Normal distributions on vector spaces

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1 Univariate normal distributions

The standard normal distribution is given by the density function:

$$\phi(z) = \frac{1}{\sqrt{2\pi}} e^{-z^2/2} \tag{1.1}$$

This, like all density functions in this section, is understood to be with respect to the Lebesgue integral on \mathbb{R} . A random variable Z that follows this distribution is said to be standard normally distributed, and we write $Z \sim N(0,1)$.

A general, univariate normal distribution with parameters μ and σ , is given by the distribution of $X = \mu + \sigma Z$. If $\sigma \neq 0$, we can invert to find $z = \frac{x-\mu}{\sigma}$. So $\frac{dz}{dx} = \frac{1}{\sigma}$. According to the usual transformation rules, x has the density function:

$$f(x) = \phi\left(\frac{x-\mu}{\sigma}\right)\frac{1}{\sigma} = \frac{1}{\sqrt{2\pi\sigma^2}}e^{-(x-\mu)^2/2\sigma^2}$$
(1.2)

We write $X \sim N(\mu, \sigma^2)$.

1.1 What if $\sigma = 0$?

The derivation above assumes σ to be non-zero. But even if this is the case, $X = \mu$ still has a distribution - it is simply μ all the time. However, this random variable does not have a density function with respect to the Lebesgue measure. In light of the Radon-Nikodym theorem, this is because the Lebesgue measure does not dominate the probability measure of X: $P_X(\{\mu\}) = 1, m_1(\{\mu\}) = 0$.

In the multivariate case, we will often run into similar problems.

2 Random vectors

Before tackling the multivariate case, we need some basic tools. In this section, we consider vectors of random variables. So a random vector of dimension n is:

$$X = \begin{pmatrix} X_1 \\ X_2 \\ \vdots \\ X_n \end{pmatrix} \tag{2.1}$$

Here, each X_i is a random variable.

2.1 Variance

The variance of a *n*-dimensional vector is the $n \times n$ matrix:

$$Var(X) = E[(X - \mu_X)(X - \mu_X)^t]$$
 (2.2)

Here $\mu_X = E[X]$, i.e. the vector of expectation values of the X_i 's. From the usual definitions of variances and covariances between random variables, we see that the diagonal of Var(X) contains the variances of each X_i , while the off diagonal elements are the covariances between variables:

$$[\operatorname{Var}(X)]_{ij} = \operatorname{Cov}(X_i, X_j) \tag{2.3}$$

Due to the symmetry of covariance, this means that Var(X) is a symmetric matrix.

2.1.1 Variance calculation rules

Similarly to ordinary random variables, we might calculate the variance matrix as follows:

$$Var(X) = E[(X - \mu_X)(X - \mu_X)^t] =$$
(2.4)

$$E(XX^{t}) - E(X)\mu_{X}^{t} - \mu_{X}E(X)^{t} + \mu_{X}\mu_{X}^{t} =$$
 (2.5)

$$E(XX^t) - \mu_X \mu_X^t \tag{2.6}$$

Here, we've used the linearity of the expectation value and the definition of μ_X .

Adding a constant vector b does not change the variance, since $E[X+b] = \mu_X + b$:

$$Var(X+b) = E[(X+b-(\mu_X+b))(X+b-(\mu_X+b))^t] = E[(X-\mu_X)(X-\mu_X)^t]$$
(2.7)

This is just the variance of X.

If A is a constant $m \times n$ matrix and X is an n-dimensional random vector, then:

$$Var(AX) = E[(AX - A\mu_X)(AX - A\mu_X)^t] =$$
 (2.8)

$$E[(A(X - \mu_X))(A(X - \mu_X))^t] =$$
 (2.9)

$$E[A(X - \mu_X)(X - \mu_X)^t A^t] =$$
 (2.10)

$$A[(X - \mu_X)(X - \mu_X)^t]A^t \tag{2.11}$$

So we have $Var(AX) = A Var(X)A^t$.

2.2 Covariance

The covariance matrix between two variable vectors X and Y is:

$$Cov(X,Y) = E[(X - \mu_X)(Y - \mu_Y)^t]$$
 (2.12)

If X has dimension m and Y dimension n, then Cov(X, Y) has dimension $m \times n$. Here, the matrix elements reduce to ordinary covariances between X_i 's and Y_j s:

$$[\operatorname{Cov}(X,Y)]_{ij} = \operatorname{Cov}(X_i,Y_j) \tag{2.13}$$

This also means, that $Cov(X, Y) = (Cov(Y, X))^t$

We note, that the variance could have been defined as a special case of covariance, since Var(X) = Cov(X, X).

2.2.1 Covariance calculation rules

Similarly to the rule for variances, we have:

$$Cov(X,Y) = E[XY^t] - \mu_X \mu_Y^t$$
(2.14)

The proof is essentially the same.

If A and B are constant matrices of appropriate dimesion, we also have:

$$Cov(AX, BY) = A Cov(X, Y)B^{t}$$
(2.15)

Again, the proof is entirely analogous to the corresponding variance formula. The covariance is bilinear:

$$Cov(X + Y, Z) = Cov(X, Z) + Cov(Y, Z)$$
(2.16)

$$Cov(X, Y + Z) = Cov(X, Y) + Cov(X, Z)$$
(2.17)

This follows from the bilinearity of the ordinary covariance.

2.2.2 Addiational variance formulas

Since we noted that Var(X) = Cov(X, X), we may use these rules to derive further properties of variances.

For instance, the variance of a sum:

$$Var(X+Y) = Cov(X+Y,X+Y) =$$
(2.18)

$$Cov(X, X) + Cov(X, Y) + Cov(Y, X) + Cov(Y, Y) = (2.19)$$

$$Var(X) + Var(Y) + Cov(X, Y) + Cov(Y, X)$$
(2.20)

This mirrors the formula for the covariance of sums ordinary random variables, but is complicated by the fact that the vector covariance is not symmetric.

2.3 Quadratic forms

If X is an n-dimensional random variable and A an $n \times n$ matrix, then the corresponding quadratic form is $Q = X^t A X$. I.e. a scalar. What is the expectation value of the quadratic form? We can use a trick here. Since Q is a scalar, we can trivially write this as a trace:

$$Q = X^t A X = \operatorname{tr}(X^t A X) = \operatorname{tr}(A X X^t)$$
 (2.21)

Here, we've used the cyclic property of traces. Now, the expectation value is:

$$E[Q] = E[\operatorname{tr}(AXX^{t})] = \operatorname{tr}(E[AXX^{t}]) = \operatorname{tr}(A\ E[XX^{t}])$$
 (2.22)

But we know, that $Var(X) = E(XX^t) - \mu_X \mu_X^t$, so $E[XX^t] = Var(X) + \mu_X \mu_X^t$:

$$E[Q] = \operatorname{tr}(A(\operatorname{Var}(X) + \mu_X \mu_X^t)) = \operatorname{tr}(A \operatorname{Var}(X)) + \operatorname{tr}(A\mu_X \mu_X^t)$$
 (2.23)

The last term may be rewritten:

$$\operatorname{tr}(A\mu_X\mu_X^t) = \operatorname{tr}(\mu_X^t A\mu_X) = \mu_X^t A\mu_X \tag{2.24}$$

In the last step we've used that the contents of the parenthesis is a scalar. So all in all:

$$E[X^t A X] = \operatorname{tr}(A \operatorname{Var}(X)) + \mu_X^t A \mu_X$$
 (2.25)

3 Multivariate normal distribution

The standard normal distribution in n dimensions is the distribution of n independent standard normals. Hence, the density function in \mathbb{R}^n is simply

a product of terms like equation 1.1:

$$\phi(z) = \prod_{i=1}^{n} \frac{1}{\sqrt{2\pi}} e^{-z_i^2/2} = (2\pi)^{-n/2} e^{||z||^2/2}$$
(3.1)

Here $z \in \mathbb{R}^n$, and $||\cdot||$ is the standard Euclidean norm. If a *n*-dimensional dimensional random vector Z follows this distribution we write $Z \sim N(0, I_n)$, where I_n is the identity matrix in n dimensions. The reason for this will soon be clear.

4 Affine transformations of euclidean spaces

In order to get to the general, multidimensional normal distribution, we need yet another component:

Let $s: \mathbb{R}^n \to \mathbb{R}^m$ be a linear transformation. This means that the is an $m \times n$ matrix A so s(x) = Ax.

An *affine* transformation t is formed by following this linear map by a translation:

$$t: \mathbb{R}^n \to \mathbb{R}^m, t(x) = Ax + v \tag{4.1}$$

Here, $v \in \mathbb{R}^m$. Since translations are always bijective, we note that t is bijective iff A is invertible.

Each component of an affine transformation is composed from measurable function - is is understood that we mean with respect to the Borel algebras of each space) - so the affine transformation itself is measurable as well.

4.1 Transformation properties of the Lebesgue measure

Recall that the Lebesgue measure in n dimensions m_n is invariant under translation: If t is a translation $t: \mathbb{R}^n \to \mathbb{R}^n, t(x) = x + x_0$, where $x_0 \in \mathbb{R}^n$ then:

$$t(m_n) = m_n (4.2)$$

Also, if $s: \mathbb{R}^n \to \mathbb{R}^n, x \mapsto Ax$ is an isomorphism, then:

$$s(m_n) = m_n |\det A^{-1}|$$
 (4.3)

Combining the two, the formula for affine transformation is the same as for linear ones.

5 General multivariate normal distribution

Given an *n*-dimension random vector $Z \sim N(0, I_n)$, we now define a general standard normal as the distribution of the random variable $X = \mu + \sum_{i=1}^{n} Z_i$.

Some notes are in order, before we procees. For now " $\Sigma^{\frac{1}{2}}$ " should simply be considered a matrix of dimensions $n \times p$, not necessarily a square root of anything (we will get back to that later). It plays the roles that σ played for the univariate case¹, and hence we might expect that special care needs to be taken when it is "zero", whatever that might signify here.

6 Orthogonal complement

Let V be a finite-dimensional vector space with an inner product $\langle \cdot, \cdot \rangle$. Let U be a subspace of V. Then we define the *orthogonal complement* of U as:

$$U^{\perp} = \{ v \in V | \forall u \in U : \langle u, v \rangle = 0 \}$$

$$(6.1)$$

Theorem 6.1. U^{\perp} is a subspace of V.

Proof. According to the subspace theorem, we need to show three things:

- U^{\perp} is not empty: Clearly $0 \in U^{\perp}$.
- Closed under addition: If $v_1, v_2 \in U^{\perp}$, then for all $u \in U^{\perp}$:

$$\langle v_1 + v_2, u \rangle = \langle v_1, u \rangle + \langle v_2, u \rangle = 0 \tag{6.2}$$

• Closed under scalar multiplication: If $v \in U^{\perp}$ and $c \in \mathbb{R}$ then for all $u \in U^{\perp}$:

$$\langle cv, u \rangle = c \langle v, u \rangle = 0$$
 (6.3)

Since the only vector perpendicular to itself is 0, we further conclude that $U \cap U^{\perp} = \{0\}.$

Theorem 6.2. If e_1, e_2, \ldots, e_m is an orthonormal basis for U, then for any $v \in V$:

$$v - \sum_{i=1}^{m} \langle v, e_i \rangle e_i \in U^{\perp}$$
 (6.4)

¹Hence, simply calling it Σ would be more be clearer, but such is the tradition notation.

Proof. Let $u \in U$. Then we can write $u = \sum_{j=1}^{m} \lambda_j e_j$ for some coefficients λ_j . Now calculate the inner product with the vector above:

$$\langle v - \sum_{i=1}^{m} \langle v, e_i \rangle e_i , \sum_{j=1}^{m} \lambda_j e_j \rangle = \sum_{i=j}^{m} \lambda_j \langle v, e_j \rangle - \sum_{i=1}^{m} \sum_{j=1}^{m} \lambda_j \langle v, e_i \rangle \langle e_i, e_j \rangle$$
 (6.5)

Since $\langle e_i, e_j \rangle = \delta_{ij}$ this vanishes.

This means that we may write any $v \in V$ as a sum of vectors from U and U^{\perp} respectively:

$$v = \underbrace{\sum_{i=1}^{m} \langle v, e_i \rangle e_i}_{\in U} + \underbrace{v - \sum_{i=1}^{m} \langle v, e_i \rangle e_i}_{\in U^{\perp}}$$
(6.6)

Theorem 6.3. The decomposition into elements from U and U^{\perp} from equation 6.6 is unique.

Proof. Let $v=u_1+u_1^{\perp}$ and $v=u_2+u_2^{\perp}$ be two such decompositions. Then $u_1+u_1^{\perp}=u_2+u_2^{\perp}$ and hence $u_1-u_2=u_2^{\perp}-u_1^{\perp}$. But this means that this vector is a member of both U and U^{\perp} , and hence it must be 0. This means $u_1=u_2$ and $u_1^{\perp}=u_2^{\perp}$.

6.1 The orthogonal projection

The previous section motivates the following:

Definition 6.1. Let V be a finite-dimensional inner product vector space and U a subspace of V. The orthogonal projection from V onto U is the map $p:V \to V$ which satisfies:

$$\forall v \in V: \quad p(v) \in U, \quad v - p(v) \in U^{\perp} \tag{6.7}$$

As we see, one could also define the co-domain of p to be U. Usually, the distinction will not matter much.

Theorem 6.4. The orthogonal projection operator is linear.

Proof. We need to show additivity and homogeneity:

• Additivity: Let $v_1, v_2 \in V$. Then $p(v_1) + p(v_2) \in U$ and:

$$v_1 - p(v_1) + v_2 - p(v_2) = v_1 + v_2 - (p(v_1) + p(v_2)) \in U^{\perp}$$
 (6.8)

Adding the two we get $v_1 + v_2$. So $p(v_1 + v_2) = p(v_1) + p(v_2)$.

• Homogeneity. Let $v \in V$ and $c \in \mathbb{R}$. Then $cp(v) \in U$ and $c(v - p(v)) = cv - cp(v) \in U^{\perp}$. Adding the two we get cv, so p(cv) = cp(v).

Theorem 6.5. The orthogonal projection operator $p: V \to V$ is idempotent. *I.e.* $p \circ p = p$.

Proof. Let $v \in V$. Then $p(v) \in U$. But this means that the decomposition of p(v) is p(v) + 0. So $p \circ p(v) = p(v)$.

7 Lebesgue measures on vector spaces

7.1 Coordinate maps

Let V be a finite-dimensional vector space of dimension n. Our question is, if we can turn V into a measure space in a natural way. Since we know that V is isomorphic to \mathbb{R}^n , it makes sense to tweak the usual Lebesgue measure in N dimensions:

Let e_1, e_2, \ldots, e_n be a basis for V. Then we can define the *coordinate map* as follows:

$$\phi: \mathbb{R}^n \to V, \quad \begin{pmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{pmatrix} \mapsto \sum_{i=1}^n x_i e_i$$
(7.1)

This is obviously an isomorphism. Specifically, it is invertible with inverse $\phi^{-1}: V \to \mathbb{R}^n$.

The coordinate map depends on the chosen basis. If we had chosen another basis $e_1^*, e_2^*, \ldots, e_n^*$ we would get another isomorphism ϕ^* .

7.2 Borel algebra on V

We can now use ϕ^{-1} to induce a σ -algebra on V. Set \mathbb{B}_V to the smallest σ -algebra that makes ϕ^{-1} measurable when \mathbb{R}^n is equipped with the Borel algebra \mathbb{B}_n . We call \mathbb{B}_V the Borel algebra on V.

At first this object seems to depend of the choice of basis for V. But it turns out that the use of definite article in the definition is justified:

Theorem 7.1. If e_1, e_2, \ldots, e_n and $e_1^*, e_2^*, \ldots, e_n^*$ are bases for V, then the induced σ -algebra \mathbb{B}_V and \mathbb{B}_V^* is the same thing.

Proof. We know that ϕ^{-1} is $\mathbb{B}_V - \mathbb{B}_n$ measurable by definition. We have:

$$(\phi^*)^{-1} = (\phi^*)^{-1} \circ id_V = (\phi^*)^{-1} \circ (\phi \circ \phi^{-1}) = ((\phi^*)^{-1} \circ \phi) \circ \phi^{-1}$$
 (7.2)

 $((\phi^*)^{-1} \circ \phi)$ is a linear operator on \mathbb{R}^n and so according to section 4.1 is measurable. So $(\phi^*)^{-1}$ must be $\mathbb{B}_V - \mathbb{B}_n$ -measurable. Since \mathbb{B}_V^* is the smallest σ -algebra to make $(\phi^*)^{-1} \mathbb{B}_V - \mathbb{B}_n$ -measurable, we must have $\mathbb{B}_V^* \subseteq \mathbb{B}_V$.

But by a totally symmetric argument, we must also have $\mathbb{B}_V \subseteq \mathbb{B}_V^*$. Hence $\mathbb{B}_V = \mathbb{B}_V^*$.

It turns out, that ϕ must be measurable too. This is a direct consequence of the pipeline lemma.

Theorem 7.2. Given two finite-dimensional vector spaces V and W, then:

$$\mathbb{B}_{V \times W} = \mathbb{B}_V \otimes \mathbb{B}_W \tag{7.3}$$

Proof. Let e_1, e_2, \ldots, e_n be a basis for V with corresponding coordinate map ϕ . And f_1, f_2, \ldots, f_m a basis for W with corresponding coordinate map ψ . Then $(e_1, 0), (e_2, 0), \ldots, (e_n, 0), (0, f_1), (0, f_2), \ldots, (0, f_m)$ is a basis for $V \times W$. The corresponding coordinate map is:

$$\phi \times \psi : (x_1, x_2, \dots, x_{n+m}) \mapsto \left(\sum_{i=1}^n x_i e_i, \sum_{j=1}^m x_{n+j} f_j\right)$$
 (7.4)

The inverse is $(\phi \times \psi)^{-1} = \phi^{-1} \times \psi^{-1}$. Since ϕ^{-1} is $\mathbb{B}_V - \mathbb{B}_n$ -measurable and ψ^{-1} is $\mathbb{B}_W - \mathbb{B}_m$ -measurable, $\phi^{-1} \times \psi^{-1}$ must be $\mathbb{B}_V \otimes \mathbb{B}_W - \mathbb{B}_n \otimes \mathbb{B}_n$ -measurable. But $\mathbb{B}_n \otimes \mathbb{B}_m = \mathbb{B}_{n+m}$. Since $\mathbb{B}_{V \times W}$ is the smallest σ -algebra to make $\phi^{-1} \times \psi^{-1}$ measurable, we must have $\mathbb{B}_{V \times W} \subseteq \mathbb{B}_V \otimes \mathbb{B}_W$.

On the other hand, consider the projection operators:

$$\pi_V : V \times W \to V, (v, w) \mapsto v$$
 (7.5)

$$\pi_n : \mathbb{R}^{n+m} \to \mathbb{R}^n, (x_1, \dots, x_{n+m}) \mapsto (x_1, \dots, x_n)$$
 (7.6)

Now consider $\pi_V \circ (\phi \times \psi)$. Applied to an $x \in \mathbb{R}^{n+m}$ we have:

$$\pi_V \circ (\phi \times \psi)(x) = \pi_v \left(\left(\sum_{i=1}^n x_i e_i, \sum_{j=1}^m x_{n+j} f_j \right) \right) = \sum_{i=1}^n x_i e_i$$
 (7.7)

But this is the same as:

$$\phi \circ \pi_1(x) = \phi((x_1, \dots x_n)) = \sum_{i=1}^n x_i e_i$$
 (7.8)

So $\pi_V \circ (\phi \times \psi) = \phi \circ \pi_1$. Now apply $\phi^{-1} \times \psi^{-1}$ from the right to get:

$$\pi_V = \phi \circ \pi_1 \circ (\phi^{-1} \times \psi^{-1}) \tag{7.9}$$

Since all the three functions on the right side are measurable, π_V must be $\mathbb{B}_{V\times W} - \mathbb{B}_V$ -measurable. By a similar argument the corresponding projection operator $\pi_W: V\times W\to W$ is $\mathbb{B}_{V\times W} - \mathbb{B}_W$ -measurable. Since $\mathbb{B}_V\otimes \mathbb{B}_W$ is the smallest σ -algebra to make both π_V and π_W measurable, we must have: $\mathbb{B}_V\otimes \mathbb{B}_W\subseteq \mathbb{B}_{V\times W}$.

7.3 Lebesgue measures on V

We now want to define a measure on the measurable space (V, \mathbb{B}_V) . If e_1, e_2, \ldots, e_n is a basis for V, we will use the associated coordinate map ϕ to define a measure:

$$\lambda_V = \phi(m_n) \tag{7.10}$$

Here, m_n is the usual Lebesgue measure in n dimensions. The problem is, that this measure depends on the chosen basis! Consider another basis $e_1^*, e_2^*, \ldots, e_n^*$ and associated coordinate map ϕ^* . Then the measure is:

$$\lambda_V^* = \phi^*(m_n) = (\phi \circ \phi^{-1}) \circ \phi^*(m_n) = \phi \circ (\phi^{-1} \circ \phi^*(m_n))$$
 (7.11)

Now $\phi^{-1} \circ \phi^*$ is an isomorphism $\mathbb{R}^n \to \mathbb{R}^n$, so according to section 4.1, there is a constant c such that $(\phi^{-1} \circ \phi^*(m_n)) = cm_n$. So:

$$\lambda_V^* = c\phi(m_n) = c\lambda_V \tag{7.12}$$

So while there are many Lebesgue measures on V they only differ from each other by a constant factor. This means that they all agree on what constitutes a null set, and on which functions are integrable. They disagree on the integral, but agree on whether it is finite or not. The also agree on whether a measure μ has a density with respect to λ_V or not.