^2} d\lambda_1(x) = \int_{\mathbb{R}} n(n-1) \Phi(x)^{n-2} \frac{1}{2\pi} e^{-x^2} d\lambda_1(x)$$

where the second equality follows from the integration by parts formula (1). In particular, if  $n = 2$ ,  $\mathbb{E}(\max(X_1, X_2)) = \frac{1}{2\sqrt{\pi}}$ , so that in the following it may be assumed that  $n$  is large

enough (larger than some fixed  $n_0$ ). Now, for any  $\alpha \in \mathbb{R}$ ,

$$\begin{aligned} \mathbb{E}\left(\max_{1 \leq k \leq n} X_k\right) &\geq \int_{\alpha}^{\infty} n(n-1)\Phi(x)^{n-2} \frac{1}{2\pi} e^{-x^2} d\lambda_1(x) \\ &\geq n(n-1)\Phi(\alpha)^{n-2} \int_{\alpha}^{\infty} \frac{1}{2\pi} e^{-x^2} d\lambda_1(x) \\ &= n(n-1)\Phi(\alpha)^{n-2} \frac{1}{2\sqrt{\pi}} [1 - \Phi(\sqrt{2}\alpha)]. \end{aligned}$$

Take then  $\alpha = \alpha_n = \sqrt{2 \log n - \log \log n}$ ,  $n \geq n_0$ . Recall the classical tail estimates (cf. [1])

$$\left(\frac{1}{t} - \frac{1}{t^3}\right) \frac{1}{\sqrt{2\pi}} e^{-\frac{1}{2}t^2} \leq 1 - \Phi(t) \leq \frac{1}{t} \frac{1}{\sqrt{2\pi}} e^{-\frac{1}{2}t^2}$$

for every  $t > 0$ . By the upper-bound,  $\Phi(\alpha_n) \geq 1 - \frac{c}{n}$  for some numerical  $c > 0$ , while by the lower-bound,  $1 - \Phi(\sqrt{2}\alpha_n) \geq \frac{c}{n^2} \sqrt{\log n}$ . The claim follows.

## References

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