System Identification for Control and Simulation.



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Two goals

A general overview of basic steps in system identification





Two goals

- A general overview of basic steps in system identification
- A more technical account of modeling goals for linear systems



Modeling Approaches: View from the Mathworks 3(27)



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The System and the Model



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Pitch rate, Canard, Elevator, Leading Edge Flap

How do the control surface angles affect the pitch rate?

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Pitch rate, Canard, Elevator, Leading Edge Flap

- How do the control surface angles affect the pitch rate?
- Simulation of the aircraft

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Pitch rate, Canard, Elevator, Leading Edge Flap

- How do the control surface angles affect the pitch rate?
- Simulation of the aircraft
- Design of autopilot (regulator)

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y(t) pitch rate at time t. $u_1(t)$ canard angle at time t. T = 1/60. Try

$$y(t) = +b_1u_1(t-T) + b_2u_1(t-2T) + b_3u_1(t-3T) + b_4u_1(t-4T)$$



Dashed line: Measured output (Pitch rate). Solid line: Model output, simulated from the fourth order model from canard angle only.

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Using All Inputs

 u_1 canard angle; u_2 Elevator angle; u_3 Leading edge flap;

$$y(t) = -a_1 y(t - T) - a_2 y(t - 2T) - a_3 y(t - 3T) - a_4 y(t - 4T) + b_1^1 u_1(t - T) + \dots + b_1^4 u_1(t - 4T) + b_2^1 u_2(t - T) + \dots + b_1^3 u_3(t - T) + \dots + b_4^3 u_3(t - 4T)$$



Dashed line: Measured output (Pitch rate). Solid line: Model output, simulated from the fourth order model with all three inputs

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- Select a class of candidate models
- Select a member in this class using the observed data
- Evaluate the quality of the obtained model
- Design the experiment so that the model will be "good".



A Typical Problem

Given Observed Input-Output Data: Find a Description of the System that Generated the Data.





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Techniques

Estimate the parameters by ML techniques/PEM (prediction error methods). Find the model structure by Cross Validation or other validation techniques

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 \mathcal{X} : The Experiment \mathcal{D} : The Measured Data \mathcal{M} : The Model Set \mathcal{I} : The Identification Method \mathcal{V} : The Validation Procedure

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The SI Flow; Model Structures ${\cal M}$



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- They can come in many different forms
- Individual models in the structure are labeled with a parameter vector θ
- A common framework is to describe the model as a predictor of the next output, based on observations of past input-output data.

Observed input–output (u, y) data up to time $t: Z^t$ Model described by predictor: $\mathcal{M}(\theta) : \hat{y}(t|\theta) = g(t, \theta, Z^{t-1}).$

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The SI Flow: Estimation: ${\cal I}$





If a model, $\hat{y}(t|\theta)$, essentially is a predictor of the next output, is is natural to evaluate its quality by assessing how well it predicts: Form the *Prediction error* and measure its size:

 $\varepsilon(t,\theta) = y(t) - \hat{y}(t|\theta), \quad \ell(\varepsilon(t,\theta))$





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$$V_N(\theta) = \sum_{t=1}^N \varepsilon^2(t,\theta)$$



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Which model in the structure performed best (Prediction Error Method, PEM)?

$$\hat{ heta}_N = rg\min_{ heta \in D_{\mathcal{M}}} V_N(heta)$$

(This is often also the Maximum Likelihood Estimate (MLE).)

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Model Estimate Properties $\mathcal{M}(\hat{\theta}_N)$





$$\hat{\theta}_N \to \theta^* \sim \arg\min_{\theta} E\varepsilon^2(t,\theta)$$

 $\blacksquare \ \mathcal{M}(\theta^*)$ is the best possible predictor in \mathcal{M}

E: Expectation.





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- The model structure is not large enough: The ML/PEM estimate converges to the best possible approximation of the system.
- "Best possible approximation" ...
- ... under the conditions of the experiment
- (If the model structure is large enough to contain a true description of the system, then the ML/PEM estimated model has (asymptotically) the best accuracy).





Experiment Design: \mathcal{X}





 $\ensuremath{\mathcal{X}}$: The design variables: Input, Sampling Interval, Feedback,... Then we just saw

$$\hat{\theta}_N \to \theta^*(\mathcal{X})$$



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■ The model $\mathcal{M}(\theta^*(\mathcal{X}))$ is the best approximation of the system under \mathcal{X}

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$$\hat{\theta}_N \to \theta^*(\mathcal{X})$$

- The model $\mathcal{M}(\theta^*(\mathcal{X}))$ is the best approximation of the system under \mathcal{X}
- Let the experimental conditions resemble those under which the model is to be used!



Model Validation: $\!\mathcal{V}$





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Essentially a subjective decision. Several basic techniques are available:

- 1. Simulation (prediction) cross validation
 - Check how well the estimated model can reproduce new, validation data Can compare different models in that way.





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Essentially a subjective decision. Several basic techniques are available:

- 1. Simulation (prediction) cross validation
 - Check how well the estimated model can reproduce new, validation data Can compare different models in that way.
- 2. Residual Analysis
 - Are the residuals $\varepsilon(t, \hat{\theta}_N)$ (the "leftovers") unpredictable? They should not be correlated with anything we knew when estimating the model. Check correlation of the residuals with old inputs.





So this is the System Identification Flow or Loop

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- Several essential choices that have to be made, and often revised.
- Many of the choices have to be taken with the intended model use in mind and thus have a subjective flavour.
- Let us now turn to some specific illustrations for linear models (Requires some more mathematical background.)





$$y(t) = G(q,\theta)u(t) + v(t); \quad G(q,\theta)u(t) = \sum_{k=1}^{\infty} g_k u(t-k),$$

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$$v$$
 has spectrum $\Phi_v(\omega) = \lambda |H(e^{i\omega}, \theta)|^2$
 $v(t) = H(q, \theta)e(t) \quad e(t)$ white noise

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General Description of a Linear Model

 $y(t) = G(q, \theta)u(t) + H(q, \theta)e(t)$

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General Description of a Linear Model

$$y(t) = G(q, \theta)u(t) + H(q, \theta)e(t)$$

The Prediction Errors

$$\varepsilon(t,\theta) = H^{-1}(q,\theta)[y(t) - G(q,\theta)u(t)]$$

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Suppose the data is generated by a true linear system $G_0(q)$ and that the prediction errors are pre-filtered by a filter L(q),

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Then

$$\hat{\theta}_{N} \to \theta^{*} = \arg\min_{\theta} \int_{-\pi}^{\pi} |G(e^{i\omega}, \theta) - G_{0}(e^{i\omega})|^{2}Q(\omega)d\omega$$
$$Q(\omega) = \frac{|L(e^{i\omega})|^{2}\Phi_{u}(\omega)}{|H(e^{i\omega}, \theta)|^{2}}$$

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$$\begin{split} \hat{\theta}_N &\to \theta^* = \arg\min_{\theta} \int_{-\pi}^{\pi} |G(e^{i\omega}, \theta) - G_0(e^{i\omega})|^2 Q(\omega) d\omega \\ Q(\omega) &= \frac{|L(e^{i\omega})|^2 \Phi_u(\omega)}{|H(e^{i\omega}, \theta)|^2} \end{split}$$

So the resulting model is closest to the true system in a norm defined by Q.

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Note that we can affect $Q(\omega) = \frac{|L(e^{i\omega})|^2 \Phi_u(\omega)}{|H(e^{i\omega},\theta)|^2}$ by choosing $L - -(\mathcal{I})$, the input spectrum $\Phi_u - -(\mathcal{X})$ and the noise model $H - -(\mathcal{M})$, so it depends on all boxes in the identification chart.



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■ Identification for simulation with an input $u_* \Rightarrow$ Make Q equal to the spectrum of u_*



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- Identification for simulation with an input $u_* \Rightarrow$ Make Q equal to the spectrum of u_*
- Identification for control ⇒ Make Q large at the intended cross-over frequency (≈ the intended bandwidth)



Identification is a work-flow loop with nodes that contain essential user choices





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- Identification is a work-flow loop with nodes that contain essential user choices
- These may (should) depend on the intended use of the final model
- Concrete illustration for linear system models

