

1 The Correlation Coefficient

The **Pearson Correlation Coefficient** $\rho(X, Y)$ is a way to measure how linearly related (in other words, how well a linear model captures the relationship between) random variables X and Y .

$$\rho(X, Y) = \frac{\text{cov}(X, Y)}{\sqrt{\text{var}(X) \text{var}(Y)}}$$

Here are some important facts about it:

- It is commutative: $\rho(X, Y) = \rho(Y, X)$
- It always lies between -1 and 1: $-1 \leq \rho(X, Y) \leq 1$
- It is completely invariant to affine transformations: for any $a, b, c, d \in \mathbb{R}$,

$$\begin{aligned} \rho(aX + b, cY + d) &= \frac{\text{cov}(aX + b, cY + d)}{\sqrt{\text{var}(aX + b) \text{var}(cY + d)}} \\ &= \frac{\text{cov}(aX, cY)}{\sqrt{\text{var}(aX) \text{var}(cY)}} \\ &= \frac{a \cdot c \cdot \text{cov}(X, Y)}{\sqrt{a^2 \text{var}(X) \cdot c^2 \text{var}(Y)}} \\ &= \frac{\text{cov}(X, Y)}{\sqrt{\text{var}(X) \text{var}(Y)}} \\ &= \rho(X, Y) \end{aligned}$$

The correlation is defined in terms of random variables rather than observed data. Assume now that $\vec{x}, \vec{y} \in \mathbb{R}^n$ are vectors containing n independent observations of X and Y , respectively. Recall the **law of large numbers**, which states that for i.i.d. X_i with mean μ ,

$$\frac{1}{n} \sum_{i=1}^n X_i \xrightarrow{\text{a.s.}} \mu \quad \text{as } n \rightarrow \infty$$

We can use this law to justify a sample-based approximation to the mean:

$$\text{cov}(X, Y) = \mathbb{E}[(X - \mathbb{E}[X])(Y - \mathbb{E}[Y])] \approx \frac{1}{n} \sum_{i=1}^n (x_i - \bar{x})(y_i - \bar{y})$$

where the bar indicates the sample average, i.e. $\bar{x} = \frac{1}{n} \sum_{i=1}^n x_i$. Then as a special case we have

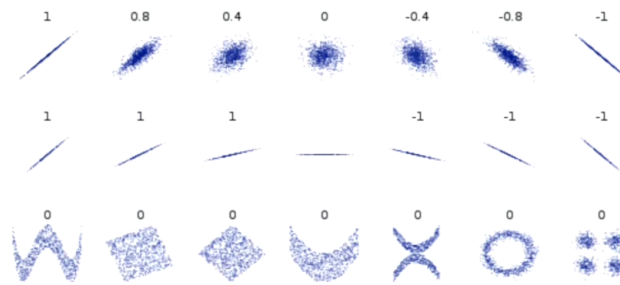
$$\text{var}(X) = \text{cov}(X, X) = \mathbb{E}[(X - \mathbb{E}[X])^2] \approx \frac{1}{n} \sum_{i=1}^n (x_i - \bar{x})^2$$

$$\text{var}(Y) = \text{cov}(Y, Y) = \mathbb{E}[(Y - \mathbb{E}[Y])^2] \approx \frac{1}{n} \sum_{i=1}^n (y_i - \bar{y})^2$$

Plugging these estimates into the definition for correlation and canceling the factor of $1/n$ leads us to the **sample Pearson Correlation Coefficient** $\hat{\rho}$:

$$\begin{aligned} \hat{\rho}(x, y) &= \frac{\sum_{i=1}^n (x_i - \bar{x})(y_i - \bar{y})}{\sqrt{\sum_{i=1}^n (x_i - \bar{x})^2 \cdot \sum_{i=1}^n (y_i - \bar{y})^2}} \\ &= \frac{\tilde{x}^\top \tilde{y}}{\sqrt{\tilde{x}^\top \tilde{x} \cdot \tilde{y}^\top \tilde{y}}} \quad \text{where } \tilde{x} = x - \bar{x}, \tilde{y} = y - \bar{y} \end{aligned}$$

Here are some 2-D scatterplots and their corresponding correlation coefficients:



You should notice that:

- The magnitude of $\hat{\rho}$ increases as X and Y become more linearly correlated.
- The sign of $\hat{\rho}$ tells whether X and Y have a positive or negative relationship.
- The correlation coefficient is undefined if either X or Y has 0 variance (horizontal line).

1.1 Correlation and Gaussians

Here's a neat fact: if X and Y are jointly Gaussian, i.e.

$$\begin{bmatrix} X \\ Y \end{bmatrix} \sim \mathcal{N}(0, \Sigma)$$

then we can define a distribution on *normalized* X and Y and have their relationship entirely captured by $\rho(X, Y)$. First write

$$\rho(X, Y) = \frac{\sigma_{xy}}{\sigma_x \sigma_y}$$

Then

$$\Sigma = \begin{bmatrix} \sigma_x^2 & \sigma_{xy} \\ \sigma_{xy} & \sigma_y^2 \end{bmatrix} = \begin{bmatrix} \sigma_x^2 & \rho \sigma_x \sigma_y \\ \rho \sigma_x \sigma_y & \sigma_y^2 \end{bmatrix}$$

so

$$\begin{aligned} \begin{bmatrix} \sigma_x^{-1} & 0 \\ 0 & \sigma_y^{-1} \end{bmatrix} \begin{bmatrix} X \\ Y \end{bmatrix} &\sim \mathcal{N} \left(0, \begin{bmatrix} \sigma_x^{-1} & 0 \\ 0 & \sigma_y^{-1} \end{bmatrix} \Sigma \begin{bmatrix} \sigma_x^{-1} & 0 \\ 0 & \sigma_y^{-1} \end{bmatrix}^\top \right) \\ &\sim \mathcal{N} \left(0, \begin{bmatrix} 1 & \rho \\ \rho & 1 \end{bmatrix} \right) \end{aligned}$$

2 Canonical Correlation Analysis

Canonical Correlation Analysis (CCA) is a method of modeling the relationship between two point sets by making use of the correlation coefficient. Formally, given zero-mean random vectors $X_{rv} \in \mathbb{R}^p$ and $Y_{rv} \in \mathbb{R}^q$, we want to find projection vectors $u \in \mathbb{R}^p$ and $v \in \mathbb{R}^q$ that maximizes the correlation between $X_{rv}^\top u$ and $Y_{rv}^\top v$:

$$\max_{u,v} \rho(X_{rv}^\top u, Y_{rv}^\top v) = \max_{u,v} \frac{\text{cov}(X_{rv}^\top u, Y_{rv}^\top v)}{\sqrt{\text{var}(X_{rv}^\top u) \text{var}(Y_{rv}^\top v)}}$$

Observe that

$$\begin{aligned} \text{cov}(X_{rv}^\top u, Y_{rv}^\top v) &= \mathbb{E}[(X_{rv}^\top u - \mathbb{E}[X_{rv}^\top u])(Y_{rv}^\top v - \mathbb{E}[Y_{rv}^\top v])] \\ &= \mathbb{E}[u^\top (X_{rv} - \mathbb{E}[X_{rv}])(Y_{rv} - \mathbb{E}[Y_{rv}])^\top v] \\ &= u^\top \mathbb{E}[(X_{rv} - \mathbb{E}[X_{rv}])(Y_{rv} - \mathbb{E}[Y_{rv}])^\top] v \\ &= u^\top \text{cov}(X_{rv}, Y_{rv}) v \end{aligned}$$

which also implies (since $\text{var}(Z) = \text{cov}(Z, Z)$ for any random variable Z) that

$$\begin{aligned} \text{var}(X_{rv}^\top u) &= u^\top \text{cov}(X_{rv}, X_{rv}) u \\ \text{var}(Y_{rv}^\top v) &= v^\top \text{cov}(Y_{rv}, Y_{rv}) v \end{aligned}$$

so the correlation writes

$$\rho(X_{rv}^\top u, Y_{rv}^\top v) = \frac{u^\top \text{cov}(X_{rv}, Y_{rv}) v}{\sqrt{u^\top \text{cov}(X_{rv}, X_{rv}) u \cdot v^\top \text{cov}(Y_{rv}, Y_{rv}) v}}$$

Unfortunately, we do not have access to the true distributions of X_{rv} and Y_{rv} , so we cannot compute these covariances matrices. However, we can estimate them from data. Assume now that we are given zero-mean data matrices $X \in \mathbb{R}^{n \times p}$ and $Y \in \mathbb{R}^{n \times q}$, where the rows of the matrix X are i.i.d. samples $x_i \in \mathbb{R}^p$ from the random variable X_{rv} , and correspondingly for Y_{rv} . Then

$$\text{cov}(X_{rv}, Y_{rv}) = \mathbb{E}[\underbrace{(X_{rv} - \mathbb{E}[X_{rv}])}_0 \underbrace{(Y_{rv} - \mathbb{E}[Y_{rv}])^\top}_0] = \mathbb{E}[X_{rv} Y_{rv}^\top] \approx \frac{1}{n} \sum_{i=1}^n x_i y_i^\top = \frac{1}{n} X^\top Y$$

where again the sample-based approximation is justified by the law of large numbers. Similarly,

$$\begin{aligned} \text{cov}(X_{rv}, X_{rv}) &= \mathbb{E}[X_{rv} X_{rv}^\top] \approx \frac{1}{n} \sum_{i=1}^n x_i x_i^\top = \frac{1}{n} X^\top X \\ \text{cov}(Y_{rv}, Y_{rv}) &= \mathbb{E}[Y_{rv} Y_{rv}^\top] \approx \frac{1}{n} \sum_{i=1}^n y_i y_i^\top = \frac{1}{n} Y^\top Y \end{aligned}$$

Plugging these estimates in for the true covariance matrices, we arrive at the problem

$$\max_{u,v} \frac{u^\top \left(\frac{1}{n} X^\top Y\right) u}{\sqrt{u^\top \left(\frac{1}{n} X^\top X\right) u \cdot v^\top \left(\frac{1}{n} Y^\top Y\right) v}} = \max_{u,v} \frac{u^\top X^\top Y v}{\underbrace{\sqrt{u^\top X^\top X u \cdot v^\top Y^\top Y v}}_{\hat{\rho}(Xu, Yv)}}$$

Let's try to massage the maximization problem into a form that we can reason with more easily. Our strategy is to choose matrices to transform X and Y such that the maximization problem is equivalent but easier to understand.

1. First, let's choose matrices W_x, W_y to **whiten** X and Y . This will make the (co)variance matrices $(XW_x)^\top (XW_x)$ and $(YW_y)^\top (YW_y)$ become identity matrices and simplify our expression. To do this, note that $X^\top X$ is positive definite (and hence symmetric), so we can employ the eigendecomposition

$$X^\top X = U_x S_x U_x^\top$$

Since

$$S_x = \text{diag}(\lambda_1(X^\top X), \dots, \lambda_d(X^\top X))$$

where all the eigenvalues are positive, we can define the "square root" of this matrix by taking the square root of every diagonal entry:

$$S_x^{1/2} = \text{diag}\left(\sqrt{\lambda_1(X^\top X)}, \dots, \sqrt{\lambda_d(X^\top X)}\right)$$

Then, defining $W_x = U_x S_x^{-1/2} U_x^\top$, we have

$$\begin{aligned} (XW_x)^\top (XW_x) &= W_x^\top X^\top X W_x \\ &= U_x S_x^{-1/2} U_x^\top U_x S_x U_x^\top U_x S_x^{-1/2} U_x^\top \\ &= U_x S_x^{-1/2} S_x S_x^{-1/2} U_x^\top \\ &= U_x U_x^\top \\ &= I \end{aligned}$$

which shows that W_x is a whitening matrix for X . The same process can be repeated to produce a whitening matrix $W_y = U_y S_y^{-1/2} U_y^\top$ for Y .

Let's denote the whitened data $X_w = XW_x$ and $Y_w = YW_y$. Then by the change of variables

$$u_w = W_x^{-1}u, v_w = W_y^{-1}v,$$

$$\begin{aligned} \max_{u,v} \hat{\rho}(Xu, Yv) &= \max_{u,v} \frac{(Xu)^\top Yv}{\sqrt{(Xu)^\top Xu (Yv)^\top Yv}} \\ &= \max_{u,v} \frac{(XW_x W_x^{-1}u)^\top YW_y W_y^{-1}v}{\sqrt{(XW_x W_x^{-1}u)^\top XW_x W_x^{-1}u (YW_y W_y^{-1}v)^\top YW_y W_y^{-1}v}} \\ &= \max_{u_w, v_w} \frac{(X_w u_w)^\top Y_w v_w}{\sqrt{(X_w u_w)^\top X_w u_w (Y_w v_w)^\top Y_w v_w}} \\ &= \max_{u_w, v_w} \frac{u_w^\top X_w^\top Y_w v_w}{\sqrt{u_w^\top X_w^\top X_w u_w \cdot v_w^\top Y_w^\top Y_w v_w}} \\ &= \max_{u_w, v_w} \frac{u_w^\top X_w^\top Y_w v_w}{\underbrace{\sqrt{u_w^\top u_w \cdot v_w^\top v_w}}_{\hat{\rho}(X_w u_w, Y_w v_w)}} \end{aligned}$$

Note we have used the fact that $X_w^\top X_w$ and $Y_w^\top Y_w$ are identity matrices by construction.

- Second, let's choose matrices D_x, D_y to **decorrelate** X_w and Y_w . This will let us simplify the covariance matrix $(X_w D_x)^\top (Y_w D_y)$ into a **diagonal** matrix. To do this, we'll make use of the SVD:

$$X_w^\top Y_w = USV^\top$$

The choice of U for D_x and V for D_y accomplishes our goal, since

$$(X_w U)^\top (Y_w V) = U^\top X_w^\top Y_w V = U^\top (USV^\top) V = S$$

Let's denote the decorrelated data $X_d = X_w D_x$ and $Y_d = Y_w D_y$. Then by the change of variables $u_d = D_x^{-1}u_w = D_x^\top u_w, v_d = D_y^{-1}v_w = D_y^\top v_w,$

$$\begin{aligned} \max_{u_w, v_w} \hat{\rho}(X_w u_w, Y_w v_w) &= \max_{u_w, v_w} \frac{(X_w u_w)^\top Y_w v_w}{\sqrt{u_w^\top u_w \cdot v_w^\top v_w}} \\ &= \max_{u_w, v_w} \frac{(X_w D_x D_x^{-1} u_w)^\top Y_w D_y D_y^{-1} v_w}{\sqrt{(D_x u_w)^\top D_x u_w \cdot (D_y v_w)^\top D_y v_w}} \\ &= \max_{u_d, v_d} \frac{(X_d u_d)^\top Y_d v_d}{\sqrt{u_d^\top u_d \cdot v_d^\top v_d}} \\ &= \max_{u_d, v_d} \frac{u_d^\top X_d Y_d v_d}{\underbrace{\sqrt{u_d^\top u_d \cdot v_d^\top v_d}}_{\hat{\rho}(X_d u_d, Y_d v_d)}} \\ &= \max_{u_d, v_d} \frac{u_d^\top S v_d}{\sqrt{u_d^\top u_d \cdot v_d^\top v_d}} \end{aligned}$$

Without loss of generality, suppose u_d and v_d are unit vectors¹ so that the denominator becomes 1, and we can ignore it:

$$\max_{u_d, v_d} \frac{u_d^\top S v_d}{\sqrt{u_d^\top u_d \cdot v_d^\top v_d}} = \max_{\substack{\|u_d\|=1 \\ \|v_d\|=1}} \frac{u_d^\top S v_d}{\|u_d\| \|v_d\|} = \max_{\substack{\|u_d\|=1 \\ \|v_d\|=1}} u_d^\top S v_d$$

The diagonal nature of S implies $S_{ij} = 0$ for $i \neq j$, so our simplified objective expands as

$$u_d^\top S v_d = \sum_i \sum_j (u_d)_i S_{ij} (v_d)_j = \sum_i S_{ii} (u_d)_i (v_d)_i$$

where S_{ii} , the singular values of $X_w^\top Y_w$, are arranged in descending order. Thus we have a weighted sum of these singular values, where the weights are given by the entries of u_d and v_d , which are constrained to have unit norm. To maximize the sum, we “put all our eggs in one basket” and extract S_{11} by setting the first components of u_d and v_d to 1, and the rest to 0:

$$u_d = \begin{bmatrix} 1 \\ 0 \\ \vdots \\ 0 \end{bmatrix} \in \mathbb{R}^p \qquad v_d = \begin{bmatrix} 1 \\ 0 \\ \vdots \\ 0 \end{bmatrix} \in \mathbb{R}^q$$

Any other arrangement would put weight on S_{ii} at the expense of taking that weight away from S_{11} , which is the largest, thus reducing the value of the sum.

Finally we have an analytical solution, but it is in a different coordinate system than our original problem! In particular, u_d and v_d are the best weights in a coordinate system where the data has been whitened and decorrelated. To bring it back to our original coordinate system and find the vectors we actually care about (u and v), we must invert the changes of variables we made:

$$u = W_x u_w = W_x D_x u_d \qquad v = W_y v_w = W_y D_y v_d$$

More generally, to get the best k directions, we choose

$$U_d = \begin{bmatrix} I_k \\ 0_{p-k, k} \end{bmatrix} \in \mathbb{R}^{p \times k} \qquad V_d = \begin{bmatrix} I_k \\ 0_{q-k, k} \end{bmatrix} \in \mathbb{R}^{q \times k}$$

where I_k denotes the k -dimensional identity matrix. Then

$$U = W_x D_x U_d \qquad V = W_y D_y V_d$$

Note that U_d and V_d have orthogonal columns. The columns of U and V , which are the projection directions we seek, will in general not be orthogonal, but they will be linearly independent (since they come from the application of invertible matrices to the columns of U_d, V_d).

¹ Why can we assume this? Observe that the value of the objective does not change if we replace u_d by αu_d and v_d by βv_d , where α and β are any positive constants. Thus if there are maximizers u_d, v_d which are not unit vectors, then $u_d/\|u_d\|$ and $v_d/\|v_d\|$ (which are unit vectors) are also maximizers.

2.1 Comparison with PCA

An advantage of CCA over PCA is that it is invariant to scalings and affine transformations of X and Y . Consider a simplified scenario in which two matrix-valued random variables X, Y satisfy $Y = X + \varepsilon$ where the noise ε has huge variance. What happens when we run PCA on Y ? Since PCA maximizes variance, it will actually project Y (largely) into the column space of ε ! However, we're interested in Y 's relationship to X , not its dependence on noise. How can we fix this? As it turns out, CCA solves this issue. Instead of maximizing variance of Y , we maximize correlation between X and Y . In some sense, we want to maximize “predictive power” of information we have.

3 CCA regression

Once we've computed the CCA coefficients, one application is to use them for regression tasks, predicting Y from X (or vice-versa). Recall that the correlation coefficient attains a greater value when the two sets of data are *more linearly correlated*. Thus, it makes sense to find the $k \times k$ weight matrix A that linearly relates XU and YV . We can accomplish this with ordinary least squares.

Denote the projected data matrices by $X_c = XU$ and $Y_c = YV$. Observe that X_c and Y_c are zero-mean because they are linear transformations of X and Y , which are zero-mean. Thus we can fit a linear model relating the two:

$$Y_c \approx X_c A$$

The least-squares solution is given by

$$\begin{aligned} A &= (X_c^\top X_c)^{-1} X_c^\top Y_c \\ &= (U^\top X^\top X U)^{-1} U^\top X^\top Y V \end{aligned}$$

However, since what we *really* want is an estimate of Y given new (zero-mean) observations \tilde{X} (or vice-versa), it's useful to have the entire series of transformations that relates the two. The predicted canonical variables are given by

$$\hat{Y}_c = \tilde{X}_c A = \tilde{X} U (U^\top X^\top X U)^{-1} U^\top X^\top Y V$$

Then we use the canonical variables to compute the actual values:

$$\begin{aligned} \hat{Y} &= \hat{Y}_c (V^\top V)^{-1} V^\top \\ &= \tilde{X} U (U^\top X^\top X U)^{-1} (U^\top X^\top Y V) (V^\top V)^{-1} V^\top \end{aligned}$$

We can collapse all these terms into a single matrix A_{eq} that gives the prediction \hat{Y} from \tilde{X} :

$$A_{\text{eq}} = \underbrace{U}_{\text{projection}} \underbrace{(U^\top X^\top X U)^{-1}}_{\text{whitening}} \underbrace{(U^\top X^\top Y V)}_{\text{decorrelation}} \underbrace{(V^\top V)^{-1} V^\top}_{\text{projection back}}$$