

# System Modeling, part 2

Marc Claesen

July 9, 2015

## 1 Nonlinear systems & linearization

## 2 System identification (cont)

- Grey box identification
- Black box identification

# Outline

- 1 Nonlinear systems & linearization
- 2 System identification (cont)
  - Grey box identification
  - Black box identification

# Nonlinear systems

In this course we focus on the linear state-space representation:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases} \quad \begin{cases} x[k+1] = Ax[k] + Bu[k], \\ y[k] = Cx[k] + Du[k]. \end{cases}$$

# Nonlinear systems

In this course we focus on the linear state-space representation:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases} \quad \begin{cases} x[k+1] = Ax[k] + Bu[k], \\ y[k] = Cx[k] + Du[k]. \end{cases}$$

Most real life systems involve nonlinearity:

$$\begin{cases} \dot{x}(t) = f(x(t), u(t)), \\ y(t) = g(x(t), u(t)), \end{cases}$$

# Nonlinear systems

In this course we focus on the linear state-space representation:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases} \quad \begin{cases} x[k+1] = Ax[k] + Bu[k], \\ y[k] = Cx[k] + Du[k]. \end{cases}$$

Most real life systems involve nonlinearity:

$$\begin{cases} \dot{x}(t) = f(x(t), u(t)), \\ y(t) = g(x(t), u(t)), \end{cases}$$

where  $f$  and/or  $g$  contain some nonlinearity, such as:

# Nonlinear systems

In this course we focus on the linear state-space representation:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases} \quad \begin{cases} x[k+1] = Ax[k] + Bu[k], \\ y[k] = Cx[k] + Du[k]. \end{cases}$$

Most real life systems involve nonlinearity:

$$\begin{cases} \dot{x}(t) = f(x(t), u(t)), \\ y(t) = g(x(t), u(t)), \end{cases}$$

where  $f$  and/or  $g$  contain some nonlinearity, such as:

- *powers*: e.g.  $\dot{x}(t) = Ax(t) + Bu(t) + \gamma u(t)^2$ ,

# Nonlinear systems

In this course we focus on the linear state-space representation:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases} \quad \begin{cases} x[k+1] = Ax[k] + Bu[k], \\ y[k] = Cx[k] + Du[k]. \end{cases}$$

Most real life systems involve nonlinearity:

$$\begin{cases} \dot{x}(t) = f(x(t), u(t)), \\ y(t) = g(x(t), u(t)), \end{cases}$$

where  $f$  and/or  $g$  contain some nonlinearity, such as:

- *powers*: e.g.  $\dot{x}(t) = Ax(t) + Bu(t) + \gamma u(t)^2$ ,
- *interactions*: e.g.  $\dot{x}(t) = Ax(t) + Bu(t) + \gamma x(t)u(t)$ ,



# Nonlinear systems

In this course we focus on the linear state-space representation:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases} \quad \begin{cases} x[k+1] = Ax[k] + Bu[k], \\ y[k] = Cx[k] + Du[k]. \end{cases}$$

Most real life systems involve nonlinearity:

$$\begin{cases} \dot{x}(t) = f(x(t), u(t)), \\ y(t) = g(x(t), u(t)), \end{cases}$$

where  $f$  and/or  $g$  contain some nonlinearity, such as:

- *powers*: e.g.  $\dot{x}(t) = Ax(t) + Bu(t) + \gamma u(t)^2$ ,
- *interactions*: e.g.  $\dot{x}(t) = Ax(t) + Bu(t) + \gamma x(t)u(t)$ ,
- *clipping*: e.g.  $\alpha \leq x(t) \leq \beta$ .

# Linearization around equilibrium point

# Linearization around equilibrium point

Nonlinear systems have (several) equilibrium points  $x_e$ ,  $u_e$ ,  $y_e$ :

$$\begin{cases} \dot{x}_e = f(x_e, u_e) = 0, \\ y_e = g(x_e, u_e). \end{cases}$$

# Linearization around equilibrium point

Nonlinear systems have (several) equilibrium points  $x_e$ ,  $u_e$ ,  $y_e$ :

$$\begin{cases} \dot{x}_e = f(x_e, u_e) = 0, \\ y_e = g(x_e, u_e). \end{cases}$$

Linearizing in the region of  $(x_e, u_e, y_e)$ :

$$x = x_e + \Delta x, \quad u = u_e + \Delta u, \quad y = y_e + \Delta y,$$

# Linearization around equilibrium point

Nonlinear systems have (several) equilibrium points  $x_e$ ,  $u_e$ ,  $y_e$ :

$$\begin{cases} \dot{x}_e = f(x_e, u_e) = 0, \\ y_e = g(x_e, u_e). \end{cases}$$

Linearizing in the region of  $(x_e, u_e, y_e)$ :

$$x = x_e + \Delta x, \quad u = u_e + \Delta u, \quad y = y_e + \Delta y,$$

with  $\Delta x$ ,  $\Delta u$  and  $\Delta y$  *sufficiently* small.

# Linearization around equilibrium point

Nonlinear systems have (several) equilibrium points  $x_e, u_e, y_e$ :

$$\begin{cases} \dot{x}_e = f(x_e, u_e) = 0, \\ y_e = g(x_e, u_e). \end{cases}$$

Linearizing in the region of  $(x_e, u_e, y_e)$ :

$$x = x_e + \Delta x, \quad u = u_e + \Delta u, \quad y = y_e + \Delta y,$$

with  $\Delta x$ ,  $\Delta u$  and  $\Delta y$  *sufficiently* small.

Linearizing is done via first order Taylor expansions:

$$\begin{cases} \frac{dx}{dt} = \frac{d\Delta x}{dt} = f(x, u) = f(x_e + \Delta x, u_e + \Delta u), \\ y_e + \Delta y = g(x, u) = g(x_e + \Delta x, u_e + \Delta u). \end{cases}$$

## Example: decalcification plant

Used to reduce concentration of calcium hydroxide in water:

- chemical reaction:  $\text{Ca}(\text{OH})_2 + \text{CO}_2 \rightarrow \text{CaCO}_3 + \text{H}_2\text{O}$
- reaction speed:  $r = c[\text{Ca}(\text{OH})_2][\text{CO}_2]$
- rate of change of concentration:

$$\begin{aligned}\frac{d[\text{Ca}(\text{OH})_2]}{dt} &= \frac{k}{V} - \frac{r}{V}, \\ \frac{d[\text{CO}_2]}{dt} &= \frac{u}{V} - \frac{r}{V},\end{aligned}$$

with inflow rates  $k$  and  $u$  in mol/s and tank volume  $V$  in L.

- input  $u$ : inflow of  $\text{CO}_2$ , output:  $[\text{Ca}(\text{OH})_2]$

# Nonlinear model and equilibrium point

**Nonlinear** model for the given reactor:



# Nonlinear model and equilibrium point

**Nonlinear** model for the given reactor:

$$\begin{aligned}\frac{d[Ca(OH)_2]}{dt} &= \frac{k}{V} - \frac{c}{V}[Ca(OH)_2][CO_2], \\ \frac{d[CO_2]}{dt} &= \frac{u}{V} - \frac{c}{V}[Ca(OH)_2][CO_2], \\ y &= [Ca(OH)_2],\end{aligned}$$

with two state variables:  $x_1 = [Ca(OH)_2]$  and  $x_2 = [CO_2]$ .

# Nonlinear model and equilibrium point

**Nonlinear** model for the given reactor:

$$\begin{aligned}\frac{d[\text{Ca}(\text{OH})_2]}{dt} &= \frac{k}{V} - \frac{c}{V} [\text{Ca}(\text{OH})_2][\text{CO}_2], \\ \frac{d[\text{CO}_2]}{dt} &= \frac{u}{V} - \frac{c}{V} [\text{Ca}(\text{OH})_2][\text{CO}_2], \\ y &= [\text{Ca}(\text{OH})_2],\end{aligned}$$

with two state variables:  $x_1 = [\text{Ca}(\text{OH})_2]$  and  $x_2 = [\text{CO}_2]$ .

The equilibrium point  $(k_{eq}, u_{eq}, x_{1,eq}, x_{2,eq}, y_{eq})$  of this system is:

# Nonlinear model and equilibrium point

**Nonlinear** model for the given reactor:

$$\begin{aligned}\frac{d[\text{Ca}(\text{OH})_2]}{dt} &= \frac{k}{V} - \frac{c}{V} [\text{Ca}(\text{OH})_2][\text{CO}_2], \\ \frac{d[\text{CO}_2]}{dt} &= \frac{u}{V} - \frac{c}{V} [\text{Ca}(\text{OH})_2][\text{CO}_2], \\ y &= [\text{Ca}(\text{OH})_2],\end{aligned}$$

with two state variables:  $x_1 = [\text{Ca}(\text{OH})_2]$  and  $x_2 = [\text{CO}_2]$ .

The equilibrium point  $(k_{eq}, u_{eq}, x_{1,eq}, x_{2,eq}, y_{eq})$  of this system is:

$$\begin{aligned}\frac{k_{eq}}{V} - \frac{c}{V} [\text{Ca}(\text{OH})_2]_{eq} [\text{CO}_2]_{eq} &= 0, \\ \frac{u_{eq}}{V} - \frac{c}{V} [\text{Ca}(\text{OH})_2]_{eq} [\text{CO}_2]_{eq} &= 0.\end{aligned}$$

# Linearization of the decalcification plant

For small deviations near the equilibrium:

$$\begin{aligned}\frac{d\Delta x_1}{dt} &= -\frac{c}{V}[CO_2]_{eq}\Delta x_1 - \frac{c}{V}[Ca(OH)_2]_{eq}\Delta x_2, \\ \frac{d\Delta x_2}{dt} &= -\frac{c}{V}[CO_2]_{eq}\Delta x_1 - \frac{c}{V}[Ca(OH)_2]_{eq}\Delta x_2 + \frac{\Delta u}{V}, \\ \Delta y &= \Delta x_1.\end{aligned}$$

# Linearization of the decalcification plant

For small deviations near the equilibrium:

$$\begin{aligned}\frac{d\Delta x_1}{dt} &= -\frac{c}{V}[CO_2]_{eq}\Delta x_1 - \frac{c}{V}[Ca(OH)_2]_{eq}\Delta x_2, \\ \frac{d\Delta x_2}{dt} &= -\frac{c}{V}[CO_2]_{eq}\Delta x_1 - \frac{c}{V}[Ca(OH)_2]_{eq}\Delta x_2 + \frac{\Delta u}{V}, \\ \Delta y &= \Delta x_1.\end{aligned}$$

The resulting linear state-space model is  $\begin{cases} \dot{x}(t) = Ax(t) + Bu(t) \\ y(t) = Cx(t) + Du(t) \end{cases}$ .

# Linearization of the decalcification plant

For small deviations near the equilibrium:

$$\begin{aligned}\frac{d\Delta x_1}{dt} &= -\frac{c}{V}[CO_2]_{eq}\Delta x_1 - \frac{c}{V}[Ca(OH)_2]_{eq}\Delta x_2, \\ \frac{d\Delta x_2}{dt} &= -\frac{c}{V}[CO_2]_{eq}\Delta x_1 - \frac{c}{V}[Ca(OH)_2]_{eq}\Delta x_2 + \frac{\Delta u}{V}, \\ \Delta y &= \Delta x_1.\end{aligned}$$

The resulting linear state-space model is  $\begin{cases} \dot{x}(t) = Ax(t) + Bu(t) \\ y(t) = Cx(t) + Du(t) \end{cases}$ :

$$\begin{bmatrix} \frac{d[Ca(OH)_2]}{dt} \\ \frac{d[CO_2]}{dt} \end{bmatrix} = - \begin{bmatrix} \frac{c}{V}[CO_2]_{eq} & \frac{c}{V}[Ca(OH)_2]_{eq} \\ \frac{c}{V}[CO_2]_{eq} & \frac{c}{V}[Ca(OH)_2]_{eq} \end{bmatrix} \begin{bmatrix} [Ca(OH)_2] \\ [CO_2] \end{bmatrix} + \begin{bmatrix} 0 \\ \frac{1}{V} \end{bmatrix} u(t)$$

$$y(t) = [Ca(OH)_2]$$

# Outline

## 1 Nonlinear systems & linearization

## 2 System identification (cont)

- Grey box identification
- Black box identification

# Main classes of identification methods

**White box modeling:** based on first principles.



# Main classes of identification methods

**White box modeling:** based on first principles.

→ known equations (structure) & parameters (coefficients).

# Main classes of identification methods

**White box modeling:** based on first principles.

→ known equations (structure) & parameters (coefficients).

**Grey box identification:** first principles & experimentation.

# Main classes of identification methods

**White box modeling:** based on first principles.

→ known equations (structure) & parameters (coefficients).

**Grey box identification:** first principles & experimentation.

→ known equations, unknown/uncertain parameters.

# Main classes of identification methods

**White box modeling:** based on first principles.

→ known equations (structure) & parameters (coefficients).

**Grey box identification:** first principles & experimentation.

→ known equations, unknown/uncertain parameters.

**Black box identification:** based on experimentation.

# Main classes of identification methods

**White box modeling:** based on first principles.

→ known equations (structure) & parameters (coefficients).

**Grey box identification:** first principles & experimentation.

→ known equations, unknown/uncertain parameters.

**Black box identification:** based on experimentation.

→ unknown equations & unknown parameters.

# Main classes of identification methods

**White box modeling:** based on first principles.

→ known equations (structure) & parameters (coefficients).

**Grey box identification:** first principles & experimentation.

→ known equations, unknown/uncertain parameters.

**Black box identification:** based on experimentation.

→ unknown equations & unknown parameters.

Most popular approaches are forms of black box identification.

# Outline

- 1 Nonlinear systems & linearization
- 2 System identification (cont)
  - Grey box identification
  - Black box identification

# Grey box identification: conceptual



# Grey box identification: conceptual

Grey box identification starts from a known model structure but with unknown/uncertain parameters  $\leftrightarrow$  **parametric statistics**.

# Grey box identification: conceptual

Grey box identification starts from a known model structure but with unknown/uncertain parameters  $\leftrightarrow$  **parametric statistics**.

We assume linear, continuous time state space representation:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases}$$

# Grey box identification: conceptual

Grey box identification starts from a known model structure but with unknown/uncertain parameters  $\leftrightarrow$  **parametric statistics**.

We assume linear, continuous time state space representation:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases}$$

**Given:** states, inputs, outputs and guesstimates of  $\tilde{A}$ ,  $\tilde{B}$ ,  $\tilde{C}$  &  $\tilde{D}$ .

# Grey box identification: conceptual

Grey box identification starts from a known model structure but with unknown/uncertain parameters  $\leftrightarrow$  **parametric statistics**.

We assume linear, continuous time state space representation:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases}$$

**Given:** states, inputs, outputs and guesstimates of  $\tilde{A}$ ,  $\tilde{B}$ ,  $\tilde{C}$  &  $\tilde{D}$ .

**Task:** estimate  $\hat{A}$ ,  $\hat{B}$ ,  $\hat{C}$  and  $\hat{D}$  adequately via experiments.

# Grey box identification: conceptual

Grey box identification starts from a known model structure but with unknown/uncertain parameters  $\leftrightarrow$  **parametric statistics**.

We assume linear, continuous time state space representation:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases}$$

**Given:** states, inputs, outputs and guesstimates of  $\tilde{A}$ ,  $\tilde{B}$ ,  $\tilde{C}$  &  $\tilde{D}$ .

**Task:** estimate  $\hat{A}$ ,  $\hat{B}$ ,  $\hat{C}$  and  $\hat{D}$  adequately via experiments.

*"All models are wrong, but some are useful."* – George E. P. Box

# Linear regression

Consider input matrix  $\mathbf{X}$ , output vector  $\mathbf{y}$  and residuals  $\epsilon$ :

$$\mathbf{X}\theta = \mathbf{y} + \epsilon.$$

The parameter vector  $\theta$  must be estimated, given the observations.

# Linear regression

Consider input matrix  $\mathbf{X}$ , output vector  $\mathbf{y}$  and residuals  $\epsilon$ :

$$\mathbf{X}\theta = \mathbf{y} + \epsilon.$$

The parameter vector  $\theta$  must be estimated, given the observations.

A common estimation approach is ordinary least squares (OLS):

$$\begin{aligned}(\mathbf{X}^T \mathbf{X}) \hat{\theta}_{OLS} &= \mathbf{X}^T \mathbf{y}, \\ \hat{\theta}_{OLS} &= (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{y}.\end{aligned}$$

# Linear regression

Consider input matrix  $\mathbf{X}$ , output vector  $\mathbf{y}$  and residuals  $\epsilon$ :

$$\mathbf{X}\theta = \mathbf{y} + \epsilon.$$

The parameter vector  $\theta$  must be estimated, given the observations.

A common estimation approach is ordinary least squares (OLS):

$$\begin{aligned}(\mathbf{X}^T \mathbf{X}) \hat{\theta}_{OLS} &= \mathbf{X}^T \mathbf{y}, \\ \hat{\theta}_{OLS} &= (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{y}.\end{aligned}$$

The OLS estimate minimizes the sum-of-squares of errors, i.e.:

$$\hat{\theta}_{OLS} = \arg \min_{\theta} \sum_{i=1}^N \left( y(i) - \sum_{j=1}^d X(i,j) \theta(j) \right)^2$$



# Linear regression with ordinary least squares

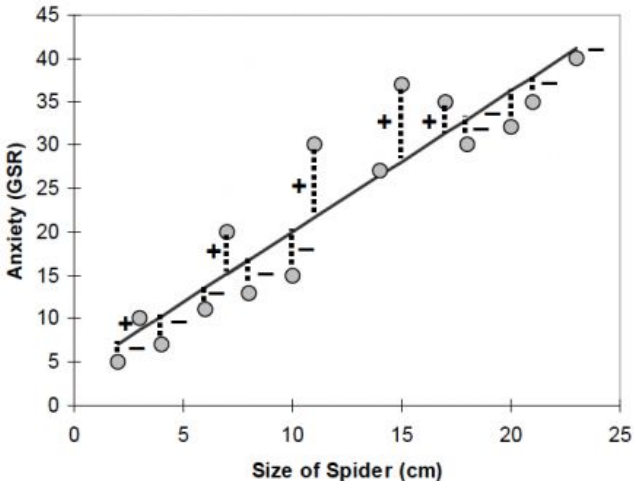


Image taken from <http://freakonometrics.hypotheses.org/2348>.

# Maximum likelihood estimation

The maximum likelihood estimate  $\hat{\theta}_{ML}$  is the parameter vector that maximizes the likelihood  $\mathcal{L}(\cdot)$  of observing the (known) outputs  $\mathbf{y}$ , given the (known) inputs  $\mathbf{X}$ :

$$\hat{\theta}_{ML} = \arg \max_{\theta} \mathcal{L}(\mathbf{y}, \mathbf{X} \mid \theta)$$

# Maximum likelihood estimation

The maximum likelihood estimate  $\hat{\theta}_{ML}$  is the parameter vector that maximizes the likelihood  $\mathcal{L}(\cdot)$  of observing the (known) outputs  $\mathbf{y}$ , given the (known) inputs  $\mathbf{X}$ :

$$\hat{\theta}_{ML} = \arg \max_{\theta} \mathcal{L}(\mathbf{y}, \mathbf{X} \mid \theta)$$

For some structures, ML estimate can be obtained in closed form.

# Maximum likelihood estimation

The maximum likelihood estimate  $\hat{\theta}_{ML}$  is the parameter vector that maximizes the likelihood  $\mathcal{L}(\cdot)$  of observing the (known) outputs  $\mathbf{y}$ , given the (known) inputs  $\mathbf{X}$ :

$$\hat{\theta}_{ML} = \arg \max_{\theta} \mathcal{L}(\mathbf{y}, \mathbf{X} \mid \theta)$$

For some structures, ML estimate can be obtained in closed form.

**Example:** least squares estimators are the maximum likelihood estimators if the associated residuals  $\epsilon$  are normally distributed.

# Maximum a posteriori (MAP) estimation

Bayesian: maximum likelihood estimation with a *prior*  $p(\theta)$ .  
→ MAP estimation is a regularization of ML estimation

# Maximum a posteriori (MAP) estimation

Bayesian: maximum likelihood estimation with a *prior*  $p(\theta)$ .

→ MAP estimation is a regularization of ML estimation

Bayes' theorem:  $P(A | B) = P(B | A) \cdot P(A) / P(B)$ .

# Maximum a posteriori (MAP) estimation

Bayesian: maximum likelihood estimation with a *prior*  $p(\theta)$ .

→ MAP estimation is a regularization of ML estimation

Bayes' theorem:  $P(A | B) = P(B | A) \cdot P(A) / P(B)$ .

If a prior distribution  $p(\cdot)$  is available for  $\theta$ , then the posterior distribution for  $\theta$  becomes:

$$\theta \mapsto \mathcal{L}(\theta | \mathbf{y}, \mathbf{X}) = \frac{\mathcal{L}(\mathbf{y}, \mathbf{X} | \theta)p(\theta)}{\int_{\vartheta} \mathcal{L}(\mathbf{y}, \mathbf{X} | \vartheta)p(\vartheta)d\vartheta}.$$

# Maximum a posteriori (MAP) estimation

Bayesian: maximum likelihood estimation with a *prior*  $p(\theta)$ .

→ MAP estimation is a regularization of ML estimation

Bayes' theorem:  $P(A | B) = P(B | A) \cdot P(A) / P(B)$ .

If a prior distribution  $p(\cdot)$  is available for  $\theta$ , then the posterior distribution for  $\theta$  becomes:

$$\theta \mapsto \mathcal{L}(\theta | \mathbf{y}, \mathbf{X}) = \frac{\mathcal{L}(\mathbf{y}, \mathbf{X} | \theta)p(\theta)}{\int_{\vartheta} \mathcal{L}(\mathbf{y}, \mathbf{X} | \vartheta)p(\vartheta)d\vartheta}.$$

The MAP estimate is the mode of the posterior distribution of  $\theta$ :

$$\hat{\theta}_{MAP} = \arg \max_{\theta} \mathcal{L}(\mathbf{y}, \mathbf{X} | \theta)p(\theta).$$



# Errors-in-variables approach

Additionally accounts for **measurement errors in inputs**.

↔ standard regression only accounts for *errors in outputs*

# Errors-in-variables approach

Additionally accounts for **measurement errors in inputs**.

↔ standard regression only accounts for **errors in outputs**

Typically described via *latent variables*:

$$\begin{cases} x = x^* + \eta, \\ y = y^* + \epsilon, \\ y^* = g(x^* \mid \theta), \end{cases}$$

with  $x$ ,  $y$  the observed inputs, outputs and latent variables  $x^*$ ,  $y^*$ .

# Errors-in-variables approach

Additionally accounts for **measurement errors in inputs**.

↔ standard regression only accounts for **errors in outputs**

Typically described via *latent variables*:

$$\begin{cases} x = x^* + \eta, \\ y = y^* + \epsilon, \\ y^* = g(x^* \mid \theta), \end{cases}$$

with  $x$ ,  $y$  the observed inputs, outputs and latent variables  $x^*$ ,  $y^*$ .

**Assumption:** latent variables  $x^*$  and  $y^*$  exist which follow the true functional relationship  $g(\cdot)$ .

# Errors-in-variables approach

Additionally accounts for **measurement errors in inputs**.

↔ standard regression only accounts for **errors in outputs**

Typically described via *latent variables*:

$$\begin{cases} x = x^* + \eta, \\ y = y^* + \epsilon, \\ y^* = g(x^* \mid \theta), \end{cases}$$

with  $x, y$  the observed inputs, outputs and latent variables  $x^*, y^*$ .

**Assumption:** latent variables  $x^*$  and  $y^*$  exist which follow the true functional relationship  $g(\cdot)$ .

**Task:** estimate  $\theta$ .

# Outline

- 1 Nonlinear systems & linearization
- 2 System identification (cont)
  - Grey box identification
  - Black box identification

# Black box identification

Start from unknown equations & unknown parameters.

→ related to **machine learning** and **nonparametric statistics**.

# Black box identification

Start from unknown equations & unknown parameters.

→ related to **machine learning** and **nonparametric statistics**.

If we assume a linear state space system:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases} \quad \begin{cases} x[k+1] = Ax[k] + Bu[k], \\ y[k] = Cx[k] + Du[k]. \end{cases}$$

# Black box identification

Start from unknown equations & unknown parameters.

→ related to **machine learning** and **nonparametric statistics**.

If we assume a linear state space system:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases} \quad \begin{cases} x[k+1] = Ax[k] + Bu[k], \\ y[k] = Cx[k] + Du[k]. \end{cases}$$

Black box identification deals with:



# Black box identification

Start from unknown equations & unknown parameters.

→ related to **machine learning** and **nonparametric statistics**.

If we assume a linear state space system:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases} \quad \begin{cases} x[k+1] &= Ax[k] + Bu[k], \\ y[k] &= Cx[k] + Du[k]. \end{cases}$$

Black box identification deals with:

- unknown states, both in number & physical interpretation

# Black box identification

Start from unknown equations & unknown parameters.

→ related to **machine learning** and **nonparametric statistics**.

If we assume a linear state space system:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases} \quad \begin{cases} x[k+1] &= Ax[k] + Bu[k], \\ y[k] &= Cx[k] + Du[k]. \end{cases}$$

Black box identification deals with:

- unknown states, both in number & physical interpretation  
→ dimensions of  $A$ ,  $B$  &  $C$  unknown

# Black box identification

Start from unknown equations & unknown parameters.

→ related to **machine learning** and **nonparametric statistics**.

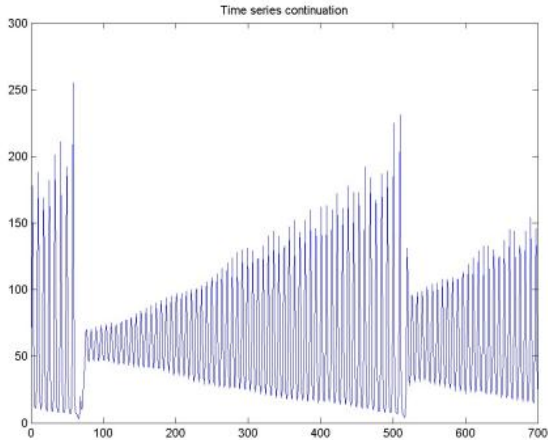
If we assume a linear state space system:

$$\begin{cases} \dot{x}(t) = Ax(t) + Bu(t), \\ y(t) = Cx(t) + Du(t). \end{cases} \quad \begin{cases} x[k+1] = Ax[k] + Bu[k], \\ y[k] = Cx[k] + Du[k]. \end{cases}$$

Black box identification deals with:

- unknown states, both in number & physical interpretation  
→ dimensions of  $A$ ,  $B$  &  $C$  unknown
- unknown parameters (values in  $A$ ,  $B$ ,  $C$ ,  $D$ )

# Time series: Santa Fe laser



# Modelling the Santa Fe laser

This laser can be treated as an autonomous discrete time system:

$$\begin{cases} x[k+1] = f(x[k-N+1], \dots, x[k]), \\ y[k] = x[k]. \end{cases}$$

The output depends on the past  $N$  states & no inputs.

# Modelling the Santa Fe laser

This laser can be treated as an autonomous discrete time system:

$$\begin{cases} x[k+1] = f(x[k-N+1], \dots, x[k]), \\ y[k] = x[k]. \end{cases}$$

The output depends on the past  $N$  states & no inputs.

→ how large is  $N$ ? → **unknown structure**

# Modelling the Santa Fe laser

This laser can be treated as an autonomous discrete time system:

$$\begin{cases} x[k+1] = f(x[k-N+1], \dots, x[k]), \\ y[k] = x[k]. \end{cases}$$

The output depends on the past  $N$  states & no inputs.

→ how large is  $N$ ? → **unknown structure**

Treat it as a regression problem with  $N$  inputs:  $y = f(X_1, \dots, X_N)$ .

# Modelling the Santa Fe laser

This laser can be treated as an autonomous discrete time system:

$$\begin{cases} x[k+1] = f(x[k-N+1], \dots, x[k]), \\ y[k] = x[k]. \end{cases}$$

The output depends on the past  $N$  states & no inputs.

→ how large is  $N$ ? → **unknown structure**

Treat it as a regression problem with  $N$  inputs:  $y = f(X_1, \dots, X_N)$ .

→ lets say linear, i.e.  $y = \mathbf{X}\theta$  → **unknown parameters**  $\theta \in \mathbb{R}^N$ .



# Modelling the Santa Fe laser

This laser can be treated as an autonomous discrete time system:

$$\begin{cases} x[k+1] = f(x[k-N+1], \dots, x[k]), \\ y[k] = x[k]. \end{cases}$$

The output depends on the past  $N$  states & no inputs.

→ how large is  $N$ ? → **unknown structure**

Treat it as a regression problem with  $N$  inputs:  $y = f(X_1, \dots, X_N)$ .

→ lets say linear, i.e.  $y = \mathbf{X}\theta$  → **unknown parameters**  $\theta \in \mathbb{R}^N$ .

→ for given  $N$ , we can estimate  $\theta$  via grey box methods.

# Modelling the Santa Fe laser

This laser can be treated as an autonomous discrete time system:

$$\begin{cases} x[k+1] = f(x[k-N+1], \dots, x[k]), \\ y[k] = x[k]. \end{cases}$$

The output depends on the past  $N$  states & no inputs.

→ how large is  $N$ ? → **unknown structure**

Treat it as a regression problem with  $N$  inputs:  $y = f(X_1, \dots, X_N)$ .

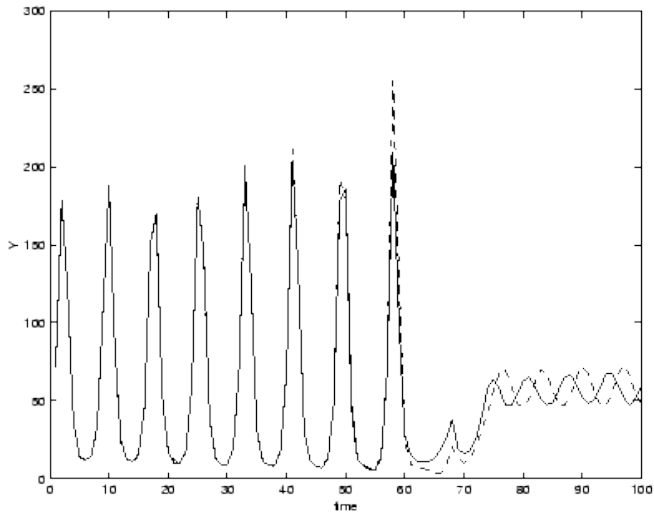
→ lets say linear, i.e.  $y = \mathbf{X}\theta$  → **unknown parameters**  $\theta \in \mathbb{R}^N$ .

→ for given  $N$ , we can estimate  $\theta$  via grey box methods.

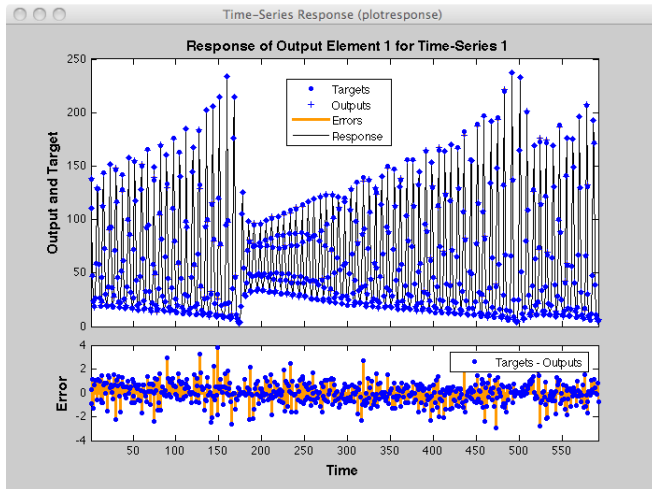
Nonlinear models can be obtained via machine learning methods.

→ neural networks, support vector machine, random forest, ...

# Predictions of a least-squares support vector machine



# Predictions of an artificial neural network



# Neural network: biological

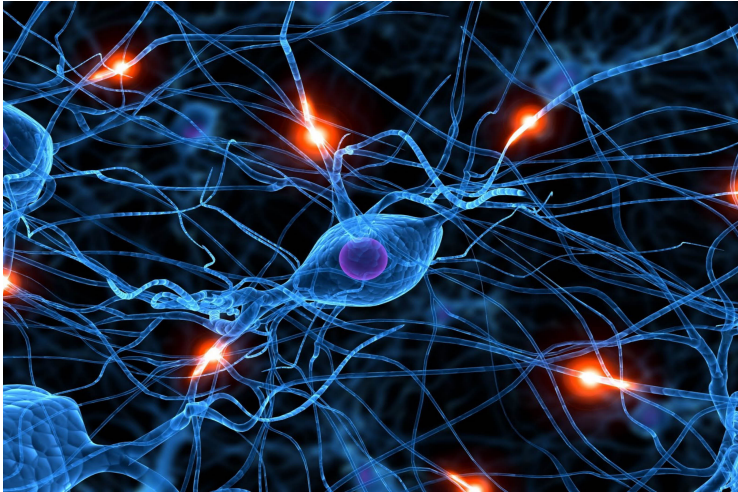


Image taken from <http://www.extremetech.com/wp-content/uploads/2013/09/340.jpg>.

# Structure of a single neuron

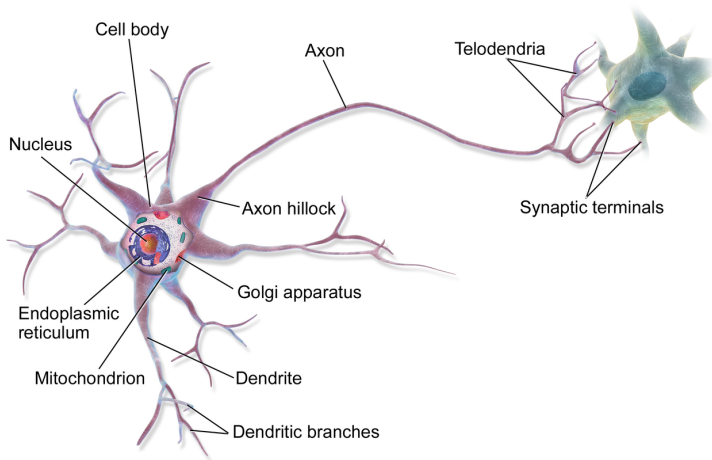


Image taken from [http://en.wikipedia.org/wiki/File:Blausen\\_0657\\_MultipolarNeuron.png](http://en.wikipedia.org/wiki/File:Blausen_0657_MultipolarNeuron.png).

# Neural network: artificial

