

MODEL REDUCTION

1. Projection of mappings

The least-squares problem

$$\min_{\mathbf{x} \in \mathbb{R}^n} \|\mathbf{y} - \mathbf{A}\mathbf{x}\| \quad (1)$$

focuses on a simpler representation of a data vector $\mathbf{y} \in \mathbb{R}^m$ as a linear combination of column vectors of $\mathbf{A} \in \mathbb{R}^{m \times n}$. Consider some phenomenon modeled as a function between vector spaces $f: X \rightarrow Y$, such that for input parameters $\mathbf{x} \in X$, the state of the system is $\mathbf{y} = f(\mathbf{x})$. For most models f is differentiable, a transcription of the condition that the system should not exhibit jumps in behavior when changing the input parameters. Then by appropriate choice of units and origin, a linearized model

$$\mathbf{y} = \mathbf{A}\mathbf{x}, \mathbf{A} \in \mathbb{R}^{m \times n},$$

is obtained if $\mathbf{y} \in C(\mathbf{A})$, expressed as (1) if $\mathbf{y} \notin C(\mathbf{A})$.

A simpler description is often sought, typically based on recognition that the inputs and outputs of the model can themselves be obtained as linear combinations $\mathbf{x} = \mathbf{B}\mathbf{u}$, $\mathbf{y} = \mathbf{C}\mathbf{v}$, involving a smaller set of parameters $\mathbf{u} \in \mathbb{R}^q$, $\mathbf{v} \in \mathbb{R}^p$, $p < m$, $q < n$. The column spaces of the matrices $\mathbf{B} \in \mathbb{R}^{n \times q}$, $\mathbf{C} \in \mathbb{R}^{m \times p}$ are vector subspaces of the original set of inputs and outputs, $C(\mathbf{B}) \leq \mathbb{R}^n$, $C(\mathbf{C}) \leq \mathbb{R}^m$. The sets of column vectors of \mathbf{B}, \mathbf{C} each form a *reduced basis* for the system inputs and outputs if they are chosen to be of full rank. The reduced bases are assumed to have been orthonormalized through the Gram-Schmidt procedure such that $\mathbf{B}^T \mathbf{B} = \mathbf{I}_q$, and $\mathbf{C}^T \mathbf{C} = \mathbf{I}_p$. Expressing the model inputs and outputs in terms of the reduced basis leads to

$$\mathbf{C}\mathbf{v} = \mathbf{A}\mathbf{B}\mathbf{u} \Rightarrow \mathbf{v} = \mathbf{C}^T \mathbf{A}\mathbf{B}\mathbf{u} \Rightarrow \mathbf{v} = \mathbf{R}\mathbf{u}.$$

The matrix $\mathbf{R} = \mathbf{C}^T \mathbf{A}\mathbf{B}$ is called the *reduced system matrix* and is associated with a mapping $\mathbf{g}: U \rightarrow V$, that is a restriction to the U, V vector subspaces of the mapping f . When f is an endomorphism, $f: X \rightarrow X$, $m = n$, the same reduced basis is used for both inputs and outputs, $\mathbf{x} = \mathbf{B}\mathbf{u}$, $\mathbf{y} = \mathbf{B}\mathbf{v}$, and the reduced system is

$$\mathbf{v} = \mathbf{R}\mathbf{u}, \mathbf{R} = \mathbf{B}^T \mathbf{A}\mathbf{B}.$$

Since \mathbf{B} is assumed to be orthogonal, the projector onto $C(\mathbf{B})$ is $\mathbf{P}_B = \mathbf{B}\mathbf{B}^T$. Applying the projector on the initial model

$$\mathbf{P}_B \mathbf{y} = \mathbf{P}_B \mathbf{A}\mathbf{x}$$

leads to $\mathbf{B}\mathbf{B}^T \mathbf{y} = \mathbf{B}\mathbf{B}^T \mathbf{A}\mathbf{x}$, and since $\mathbf{v} = \mathbf{B}^T \mathbf{y}$ the relation $\mathbf{B}\mathbf{v} = \mathbf{B}\mathbf{B}^T \mathbf{A}\mathbf{B}\mathbf{u}$ is obtained, and conveniently grouped as

$$\mathbf{B}\mathbf{v} = \mathbf{B}(\mathbf{B}^T \mathbf{A}\mathbf{B})\mathbf{u} \Rightarrow \mathbf{B}\mathbf{v} = \mathbf{B}(\mathbf{R}\mathbf{u}),$$

again leading to the reduced model $\mathbf{v} = \mathbf{B}\mathbf{u}$. The above calculation highlights that the reduced model is a projection of the full model $\mathbf{y} = \mathbf{A}\mathbf{x}$ on $C(\mathbf{B})$.

2. Reduced bases

2.1. Correlation matrices

Correlation coefficient. Consider two functions $x_1, x_2: \mathbb{R} \rightarrow \mathbb{R}$, that represent data streams in time of inputs $x_1(t)$ and outputs $x_2(t)$ of some system. A basic question arising in modeling and data science is whether the inputs and outputs are themselves in a functional relationship. This usually is a consequence of incomplete knowledge of the system, such that while x_1, x_2 might be assumed to be the most relevant input, output quantities, this is not yet fully established. A typical approach is to then carry out repeated measurements leading to a data set $D = \{(x_1(t_i), x_2(t_i)) | i = 1, \dots, N\}$, thus defining a relation. Let $\mathbf{x}_1, \mathbf{x}_2 \in \mathbb{R}^N$ denote vectors containing the input and output values. The *mean values* μ_1, μ_2 of the input and output are estimated by the statistics

$$\mu_1 \cong \bar{x}_1 = \frac{1}{N} \sum_{i=1}^N x_1(t_i) = E[x_1], \mu_2 \cong \bar{x}_2 = \frac{1}{N} \sum_{i=1}^N x_2(t_i) = E[x_2],$$

where E is the expectation seen to be a linear mapping, $E: \mathbb{R}^N \rightarrow \mathbb{R}$ whose associated matrix is

$$\mathbf{E} = \frac{1}{N} [1 \ 1 \ \dots \ 1],$$

and the means are also obtained by matrix vector multiplication (linear combination),

$$\bar{x}_1 = \mathbf{E} \mathbf{x}_1, \bar{x}_2 = \mathbf{E} \mathbf{x}_2.$$

Deviation from the mean is measured by the *standard deviation* defined for x_1, x_2 by

$$\sigma_1 = \sqrt{E[(x_1 - \mu_1)^2]}, \sigma_2 = \sqrt{E[(x_2 - \mu_2)^2]}.$$

Note that the standard deviations are no longer linear mappings of the data.

Assume that the origin is chosen such that $\bar{x}_1 = \bar{x}_2 = 0$. One tool to establish whether the relation D is also a function is to compute the *correlation coefficient*

$$\rho(x_1, x_2) = \frac{E[x_1 x_2]}{\sigma_1 \sigma_2} = \frac{E[x_1 x_2]}{\sqrt{E[x_1^2] E[x_2^2]}}$$

that can be expressed in terms of a scalar product and 2-norm as

$$\rho(x_1, x_2) = \frac{\mathbf{x}_1^T \mathbf{x}_2}{\|\mathbf{x}_1\| \|\mathbf{x}_2\|}.$$

Squaring each side of the norm property $\|\mathbf{x}_1 + \mathbf{x}_2\| \leq \|\mathbf{x}_1\| + \|\mathbf{x}_2\|$, leads to

$$(\mathbf{x}_1 + \mathbf{x}_2)^T (\mathbf{x}_1 + \mathbf{x}_2) \leq \mathbf{x}_1^T \mathbf{x}_1 + \mathbf{x}_2^T \mathbf{x}_2 + 2 \|\mathbf{x}_1\| \|\mathbf{x}_2\| \Rightarrow \mathbf{x}_1^T \mathbf{x}_2 \leq \|\mathbf{x}_1\| \|\mathbf{x}_2\|,$$

known as the Cauchy-Schwarz inequality, which implies $-1 \leq \rho(x_1, x_2) \leq 1$. Depending on the value of ρ , the variables $x_1(t), x_2(t)$ are said to be:

1. *uncorrelated*, if $\rho = 0$;
2. *correlated*, if $\rho = 1$;
3. *anti-correlated*, if $\rho = -1$.

The numerator of the correlation coefficient is known as the covariance of x_1, x_2

$$\text{cov}(x_1, x_2) = E[x_1 x_2].$$

The correlation coefficient can be interpreted as a normalization of the covariance, and the relation

$$\text{cov}(x_1, x_2) = \mathbf{x}_1^T \mathbf{x}_2 = \rho(x_1, x_2) \|\mathbf{x}_1\| \|\mathbf{x}_2\|,$$

is the two-variable version of a more general relationship encountered when the system inputs and outputs become vectors.

Patterns in data. Consider now a related problem, whether the input and output parameters $\mathbf{x} \in \mathbb{R}^n, \mathbf{y} \in \mathbb{R}^m$ thought to characterize a system are actually well chosen, or whether they are redundant in the sense that a more insightful description is furnished by $\mathbf{u} \in \mathbb{R}^q, \mathbf{v} \in \mathbb{R}^p$ with fewer components $p < m, q < n$. Applying the same ideas as in the correlation coefficient, a sequence of N measurements is made leading to data sets

$$\mathbf{X} = [\mathbf{x}_1 \ \mathbf{x}_2 \ \dots \ \mathbf{x}_n] \in \mathbb{R}^{N \times n}, \mathbf{Y} = [\mathbf{y}_1 \ \mathbf{y}_2 \ \dots \ \mathbf{y}_n] \in \mathbb{R}^{N \times m}.$$

Again, by appropriate choice of the origin the means of the above measurements is assumed to be zero

$$E[\mathbf{x}] = \mathbf{0}, E[\mathbf{y}] = \mathbf{0}.$$

Covariance matrices can be constructed by

$$\mathbf{C}_X = \mathbf{X}^T \mathbf{X} = \begin{bmatrix} \mathbf{x}_1^T \\ \mathbf{x}_2^T \\ \vdots \\ \mathbf{x}_n^T \end{bmatrix} [\mathbf{x}_1 \ \mathbf{x}_2 \ \dots \ \mathbf{x}_n] = \begin{bmatrix} \mathbf{x}_1^T \mathbf{x}_1 & \mathbf{x}_1^T \mathbf{x}_2 & \dots & \mathbf{x}_1^T \mathbf{x}_n \\ \mathbf{x}_2^T \mathbf{x}_1 & \mathbf{x}_2^T \mathbf{x}_2 & \dots & \mathbf{x}_2^T \mathbf{x}_n \\ \vdots & \vdots & \ddots & \vdots \\ \mathbf{x}_n^T \mathbf{x}_1 & \mathbf{x}_n^T \mathbf{x}_2 & \dots & \mathbf{x}_n^T \mathbf{x}_n \end{bmatrix} \in \mathbb{R}^{n \times n}.$$

Consider now the SVDs of $C_X = N \Lambda N^T$, $X = U \Sigma S^T$, and from

$$C_X = X^T X = (U \Sigma S^T)^T U \Sigma S^T = S \Sigma^T U^T U \Sigma S^T = S \Sigma^T \Sigma S^T = N \Lambda N^T,$$

identify $N = S$, and $\Lambda = \Sigma^T \Sigma$.

Recall that the SVD returns an order set of singular values $\sigma_1 \geq \sigma_2 \geq \dots \geq$, and associated singular vectors. In many applications the singular values decrease quickly, often exponentially fast. Taking the first q singular modes then gives a basis set suitable for mode reduction

$$\mathbf{x} = \mathbf{S}_q \mathbf{u} = [s_1 \ s_2 \ \dots \ s_q] \mathbf{u}.$$