# Lecture 1: Gaussian Random Element

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### 1 Gaussian Vector and its Distributions

#### 1.1 Univariate Case

• Definition (Gaussian Random Variable)

Let  $(\mathbb{R}, \mathcal{B}(\mathbb{R}))$  be a measurable space, where  $\mathcal{B}(\mathbb{R})$  is the Borel  $\sigma$ -algebra on  $\mathbb{R}$ . A real-valued random variable X is **Normally distributed** or **Gaussian** with expectation  $\mu \in \mathbb{R}$  and variance  $\sigma^2 > 0$ , if its **distribution density** with respect to Lebesgue measure is

$$p(x) = \frac{1}{\sqrt{2\pi\sigma^2}} \exp\left(-\frac{(x-\mu)^2}{2\sigma^2}\right).$$

- Remark The followings are properties to the Gaussian distribution
  - 1. The c.d.f. for the standard Normal distribution  $\mathcal{N}(0,1)$  is

$$\Phi(x) = \frac{1}{\sqrt{2\pi}} \int_{-\infty}^{x} \exp(-u^2/2) du$$

- 2. p(x) is **unimodal**, **symmetric** about the mean  $\mu$  and it is **uniformly bounded** on  $\mathbb{R}$ . which has a **unique maximum**  $\frac{1}{\sqrt{2\pi\sigma^2}}$  at the mean  $x = \mu$ .
- 3. The Normal distribution has *super-exponential decay tail*; that is, when x moves away from  $\mu$ , p(x) decreases *monotonically* and *very fast*.
- 4. The **barycenter** (or the center of gravity) of  $\mathcal{N}(\mu, \sigma^2)$  is  $x = \mu$  due to  $\int (x-\mu)p(x)dx = 0$ ; and the **second central moment**  $\int (x-\mu)^2p(x)dx = \sigma^2$ .
- 5. The characteristic function (Fourier transforms) and moment generating function (Laplace transforms)

$$\mathcal{F}\left\{p\right\} = \mathbb{E}_p\left[\exp(i\omega x)\right] = \exp\left(i\mu\omega - \frac{1}{2}\omega^2\sigma^2\right)$$
$$\mathcal{L}\left\{p\right\} = \mathbb{E}_p\left[\exp(sx)\right] = \exp\left(s\mu + \frac{1}{2}s^2\sigma^2\right)$$

6.  $\mathcal{N}(\mu_1, \sigma_1^2) * \mathcal{N}(\mu_2, \sigma_2^2) = \mathcal{N}(\mu_1 + \mu_2, \sigma_1^2 + \sigma_2^2)$ , where \* is the **convolution operation**. In other words, the family  $\{\mathcal{N}(\mu, \sigma^2)\}$  is **stable** with respect to convolutions

$$\mathcal{P}_1 * \mathcal{P}_2(A) \equiv \int_r \mathcal{P}_1(A-r)\mathcal{P}_2(dr), \ A \in \mathcal{B}^1.$$

7. The **Gaussian measure** is **convex**. (Note not the density function p(x) but the measure  $d\mathcal{P} = p(x)dx$ ). That is, for any sets  $A, B \in \mathcal{B}(\mathbb{R})$ , and each  $\gamma \in [0, 1]$ ,

$$\gamma g(\mathcal{P}(A)) + (1 - \gamma)g(\mathcal{P}(B)) \le g(\mathcal{P}(\gamma A + (1 - \gamma)B))$$

where  $g: \mathbb{R}_+ \to \mathbb{R}_+$  is a normalizing function. For Gaussian measure,  $g = \Phi^{-1}$  the inverse c.d.f.

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#### 1.2 Multivariate Case

• Definition (Standard Gaussian Random Vector) A random vector  $X = (X_j)_{j=1}^n \in \mathbb{R}^n$  is called <u>standard Gaussian</u>, if its components are independent and have a standard normal distribution. The distribution of X has a density

$$p(\boldsymbol{x}) = \frac{1}{(2\pi)^n} \exp\left(-\frac{1}{2}\boldsymbol{x}^T\boldsymbol{x}\right), \text{ for } \boldsymbol{x} \in \mathbb{R}^n.$$
 (1)

• Definition (Gaussian Random Vector)

A random vector  $Y \in \mathbb{R}^n$  is called <u>Gaussian</u>, if it can be represented as Y = a + LX, where X is a standard Gaussian vector,  $\mathbf{a} \in \mathbb{R}^n$ , and  $L : \mathbb{R}^n \to \mathbb{R}^n$  is a **linear mapping**.

Equivalently,

Definition (Gaussian Random Vector)

A random vector  $Y \in \mathbb{R}^n$  is called <u>Gaussian</u>, if  $\langle v, Y \rangle$  is a Normal random variable for each  $v \in \mathbb{R}^n$ .

• Definition (Covariance Operator for Gaussian Random Vector) Given a Gaussian random vector  $X = (X_j)_{j=1}^n$ , define the <u>covariance operator</u> as a linear mapping  $K_X : \mathbb{R}^n \to \mathbb{R}^n$  such that

$$cov(\langle u, X \rangle, \langle v, X \rangle) = \langle u, K_X(v) \rangle.$$

The matrix representation of  $K_X$  is called a **covariance matrix** 

$$\mathbf{K} = [K(i,j)]_{i,j=1}^n \in \mathbb{R}^{n \times n}, \quad \text{where } K(i,j) = \langle e_i, K_X(e_j) \rangle.$$

• Remark (The Covariance Operator is Self-Adjoint and Positive)

The covariance operator K is self-adjoint  $(K_X^* = K_X)$ , positive semi-definite  $K \succeq 0$ .

This is due to the symmetry and positive definiteness property of inner product.

Equivalently, the covariance matrix K is symmetric, positive semi-definite.

• Remark (Density for Multivariate Gaussian) In the case, when the linear mapping L is invertible (non-degenerate), the multivariate Normal distribution  $\mathcal{N}(\mu, K)$  can be defined via its density function w.r.t. the Lebesgue measure on  $\mathbb{R}^n$ 

$$p(\boldsymbol{x}) = \frac{1}{\sqrt{(2\pi)^n \det(|\Sigma|)}} \exp\left(-\frac{1}{2}(\boldsymbol{x} - \boldsymbol{\mu})^T \boldsymbol{K}^{-1}(\boldsymbol{x} - \boldsymbol{\mu})\right)$$
(2)

• Remark The expression for density in (2) holds only if the linear operator L is invertible; that is, the general definition used is the linear projection definition. [Lifshits, 2013]

If L is singular, K is singular, i.e., det K = 0; there is no proper density expression as (2). On the other hand, for every K nonnegative definite,  $L = K^{1/2}$  exists and is nonnegative definite as well.

• Remark (The Characteristic Function of Multivariate Gaussian)

The characteristic functions of  $\mathcal{N}(\boldsymbol{\mu}, K)$  is determined by its one-dimension projection

$$\varphi(\boldsymbol{v}) = \int \exp\left(i\langle \boldsymbol{x}, \boldsymbol{v}\rangle\right) \mathcal{P}(d\boldsymbol{x})$$

$$= \int \exp\left(ir\right) \mathcal{P}^{\boldsymbol{v}}(dr)$$

$$= \exp\left(i\mu^{\boldsymbol{v}}\omega - \frac{1}{2}\sigma^{2}(\boldsymbol{v})\omega^{2}\right)\Big|_{\omega=1}$$

$$= \exp\left(i\langle \boldsymbol{\mu}, \boldsymbol{v}\rangle - \frac{1}{2}\langle K\boldsymbol{v}, \boldsymbol{v}\rangle\right)$$
(3)

The equation (3) is known as the *characteristic functional* of measure  $\mathcal{P}$ .

Use the affine mapping  $\mu + L\mathcal{P}_0$ , the characteristic functional is given by

$$\varphi(\boldsymbol{v}) = \int \exp\left(i\left\langle \boldsymbol{\mu} + L\boldsymbol{x}, \, \boldsymbol{v}\right\rangle\right) \mathcal{P}_0(d\boldsymbol{x})$$

$$= \exp\left(i\left\langle \boldsymbol{\mu}, \, \boldsymbol{y}\right\rangle\right) \int \exp\left(i\left\langle L\boldsymbol{x}, \, \boldsymbol{v}\right\rangle\right) \mathcal{P}_0(d\boldsymbol{x})$$

$$= \exp\left(i\left\langle \boldsymbol{\mu}, \, \boldsymbol{y}\right\rangle\right) \int \exp\left(i\left\langle \boldsymbol{x}, \, L^*\boldsymbol{v}\right\rangle\right) \mathcal{P}_0(d\boldsymbol{x})$$

$$= \exp\left(i\left\langle \boldsymbol{\mu}, \, \boldsymbol{y}\right\rangle - \frac{1}{2}\left\langle L^*\boldsymbol{v}, \, L^*\boldsymbol{v}\right\rangle\right)$$

$$= \exp\left(i\left\langle \boldsymbol{\mu}, \, \boldsymbol{y}\right\rangle - \frac{1}{2}\left\langle LL^*\boldsymbol{v}, \, \boldsymbol{v}\right\rangle\right)$$

$$= \exp\left(i\left\langle \boldsymbol{\mu}, \, \boldsymbol{y}\right\rangle - \frac{1}{2}\left\langle LL^*\boldsymbol{v}, \, \boldsymbol{v}\right\rangle\right)$$

And the density is computed, for L invertible, by change of variable for  $y = \mu + Lx$ 

$$p_{\boldsymbol{\mu},K}(\boldsymbol{y}) = \left| \det L \right|^{-1} p(\boldsymbol{x})$$
$$= (2\pi)^{n/2} \left| \boldsymbol{K} \right|^{-1/2} \exp\left( -\left\langle K^{-1}(\boldsymbol{y} - \boldsymbol{\mu}), \, \boldsymbol{y} - \boldsymbol{\mu} \right\rangle / 2 \right)$$

- Proposition 1.1 (Existence and Uniqueness of Gaussian Distribution) [Lifshits, 2013] Let  $\mathcal{P}$  be a Gaussian distribution in  $\mathbb{R}^n$ . Then the mean value  $\mu$  and the covariance operator K of the measure  $\mathcal{P}$  exist and are uniquely defined. The operator K is symmetric and positive definite.
- Proposition 1.2 (Gaussian Random Vector from Kernel) [Lifshits, 2013] Assume  $\mu \in \mathbb{R}^n$  and  $K : \mathbb{R}^n \to \mathbb{R}^n$  is nonnegative definite linear operator. Then there exists a unique Gaussian distribution  $\mathcal{N}(\mu, K)$  with mean  $\mu$  and covariance operator K. The characteristic functional of  $\mathcal{N}(\mu, K)$  has the form of (3). If the operator K is non-singular, the distribution  $\mathcal{N}(\mu, K)$  is absolutely continuous with respect to the Lebesgue measure, and its density is of form (2). There are no other Gaussian distribution in  $\mathbb{R}^n$ , except for the form  $\mathcal{N}(\mu, K)$ .

### 2 Gaussian Random Element

#### 2.1 Gaussian Random Element in Topological Vector Space

Definition (Random Element in Topological Vector Space)
Let (Ω, F, P) be a probability space, (X, B) be a topological vector space with σ-algebra
B. A random element in X is a F/B-measurable function X : Ω → X so that

$$X^{-1}(A) \in \mathscr{F}, \quad \forall A \in \mathscr{B}.$$

We write  $X \in \mathcal{X}$ .

• Definition (Duality)

Let  $\mathcal{X}^*$  be the dual space of  $\mathcal{X}$ , i.e. the space of bounded linear functional on  $\mathcal{X}$ .

We denote  $\langle f, x \rangle$  the **duality** between the spaces X and X\*, i.e.

$$\langle f, x \rangle := f(x), \quad \forall f \in X^*, x \in X.$$

Note that we **do not confuse this notation with inner product**. In inner product  $\langle x, y \rangle$  both arguments are from the same space.

• Definition (Gaussian Random Element in Topological Vector Space) A random element  $X \in \mathcal{X}$  is called Gaussian, if

$$\langle f\,,\,X\rangle:=f(X)$$

is a *Normal random variable*, for all  $f \in \mathcal{X}^*$ .

• Definition (*Expectation*)

A vector  $a \in \mathcal{X}$  is called **expectation** of a random element  $X \in \mathcal{X}$ , if

$$\mathbb{E}\left[\langle f, X \rangle\right] = \langle f, a \rangle$$

for all  $f \in \mathcal{X}^*$ . We write  $a = \mathbb{E}[X]$ .

• Definition (Covariance Operator)

A linear operator  $K: \mathcal{X}^* \to \mathcal{X}$  is called <u>covariance operator</u> of a random vector  $X \in \mathcal{X}$ , if

$$cov(\langle f, X \rangle, \langle g, X \rangle) = \langle f, K g \rangle.$$

for all  $f, g \in X^*$ . We write K = cov(X).

Remark (Covariance as Function-Valued Linear Transformation on Dual Space ) The covariance operator  $K: \mathcal{X}^* \to \mathcal{X}$  acts on linear functional on  $\mathcal{X}$  and returns an element (function) in  $\mathcal{X}$ 

$$f(Kq) := cov(f(X), q(X))$$

• Remark (Covariance Operator is Self-Adjoint and Positive)

Covariance operator is self-adjoint, due to symmetric property of covariance in R.

$$\left\langle f\,,\,Kg\right\rangle =\left\langle g\,,\,Kf\right\rangle ,\quad\forall f,g\in X^{\ast },$$

and it is **positive** (semi-definite), i.e.

$$\langle f, Kf \rangle = \text{var}(f(X)) \ge 0, \quad \forall f \in X^*.$$

• Remark (Topological Constraints on  $\mathcal{X}$  for Gaussian Element) [Lifshits, 2012] From the definition of Gaussian element, we see that it only makes sense when the space of continuous linear functionals on  $\mathcal{X}$  is rich enough. For example, if  $\mathcal{X}^* = \{0\}$ , then any vector satisfies this definition rendering it senseless.

Therefore, usually one of three situations of increasing generality is considered.

- 1.  $\mathcal{X}$  is a **separable Banach space**, for example,  $\mathcal{C}[0,1]$ ,  $L^p[0,1]$  etc;
- 2.  $\mathcal{X}$  is a <u>complete separable locally convex metrizable</u> topological vector space, for example,  $\mathcal{C}[0,\infty)$ ,  $\mathbb{R}^{\infty}$  etc.
- 3.  $\mathcal{X}$  is a <u>locally convex topological vector space</u> and a vector X is such that its distribution is a <u>Radon measure</u>.

In cases (1) and (2) every **finite measure** is a **Radon measure**, thus case (3) is the most general one. These assumptions are called usual assumptions in [Lifshits, 2012, 2013]

- Proposition 2.1 (Existence of Covariance Operator) [Lifshits, 2013] Under usual assumptions on  $\mathcal{X}$ , every Gaussian random element in  $\mathcal{X}$  possesses an expectation and a covariance operator. In other words, the distribution of Gaussian elements in  $\mathcal{X}$  is of the form  $\mathcal{N}(a, K)$ .
- Remark (Distribution and Characteristic Function of Gaussan Random Element) The pair (a, K) determines the distribution of a Gaussian variable  $\langle f, x \rangle$  as

$$\mathcal{N}(\langle f, a \rangle, \langle f, Kf \rangle),$$

and we find  $\emph{the characteristic function}$  of  $\langle f\,,\,x\rangle$ 

$$\begin{split} \varphi(\langle f\,,\,X\rangle) &= \mathbb{E}\left[\exp\left\{i\omega\,\langle f\,,\,x\rangle\right\}\right] \\ &= \exp\left(i\omega\,\langle f\,,\,a\rangle - \frac{1}{2}\omega^2\,\langle f\,,\,Kf\rangle\right) \\ &:= \exp\left(i\omega f(a) - \frac{1}{2}\omega^2 f(Kf)\right) \end{split}$$

Any Radon distribution in  $\mathcal{X}$  is determined by its characteristic function. Therefore, distribution  $\mathcal{N}(a, K)$  is **unique**.

#### 2.2 Examples of Gaussian Random Elements

• Example (Standard Gaussian Measure in  $\mathbb{R}^{\infty}$ )
Consider the space  $\mathcal{X} = \mathbb{R}^{\infty}$  of all countable infinite sequence  $(x_1, x_2, \ldots)$  equipped with the **product topology**. The product topology induces a metric as

$$\rho(\{x_n\}_{n=1}^{\infty}, \{y_n\}_{n=1}^{\infty}) = \sup_{n} \left\{ \frac{\min |x_n - y_n|, 1}{n} \right\}.$$

 $\mathbb{R}^{\infty}$  is a **complete separable metric space** under the product topology. The dual space  $\mathcal{X}^* = c_0$  is the space of sequences  $(f_1, f_2, \ldots)$  with  $f_n = 0$  for all but finite number of n. The duality

$$\langle f, x \rangle = \sum_{n=1}^{\infty} f_n x_n < \infty.$$

Consider a sequence of i.i.d.  $\mathcal{N}(0,1)$ -distributed random variables as a vector  $X \in \mathcal{X}$ , i.e.  $X := (X_n)_{n=1}^{\infty}, X_n \sim \mathcal{N}(0,1)$ . Due to stability of normal distribution, for any  $f \in \mathcal{X}^*$  the random variable

$$\langle f, X \rangle = \sum_{n=1}^{\infty} f_n X_n \sim \mathcal{N}(0, \sigma^2)$$

where  $\sigma^2 = \sum_{n=1}^{\infty} f_n^2 < \infty$ . Therefore, X is a **Gaussian element**. It is clear that  $\mathbb{E}[X] = 0$ .

**Embedding operator** serves as **covariance operator** for X, i.e.

$$K = \iota : c_0 \hookrightarrow \mathbb{R}^{\infty}.$$

To show that

$$\operatorname{cov}(\langle f, X \rangle, \langle g, X \rangle) = \mathbb{E}\left[\langle f, X \rangle \langle g, X \rangle\right]$$

$$= \mathbb{E}\left[\left(\sum_{n=1}^{\infty} f_n X_n\right) \left(\sum_{n=1}^{\infty} g_n X_n\right)\right]$$

$$= \mathbb{E}\left[\sum_{n,m=1}^{\infty} f_n g_m X_n X_m\right]$$

$$= \sum_{n,m=1}^{\infty} f_n g_m \mathbb{E}\left[X_n X_m\right] = \sum_{n,m=1}^{\infty} f_n g_m \,\delta_{n,m}$$

$$= \sum_{n=1}^{\infty} f_n g_n := \langle f, Kg \rangle$$

We call the distribution of X a standard Gaussian measure in  $\mathbb{R}^{\infty}$ .

• Example (Gaussian Elements in a Hilbert space  $\mathcal{H}$ ) [Lifshits, 2012] Let  $\mathcal{X} = \mathcal{H}$  be a separable Hilbert space whose inner product will be denoted by  $\langle \cdot, \cdot \rangle_{\mathcal{H}}$ . By the Riesz representation theorem, we can identify its dual space  $\mathcal{H}^*$  with  $\mathcal{H}$ , i.e. for each  $f \in \mathcal{H}^*$ , there exists a unique  $x_f \in \mathcal{H}$  such that

$$\langle f, x \rangle = f(x) = \langle x, x_f \rangle_{\mathcal{H}}, \quad \forall x \in \mathcal{H}.$$

Define  $h: \mathcal{H}^* \to \mathcal{H}$  as an isometric isomorphism that maps  $f \mapsto x_f$ .

In order to construct a Gaussian element in  $\mathcal{H}$ , consider a *complete orthonormal basis*  $\{\varphi_n\}_{n=1}^{\infty}$  on  $\mathcal{H}$ , a sequence of *independent*  $\mathcal{N}(0,1)$ -distributed random variables  $\{\xi_n\}_{n=1}^{\infty}$ , and a sequence of *non-negative numbers*  $\{\sigma_n\}_{n=1}^{\infty}$  satisfying assumption  $\sum_{n=1}^{\infty} \sigma_n^2 < \infty$  so that the series

$$\sum_{n=1}^{\infty} \sigma_n \xi_n(\omega) \varphi_n$$

is **convergent** in  $\|\cdot\|_{\mathcal{H}}$ -norm almost surely in  $\mathcal{H}$ . Define a random element  $X:\Omega\to\mathcal{H}$  as the limit of the series

$$X = \sum_{n=1}^{\infty} \sigma_n \xi_n \varphi_n \tag{4}$$

This representation is called *Karhunen-Loève expansion*.

For any linear functional  $f \in \mathcal{H}^*$ , we can its corresponding vector  $x_f = h(f) \in \mathcal{H}$  and  $x_f = \sum_{n=1}^{\infty} f_n \varphi_n$ . Thus the random variable

$$\langle f, X \rangle = \langle X, x_f \rangle_{\mathcal{H}} = \left\langle \sum_{n=1}^{\infty} \sigma_n \xi_n \varphi_n, \sum_{n=1}^{\infty} f_n \varphi_n \right\rangle_{\mathcal{H}}$$

$$= \sum_{n,m=1}^{\infty} \sigma_n \bar{f_m} \xi_n \langle \varphi_n, \varphi_m \rangle_{\mathcal{H}}$$
by orthonormal  $\langle \varphi_n, \varphi_m \rangle_{\mathcal{H}} = \delta_{n,m}$ 

$$= \sum_{n=1}^{\infty} \sigma_n \bar{f_n} \xi_n \sim \mathcal{N}(0, \sigma^2)$$

where  $\sigma^2 := \sum_{n=1}^{\infty} \sigma_n^2 f_n^2 \le (\sum_{n=1}^{\infty} \sigma_n^2) \sup_n |f_n|^2 < \infty$ . Therefore, X is a **Gaussian random element** in  $\mathcal{H}$  and  $\mathbb{E}[X] = 0$ . In order to find **the covariance operator** of X, let us compute

$$\operatorname{cov}\left(\left\langle f\,,\,X\right\rangle\left\langle g\,,\,X\right\rangle\right) = \mathbb{E}\left[\left\langle f\,,\,X\right\rangle\left\langle g\,,\,X\right\rangle\right]$$

$$= \mathbb{E}\left[\left(\sum_{n=1}^{\infty}\sigma_{n}\bar{f}_{n}\xi_{n}\right)\left(\sum_{n=1}^{\infty}\sigma_{n}\bar{g}_{n}\xi_{n}\right)\right]$$

$$= \sum_{n,m=1}^{\infty}\bar{f}_{n}\bar{g}_{m}\sigma_{n}\sigma_{m}\mathbb{E}\left[\xi_{n}\xi_{m}\right]$$

$$\operatorname{since}\,\mathbb{E}\left[\xi_{n}\xi_{m}\right] = \delta_{n,m}$$

$$= \sum_{n=1}^{\infty}\sigma_{n}^{2}\bar{f}_{n}\bar{g}_{n} = \left\langle f\,,\,Kg\right\rangle$$

By plugging in the basis, we have

$$K: g \to \sum_{n=1}^{\infty} \sigma_n^2 g_n \varphi_n = \sum_{n=1}^{\infty} \sigma_n^2 \langle g, \varphi_n \rangle \varphi_n$$
 (5)

$$\Rightarrow \widetilde{K} = K \circ h^{-1} = \sum_{n=1}^{\infty} \sigma_n^2 \langle \cdot , \varphi_n \rangle_{\mathcal{H}} \varphi_n$$
 (6)

Therefore  $\sigma_n^2$  and  $\varphi_n$  are the **eigenvalues** and **eigenfunctions** of  $\widetilde{K} = K \circ h^{-1}$  and  $\widetilde{K}$  is a **positive**, **compact operator** on  $\mathcal{H}$  since  $\operatorname{tr}(\widetilde{K}) = \sum_{n=1}^{\infty} \sigma_n^2 < \infty$ .

One can show that any Gaussian element in a Hilbert space admits a representation (4) [Lifshits, 2012]. This means that a Gaussian distribution with covariance operator K exists if and only if the induced linear operator  $\widetilde{K} = K \circ h^{-1} \in \mathcal{L}(\mathcal{H})$  is a self-adjoint, positive, trace-class operator (which is compact).

• Remark (Equivalent Definition of Covariance Operator on Hilbert Space) In the previous example, we see that the covariance operator on Hilbert space can be equivalently defined via linear operator  $\widetilde{K}: \mathcal{H} \to \mathcal{H}$  so that

$$\operatorname{cov}\left(\langle f_h, X \rangle_{\mathcal{H}}, \langle g_h, X \rangle_{\mathcal{H}}\right) = \left\langle \widetilde{K} f_h, g_h \right\rangle_{\mathcal{H}}.$$

Note that  $\widetilde{K} \succeq 0$  is **self-adjoint** and **positive** and it has **finite** trace  $tr(\widetilde{K})$  so it is **traceclass operator** which is **compact**. And, conversely, for any **positive trace-class operator**  $K \in \mathcal{B}_1(\mathcal{H})$ , there exists **Gaussian element** in  $\mathcal{H}$  with distribution  $\mathcal{N}(0,K)$ .

• Remark (Identity Operator is Not Covariance Operator on Hilbert Space) For identity operator  $I: \mathcal{H} \to \mathcal{H}$ , we see that its trace  $\operatorname{tr}(I) = \infty$ , this means that it does not admit a Gaussian distribution as  $\mathcal{N}(0,I)$  on infinite dimensional space  $\mathcal{H}$ . In fact, we can see that  $\mathbb{E}\left[|X(t)|^2\right] = \infty$ .

#### 2.3 Gaussian Random Process

• Definition (Random Process)

Let  $(\Omega, \mathscr{F}, \mathcal{P})$  be a probability space and T be a parametric set called *index set*. A random process X on T is a family of random variables  $X(t, \omega), t \in T$ , defined on the common probability space  $(\Omega, \mathscr{F}, \mathcal{P})$ . For each  $\omega \in \Omega$ ,

$$X(\omega) := \{X_t(\omega) : t \in T\}$$

is called a **sample function** of  $(X_t)$  and if T is one-dimensional, they are often called **sample paths** of the process  $(X_t)$ .

- Remark Determined by index set T, we have:
  - 1. if  $T \subset \mathbb{R}$ ,  $\{X_t\}_{t \in T}$  is called a **random process**.
  - 2. if  $T \subset \mathbb{R}^n$ ,  $\{X_t\}_{t \in T}$  is called a *random field*.
  - 3. if  $T = \mathbb{N}$ ,  $\{X_t\}_{t \in T}$  is called a **random sequence**.
- ullet Definition (Gaussian Random Process)

A process  $(X_t)_{t\in T}$  is called <u>Gaussian</u> if for any  $t_1, \ldots, t_n \in T$  the distribution of the random vector

$$(X(t_1),\ldots,X(t_n))$$

is a *Gaussian distribution* in  $\mathbb{R}_n$ .

The properties of a *Gaussian process* are *completely determined* by its *expectation*  $\mathbb{E}[X(t)], t \in T$ , and *covariance*  $cov(X(s), X(t)), s, t \in T$ .

• Remark (Gaussian Random Process as Gaussian Element on Function Space) Consider the topological vector space  $\mathcal{X} \subset \mathbb{R}^T$  as a function space on T, then the Gaussian random element in  $\mathcal{X}$  is a Gaussian process:

$$X: \Omega \to \mathcal{X} \subset \mathbb{R}^T$$
  
\Rightarrow X(\omega)(t) = X(\omega, t), \forall t \in T

• Definition (Continuous Sample Path)
If T is a topological space, we say that  $\{X_t\}_{t\in T}$  has continuous sample paths, if the function  $X(\cdot,\omega)$  is continuous on T for  $\mathcal{P}$ -almost every  $\omega \in \Omega$ .

### 2.4 Examples of Gaussian Random Processes

• Example (Continuous Sample Path Gaussian Process) [Lifshits, 2012]

Let T be a <u>compact metric space</u>, let  $\mathcal{X} = \mathcal{C}(T)$  denote the Banach space of all continuous functions on T equipped with supremum norm

$$||x||_{\infty} := \sup_{t \in T} |x(t)|$$

and with the corresponding topology of uniform convergence. By Riesz-Markov theorem, the dual space  $\mathcal{X}^* = \mathcal{M}(T)$  is a space of <u>signed Radon measures</u> of finite variations on T. The duality is given by

$$\langle \mu, f \rangle = \int_T f \ d\mu, \quad \forall f \in \mathcal{X}, \forall \mu \in \mathcal{M}(T) = \mathcal{X}^*.$$

Let  $\{X(t), t \in T\}$ , be a *Gaussian random process* with <u>continuous sample paths</u> on the parametric set T. It is *completely characterized* by the functions

$$a(t) := \mathbb{E}[X(t)], \quad K(s,t) := \operatorname{cov}(X(s), X(t)).$$

Then we can view at  $X := \{X(t), t \in T\}$  as a **Gaussian random element** of  $\mathcal{X}$ . The **expectation** of X is computed as

$$\mathbb{E}\left[X\right] = a := (a(t))_{t \in T},$$

and the **covariance operator**  $K: \mathcal{M}(T) \to \mathcal{C}(T)$  can be calculated by

$$(K\nu)(s) = \int_T K(s,t)\nu(dt). \tag{7}$$

This is because

$$\begin{split} \operatorname{cov}\left(\left\langle \mu\,,\,X\right\rangle\,,\,\left\langle \nu\,,\,X\right\rangle\right) &= \mathbb{E}\left[\left\langle \mu\,,\,(X-a)\right\rangle\left\langle \nu\,,\,(X-a)\right\rangle\right] \\ &= \mathbb{E}\left[\int_{T}(X-a)d\mu\int_{T}(X-a)d\nu\right] \\ &= \mathbb{E}\left[\int_{T\times T}(X(s)-a(s))(X(t)-a(t))\mu(ds)\nu(dt)\right] \\ &= \int_{T}\int_{T}\mathbb{E}\left[(X(s)-a(s))(X(t)-a(t))\right]\mu(ds)\nu(dt) \\ &= \int_{T}\left(\int_{T}K(s,t)\nu(dt)\right)\mu(ds) := \left\langle \mu\,,\,K\nu\right\rangle, \end{split}$$

thus we have (7).

### • Example (Wiener Process) [Lifshits, 2012]

We will now consider T = [0, 1] and  $\mathcal{X} = \mathcal{C}[0, 1]$  with dual  $\mathcal{M}[0, 1]$ . Define a Gaussian element composed of the sample paths of a **Wiener process** 

$$\mathcal{W} := \mathcal{W}(t), \quad 0 \le t \le 1,$$

i.e. of a process satisfying assumptions

$$\mathbb{E}[\mathcal{W}(t)] = 0, \quad \mathbb{E}[\mathcal{W}(s)\mathcal{W}(t)] = \min\{s, t\}.$$

It is just a special case of previous example, so we can find the expectation of  $\mathcal{W}$  by

$$\mathbb{E}\left[\left\langle \mu\,,\,\mathcal{W}\right\rangle \right] = \mathbb{E}\left[\int_{[0,1]}\mathcal{W}d\mu\right] = \int_0^1 \mathbb{E}\left[\mathcal{W}(t)\right]\mu(dt) = 0$$

we have  $\mathbb{E}[\mathcal{W}] = 0$ . Moreover, the *covariance operator*  $K : \mathcal{M}([0,1]) \to \mathcal{C}([0,1])$ 

$$(K\nu)(s) = \int_0^1 K(s,t)\nu(dt)$$
$$= \int_0^1 \min\{s,t\} \nu(dt). \quad \blacksquare$$

**Remark** Finally, we recall the properties of Wiener process W(t): [Lifshits, 2012]

1. It is 1/2-self-similar, i.e. for any c > 0 the process

$$Y(t) := \frac{\mathcal{W}(ct)}{\sqrt{c}}$$

is also a *Wiener process*;

- 2. It has *stationary increments*;
- 3. It has *independent increments*;
- 4. It is a *Markov process*;
- 5. It admits *time inversion*: the process

$$Z(t) := t\mathcal{W}\left(\frac{1}{t}\right)$$

is also a Wiener process.

# References

Mikhail Lifshits. Lectures on gaussian processes. In Lectures on Gaussian Processes, pages 1–117. Springer, 2012.

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