

Design of optimal identification experiments

DISC course: System Identification for control

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1. Reminders and introduction

Consider a real-life system \mathcal{S} operated in open loop:

$$y(t) = G_0(z)u(t) + H_0(z)e(t) = G(z, \theta_0)u(t) + H(z, \theta_0)e(t)$$

with θ_0 the unknown true parameter vector (z^{-1} delay operator)

Consider also a full-order model structure \mathcal{M} for this true system \mathcal{S} ($\mathcal{S} \in \mathcal{M}$):

$$\mathcal{M} = \{ G(z, \theta); H(z, \theta) \mid \theta \in \mathbb{R}^k \}$$

Prediction error identification of \mathcal{S}

An input signal $u(t)$ ($t = 1..N$) is applied to \mathcal{S} and the corresponding output $y(t)$ is collected

$$Z^N = \{y(t) \ u(t) \mid t = 1..N\}$$

Based on these input-output data, prediction error identification can be used to obtain a consistent estimate $\hat{\theta}_N$ of θ_0 :

$$\hat{\theta}_N = \arg \min_{\theta} \frac{1}{N} \sum_{t=1}^N \left(\underbrace{H(z, \theta)^{-1} (y(t) - G(z, \theta)u(t))}_{=\epsilon(t, \theta)} \right)^2$$

Disturbance $v(t) = H_0(z)e(t) \implies \hat{\theta}_N \neq \theta_0$

Question: how can we evaluate the accuracy of $\hat{\theta}_N$ and relate it to the experimental conditions?

Choice of experimental conditions

1. the number N of data that will be collected
2. the signal $u(t)$ that will be applied

Properties of the identified parameter vector $\hat{\theta}_N$

$\hat{\theta}_N$ is (asymptotically) normally distributed around θ_0

$$\sqrt{N}(\hat{\theta}_N - \theta_0) \sim \mathcal{N}(0, P_\theta)$$

$$P_\theta = \sigma_e^2 \left(\bar{E} \psi(t, \theta_0) \psi^T(t, \theta_0) \right)^{-1}$$

with $\psi(t, \theta) = -\frac{\partial \varepsilon(t, \theta)}{\partial \theta}$

The covariance matrix $\text{cov}(\hat{\theta}_N)$ of $\hat{\theta}_N$ given by $\frac{P_\theta}{N}$ is a **measure of accuracy for $\hat{\theta}_N$**

The variance $cov(\hat{\theta}_N)$ in the parameter space induces a variance $cov(G(e^{j\omega}, \hat{\theta}_N))$ in the transfer function space

$$cov(G(e^{j\omega}, \hat{\theta}_N)) = \Lambda_G^*(e^{j\omega}, \theta_0) \frac{P_\theta}{N} \Lambda_G(e^{j\omega}, \theta_0)$$

with $\Lambda_G(z, \theta) = \frac{\partial G(z, \theta)}{\partial \theta}$

$cov(G(e^{j\omega}, \hat{\theta}_N))$ is a measure of accuracy for the identified model $G(z, \hat{\theta}_N)$

Bounding the modeling error using $cov(G(e^{j\omega}, \hat{\theta}_N))$

Using $cov(G(e^{j\omega}, \hat{\theta}_N))$, we can build a confidence region for the modeling error at each frequency:

$$|G(e^{j\omega}, \hat{\theta}_N) - G(e^{j\omega}, \theta_0)| < \alpha \sqrt{cov(G(e^{j\omega}, \hat{\theta}_N))}$$

The bound holds with an user-chosen probability that can be tuned via $\alpha \in \mathbb{R}$ (e.g. the probability is 0.95 if $\alpha = 2.45$)

Remark. $\text{cov}(G(e^{j\omega}, \hat{\theta}_N))$ can be estimated after the identification experiment based on Z^N and $\hat{\theta}_N$

$$\widehat{\text{cov}}(G(e^{j\omega}, \hat{\theta}_N)) = \Lambda_G^*(e^{j\omega}, \hat{\theta}_N) \frac{\hat{P}_\theta}{N} \Lambda_G(e^{j\omega}, \hat{\theta}_N)$$

with

$$\hat{P}_\theta = \hat{\sigma}_e^2 \left(\frac{1}{N} \sum_{t=1}^N \psi(t, \hat{\theta}_N) \psi^T(t, \hat{\theta}_N) \right)^{-1}$$

$$\hat{\sigma}_e^2 = \frac{1}{N} \sum_{t=1}^N \epsilon^2(t, \hat{\theta}_N)$$

Modeling error and experimental conditions i.e. N and $u(t)$

The bound on the modeling error can be expressed as follows:

$$\alpha \sqrt{\text{cov}(G(e^{j\omega}, \hat{\theta}_N))} = \alpha \sqrt{\Lambda_G^*(e^{j\omega}) \frac{P_\theta}{N} \Lambda_G(e^{j\omega})}$$

$$\text{with } P_\theta = \sigma_e^2 \left(\bar{E} \psi(t, \theta_0) \psi^T(t, \theta_0) \right)^{-1}$$

$$\psi(t, \theta_0) = F_u(z, \theta_0) \mathbf{u}(t) + F_e(z, \theta_0) e(t)$$

$$F_u(z, \theta_0) = \frac{1}{H_0} \left. \frac{\partial G(z, \theta)}{\partial \theta} \right|_{\theta=\theta_0} \quad \text{and} \quad F_e(z, \theta_0) = \frac{1}{H_0} \left. \frac{\partial H(z, \theta)}{\partial \theta} \right|_{\theta=\theta_0}$$

Important observations

the experimental conditions N and $u(t)$ influence (the bound on) the modeling error

the larger N , the smaller the modeling error

For a given frequency content, the most powerful $u(t)$, the smaller the modeling error

For signals $u(t)$ having the same power, different frequency contents could lead to different modeling errors

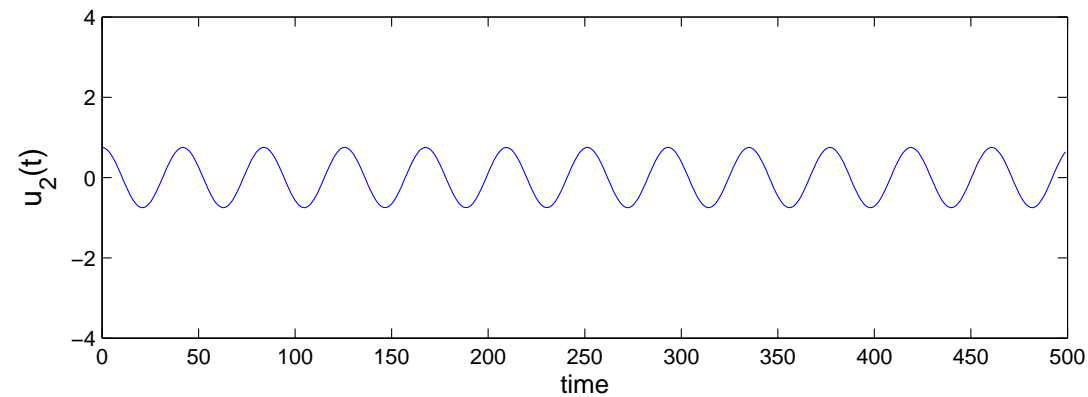
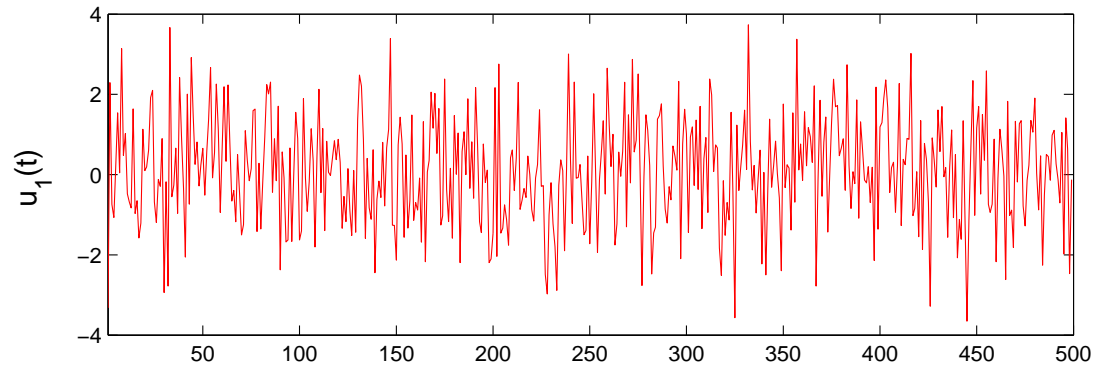
Example

$$\mathcal{S} : y(t) = \frac{3.6z^{-1}}{1 - 0.7z^{-1}}u(t) + (1 + 0.9z^{-1})e(t)$$

$$\mathcal{M} : G(z, \theta) = \frac{bz^{-1}}{1 - fz^{-1}} \quad H(z, \theta) = 1 + cz^{-1} \quad \theta = \begin{pmatrix} b \\ c \\ f \end{pmatrix}$$

Let us compare the accuracy obtained with two input signals: a white noise and a cosine at $\omega = 0.15$

$u_1(t)$ a white noise with variance 1.7 and
 $u_2(t) = 0.75\cos(0.15t)$; both of length $N = 500$



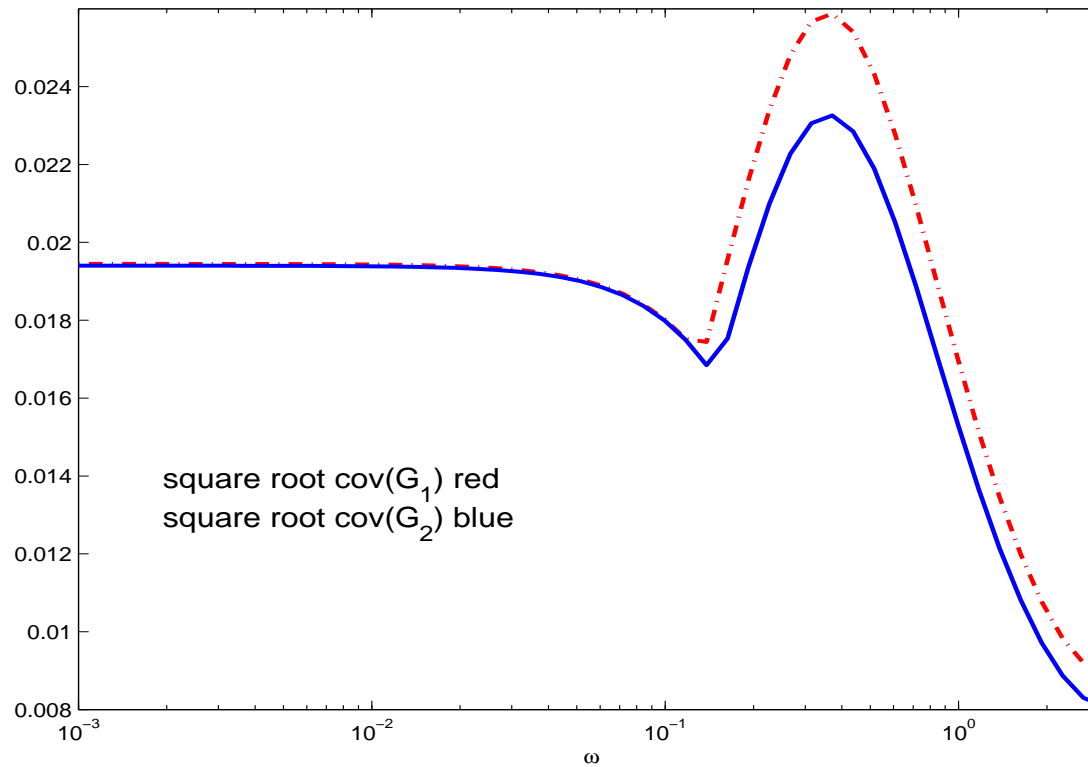
$$\mathcal{P}_{u_1} \gg \mathcal{P}_{u_2}$$

Question:

Which signal leads to the most accurate model?

While $\mathcal{P}_{u_1} \gg \mathcal{P}_{u_2}$, the model \hat{G}_2 identified with u_2 has a slightly better accuracy than the model \hat{G}_1 identified with u_1

$$\sqrt{\text{cov}(\hat{G}_1(e^{j\omega}))} \geq \sqrt{\text{cov}(\hat{G}_2(\omega))} \quad \forall \omega$$



A cosine at $\omega = 0.15$ seems a much better signal than a white noise to identify this true system

How can we determine the optimal input signal for a given true system?

→ Optimal experiment design

2. Optimal experiment design

Suppose that, for the subsequent use of the model (e.g. for control), it is required that:

$$|G(e^{j\omega}, \hat{\theta}_N) - G(e^{j\omega}, \theta_0)| < r_{adm}(\omega) \quad \forall \omega$$

where $r_{adm}(\omega)$ is a given frequency function

The function $r_{adm}(\omega)$ can e.g. be chosen equal to $0.1|G(e^{j\omega}, \hat{\theta}_N)|$ (relative error of 10%) or can be determined using robust control considerations

$$|G(e^{j\omega}, \hat{\theta}_N) - G(e^{j\omega}, \theta_0)| < r_{adm}(\omega) \quad \forall \omega$$

How can we choose $u(t)$ and N for the identification experiment in order to guarantee that the model identified with this experiment meets the accuracy constraint?

We have thus to determine N and $u(t)$ such that:

$$\alpha \sqrt{\Lambda_G^*(e^{j\omega}) \frac{P_\theta}{N} \Lambda_G(e^{j\omega})} < r_{adm}(\omega) \quad \forall \omega$$

First observation: By choosing N or the power of $u(t)$ sufficiently large, it is always possible to achieve this constraint for any given $r_{adm}(\omega)$

\implies multiple solutions !

Let us therefore

- fix the value of N

- and determine, for an experiment of that length N , the

excitation signal $u(t)$ with the least power $\mathcal{P}_u = \frac{1}{N} \sum_{t=1}^N u^2(t)$

that nevertheless leads to an identified model with the required accuracy

Optimization problem for the design of the excitation signal (N given)

Determine the sequence $u(t)$ ($t = 1 \dots N$) which solves:

$$\min_{u(t) \ (t=1 \dots N)} \quad \frac{1}{N} \sum_{t=1}^N u^2(t)$$

subject to $\alpha \sqrt{\Lambda_G^*(e^{j\omega}) \frac{P_\theta}{N} \Lambda_G(e^{j\omega})} < r_{adm}(\omega) \quad \forall \omega$

Issues with this optimization problem

non-linear relations between $u(t)$ and \mathcal{P}_u and between $u(t)$ and $P_\theta \rightarrow$ convexification

chicken-and-egg problem: dependence of $cov(G(e^{j\omega}, \hat{\theta}_N))$ on $\theta_0 \rightarrow$ replace θ_0 by an initial estimate θ_{init} obtained e.g. via an initial identification with white noise

infinite number of constraints i.e. at each frequency $\omega \rightarrow$ frequency grid

3. Convexification of the optimization problem

Approach: use of the power spectrum $\Phi_u(\omega)$ of $u(t)$ as alternative decision variable (instead of $u(t)$ itself)

3.1 Convexification of the cost function \mathcal{P}_u

for large N ,

$$\mathcal{P}_u = \frac{1}{N} \sum_{t=1}^N u^2(t) \approx \bar{E}u^2(t) = \frac{1}{2\pi} \int_{-\pi}^{\pi} \Phi_u(\omega) d\omega$$

\mathcal{P}_u linear in the new decision variable $\Phi_u(\omega)$

3.2 Convexification of the constraint $\alpha \sqrt{\Lambda_G^* \frac{P_\theta}{N} \Lambda_G} < r_{adm}(\omega)$

$$P_\theta = \sigma_e^2 \left(\bar{E} \psi(t, \theta_0) \psi^T(t, \theta_0) \right)^{-1} \text{ and } \psi(t) = F_u(z)u(t) + F_e(z)e(t)$$

$$\implies P_\theta = \sigma_e^2 \left(\frac{1}{2\pi} \int_{-\pi}^{\pi} F_u F_u^* \Phi_u(\omega) + F_e F_e^* \sigma_e^2 d\omega \right)^{-1}$$

P_θ^{-1} linear in the new decision variable $\Phi_u(\omega)$

Let us thus rewrite the constraint as a linear function of P_θ^{-1}

$$\alpha \sqrt{\Lambda_G^*(e^{j\omega}) \frac{P_\theta}{N} \Lambda_G(e^{j\omega})} < r_{adm}(\omega) \quad \forall \omega$$

is equivalent (via the Schur complement) to

$$N P_\theta^{-1} > R_{adm}(\omega) \quad \forall \omega$$

where $R_{adm}(\omega) = \frac{\alpha^2}{r_{adm}^2(\omega)} \Lambda_G(e^{j\omega}) \Lambda_G^*(e^{j\omega})$

3.3. A convex formulation of experiment design

Reminder: initial optimization problem (non-convex)

Determine the sequence $u(t)$ ($t = 1 \dots N$) which solves:

$$\min_{u(t) \ (t=1 \dots N)} \frac{1}{N} \sum_{t=1}^N u^2(t)$$

$$\text{subject to } \alpha \sqrt{\Lambda_G^*(e^{j\omega}) \frac{P_\theta}{N} \Lambda_G(e^{j\omega})} < r_{adm}(\omega) \quad \forall \omega$$

Convex alternative (fixed N):

Determine the power spectrum $\Phi_u(\omega)$ of the excitation signal $u(t)$ which solves:

$$\min_{\Phi_u(\omega)} \quad \frac{1}{2\pi} \int_{-\pi}^{\pi} \Phi_u(\omega) d\omega$$

subject to $\Phi_u(\omega) \geq 0 \quad \forall \omega$ and to

$$\frac{N}{\sigma_e^2} \left(\frac{1}{2\pi} \int_{-\pi}^{\pi} F_u F_u^* \Phi_u(\omega) + F_e F_e^* \sigma_e^2 d\omega \right) > R_{adm}(\omega) \quad \forall \omega$$

3.4. Parametrization of the power spectrum $\Phi_u(\omega)$

The decision variable $\Phi_u(\omega)$ has an infinite dimension \rightarrow a simpler and linear parametrization of $\Phi_u(\omega)$ is required

Two possible approaches:

- $\Phi_u(\omega)$ for filtered white noise u
- $\Phi_u(\omega)$ for multisine u

Parametrization of the power spectrum $\Phi_u(\omega)$: filtered white noise

Suppose we want to be able to generate our input signal $u(t)$ of spectrum $\Phi_u(\omega)$ as the realization of a white noise filtered by a FIR filter $F_m(z)$ of order m :

$$u(t) = F_m(z)w(t)$$

where $w(t)$ is a white noise of variance 1 and

$F_m(z) = f_0 + f_1z^{-1} + \dots + f_mz^{-m}$ an arbitrary FIR filter

In this case, we have to restrict the structure of $\Phi_u(\omega)$ as follows:

$$\Phi_u(\omega) = |F_m(e^{j\omega})|^2 = \sum_{r=-m}^m c_r e^{j\omega r} \quad (\text{met } c_r = c_{-r})$$

where c_r ($r = 0..m$) are the new decision variables of the optimization problem

The larger m is chosen, the more flexible is the parametrization of the spectrum

Choosing $m = 0$ is equivalent to restrict attention to a flat spectrum (white noise)

Parametrization of the power spectrum $\Phi_u(\omega)$: multisine

Suppose we want to generate our input signal $u(t)$ as:

$$u(t) = \sum_{r=1}^m A_r \sin(\omega_r t)$$

with arbitrary amplitudes A_r ($r = 1 \dots m$), but fixed frequencies ω_r ($r = 1 \dots m$) e.g. linearly distributed in the interval $[0 \ \pi]$

$$u(t) = \sum_{r=1}^m A_r \sin(\omega_r t)$$

In this case, we have to restrict the structure of $\Phi_u(\omega)$ as follows:

$$\Phi_u(\omega) = \frac{\pi}{2} \sum_{r=1}^m A_r^2 (\delta(\omega - \omega_r) + \delta(\omega + \omega_r))$$

where A_r^2 ($r = 1 \dots m$) are the new decision variables of the optimization problem

Here also, the larger m is chosen, the more "flexible" is the parametrization of the spectrum

A bit more details....

$$\Phi_u(\omega) = \frac{\pi}{2} \sum_{r=1}^m A_r^2 (\delta(\omega - \omega_r) + \delta(\omega + \omega_r)) \implies$$

$$\mathcal{P}_u = \frac{1}{2\pi} \int_{-\pi}^{\pi} \Phi_u(\omega) d\omega = \sum_{r=1}^m \frac{A_r^2}{2}$$

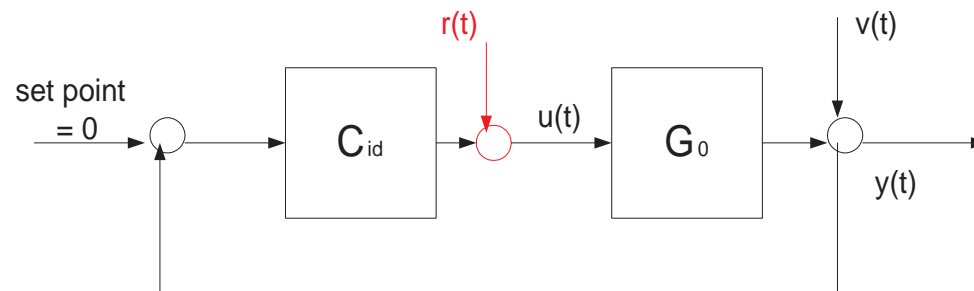
$$N P_\theta^{-1} = \frac{N}{\sigma_e^2} \left(\left(\sum_{r=1}^m \frac{A_r^2}{2} M(\omega_r) \right) + R_e \right)$$

$$M(\omega_r) = \text{Re}(F_u(e^{j\omega_r}) F_u^*(e^{j\omega_r})) \quad R_e = \frac{1}{2\pi} \int_{-\pi}^{\pi} F_e F_e^* \sigma_e^2 d\omega$$

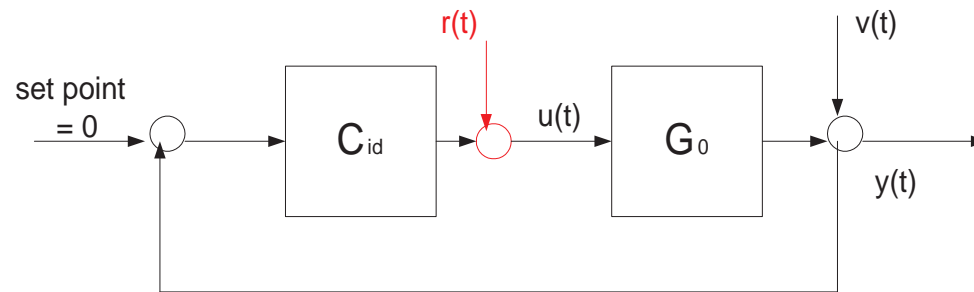
\mathcal{P}_u and P_θ^{-1} are indeed affine in A_r^2 ($r = 1 \dots m$)

4. Some alternative experiment design problems

- 1) Instead of \mathcal{P}_u , the cost function can also be the power \mathcal{P}_y of the output signal $y(t)$ (or a combination of both)
- 2) Optimal experiment design can also be formulated for **direct closed-loop identification**



$$\mathbf{Z}^N = \{y(t) \ u(t) | t = 1 \dots N\} \text{ generated via } r(t)$$



Decision variable: $\Phi_r(\omega)$

Expression for P_θ

$$P_\theta = \sigma_e^2 \left(\frac{1}{2\pi} \int_{-\pi}^{\pi} F_r F_r^* \Phi_r(\omega) + F_v F_v^* \sigma_e^2 d\omega \right)^{-1}$$

$$F_r(z, \theta_0) = \frac{S_{id}}{H_0} \left. \frac{\partial G(z, \theta)}{\partial \theta} \right|_{\theta=\theta_0} \quad F_v(z, \theta_0) = \frac{\left. \frac{\partial H(z, \theta)}{\partial \theta} \right|_{\theta=\theta_0}}{H_0} - C_{id} S_{id} \left. \frac{\partial G(z, \theta)}{\partial \theta} \right|_{\theta=\theta_0}$$

P_θ^{-1} linear in the decision variable $\Phi_r(\omega)$

Possible formulation for direct closed-loop identification

$$\min_{\Phi_r(\omega)} \quad \frac{1}{2\pi} \int_{-\pi}^{\pi} \Phi_r(\omega) d\omega$$

subject to $\Phi_r(\omega) \geq 0 \quad \forall \omega$ and to

$$\frac{N}{\sigma_e^2} \left(\frac{1}{2\pi} \int_{-\pi}^{\pi} F_r F_r^* \Phi_r(\omega) + F_v F_v^* \sigma_e^2 d\omega \right) > R_{adm}(\omega) \quad \forall \omega$$

An other formulation will be presented in the sequel

3) Other (dual) paradigms for (open-loop) input design:

$$\min_{\Phi_u(\omega)} \det(P_\theta)$$

subject to $\Phi_u(\omega) \geq 0 \quad \forall \omega$ and to

$$\frac{1}{2\pi} \int_{-\pi}^{\pi} \Phi_u(\omega) d\omega < \mathcal{P}_{max}$$

with \mathcal{P}_{max} the maximal power that is allowed at the input

5. Illustration 1 (see example slides 8-12)

$$\mathcal{S} : y(t) = \frac{3.6z^{-1}}{1 - 0.7z^{-1}}u(t) + (1 + 0.9z^{-1})e(t)$$

$$\mathcal{M} : G(z, \theta) = \frac{bz^{-1}}{1 - fz^{-1}} \quad H(z, \theta) = 1 + cz^{-1}$$

We would like to find, for an experiment of duration $N = 500$, the least powerful excitation signal $u(t)$ which leads to an identified model $G(z, \hat{\theta}_N)$ with a relative modeling error of less than 1 % at each ω

$$|G(e^{j\omega}, \hat{\theta}_N) - G(e^{j\omega}, \theta_0)| < \underbrace{0.01 |G(e^{j\omega}, \hat{\theta}_N)|}_{=r_{adm}(\omega)} \quad \forall \omega$$

Design of the optimal input signal

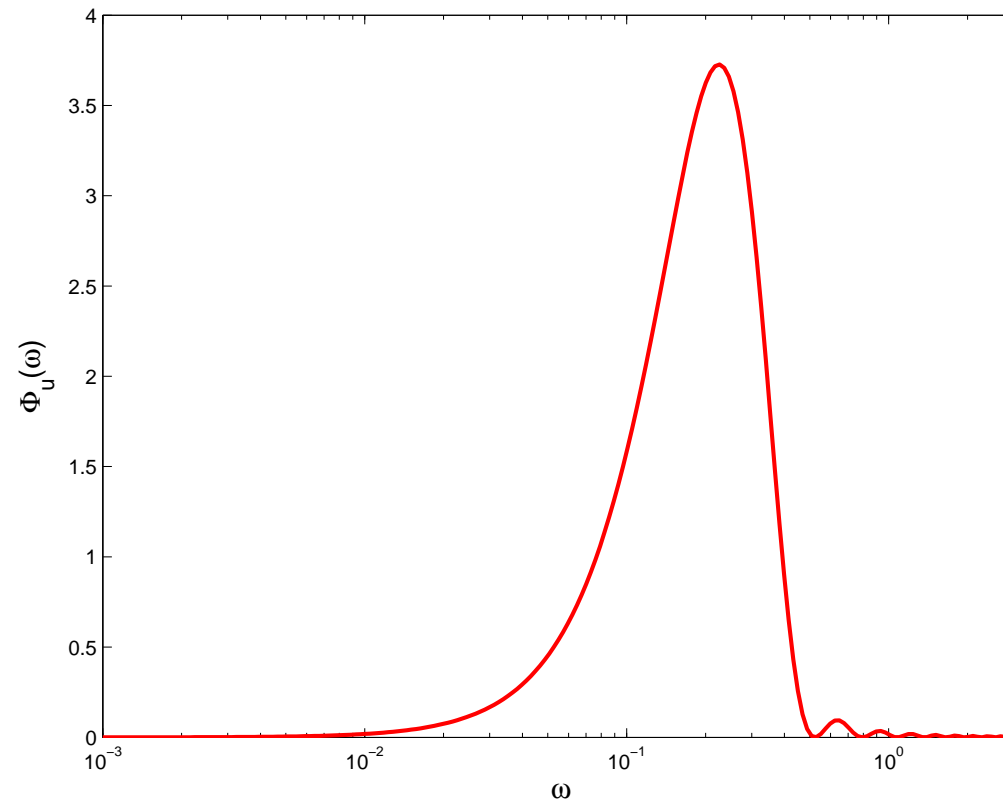
The required initial estimate θ_{init} is obtained via an initial experiment of duration $N = 100$ with white input ($\sigma_u^2 = 1$)

We solve then the convex optimization problem where θ_0 and $\hat{\theta}_N$ have been replaced by θ_{init} e.g.

$$r_{adm}(\omega) = 0.01 |G(e^{j\omega}, \theta_{init})| \quad \forall \omega$$

Other choices: $\alpha = 2.45$ $N = 500$

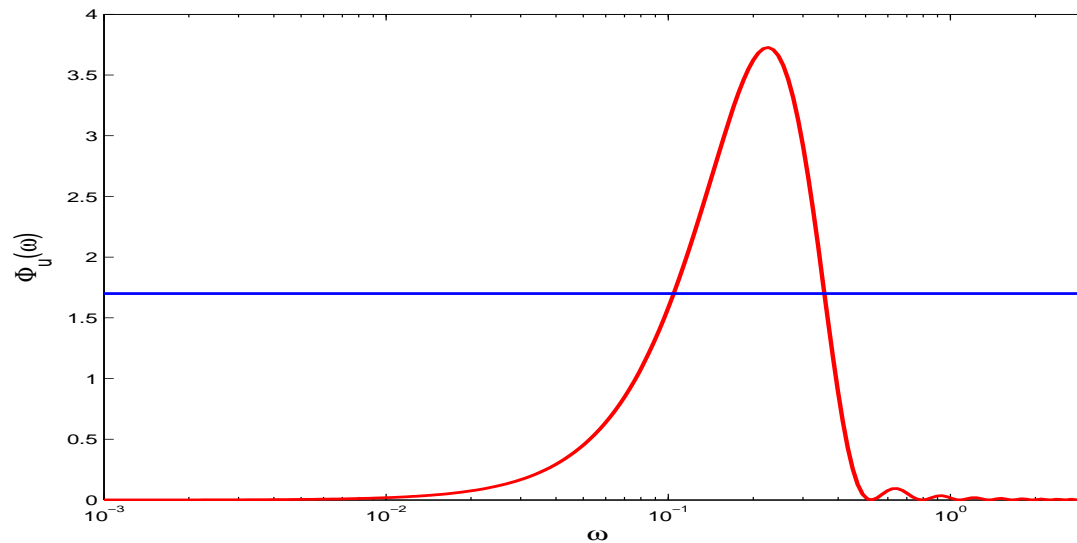
The filtered white noise parametrization of $\Phi_u(\omega)$ with $m = 20$ leads to the following optimal spectrum $\Phi_u(\omega)$:



This spectrum corresponds to a power $\mathcal{P}_{u,opt} = 0.25$

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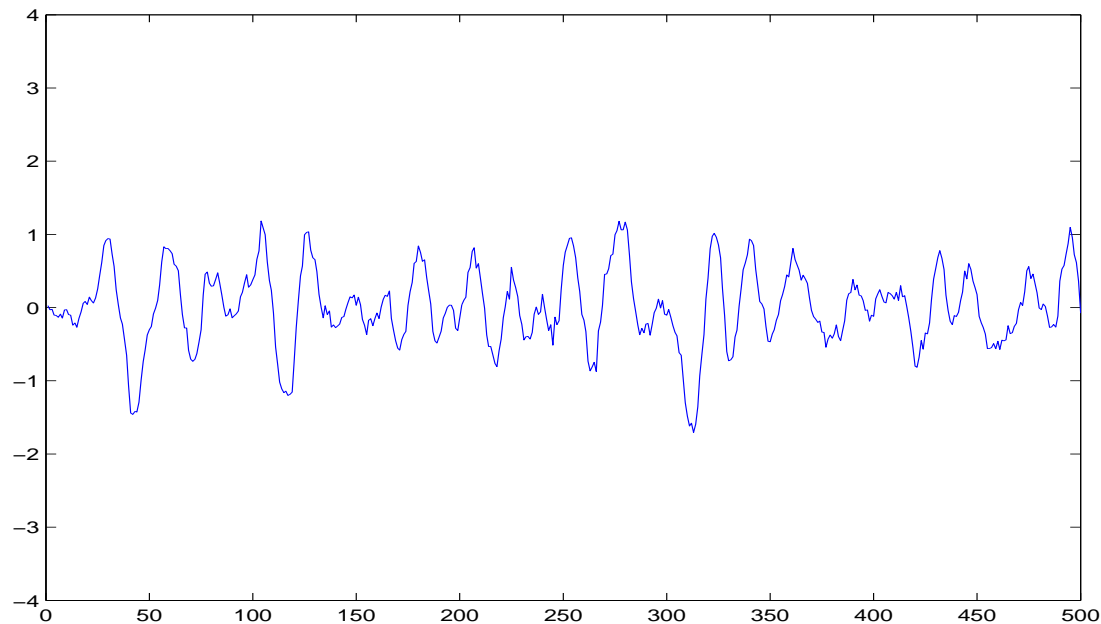
In comparison, if you want to achieve the same accuracy with a white noise ($m = 0$), you need a white noise of power 1.7



By shaping $\Phi_u(\omega)$, less power is needed to achieve the required accuracy !

Verification

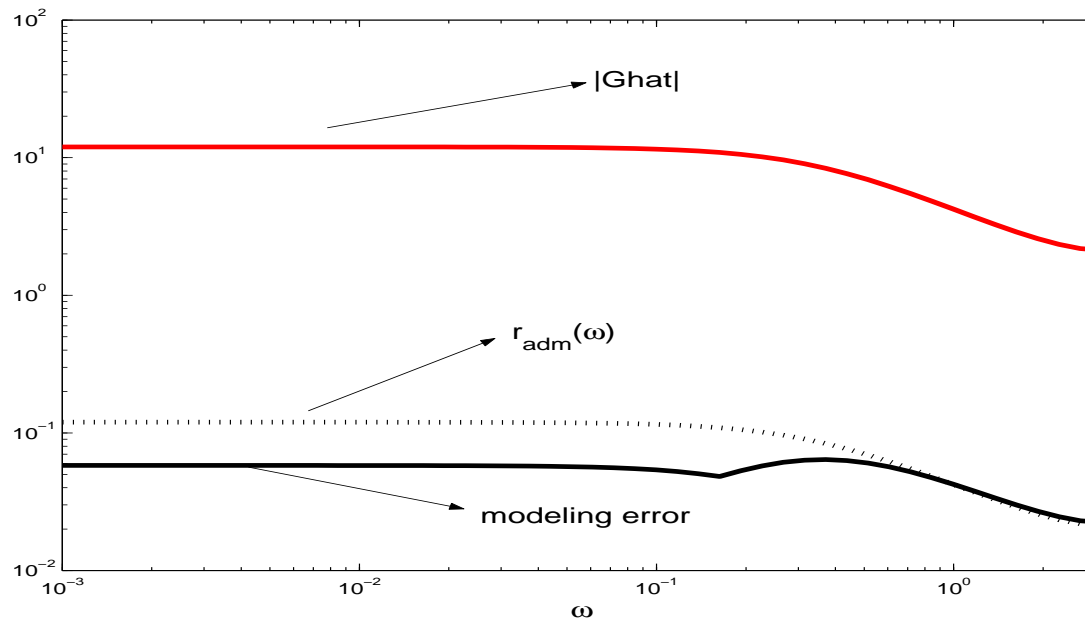
We have generated an input signal of length $N = 500$ having the optimal spectrum:



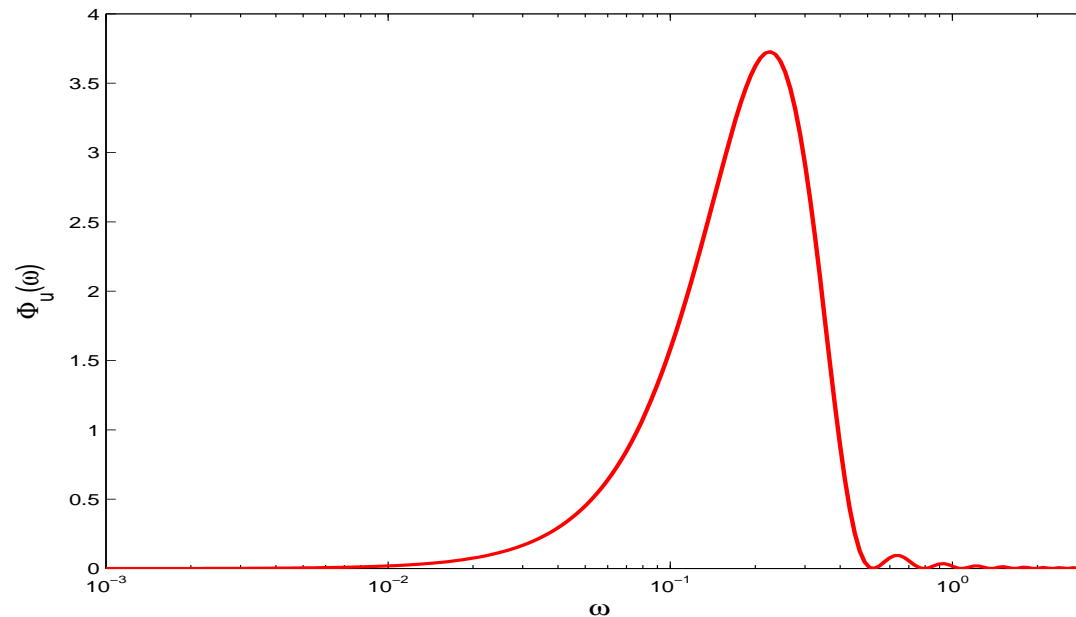
This signal is used to identify a model $G(z, \hat{\theta}_N)$ of G_0

The modeling error is computed/bounded using P_θ and we observe, as expected,

$$\alpha \sqrt{\Lambda_G^* \frac{P_\theta}{N} \Lambda_G} < r_{adm}(\omega) \quad \forall \omega$$



Analysis of the optimal spectrum $\Phi_u(\omega)$



The optimal power spectrum concentrates the power around the frequency $\omega = 0.15 \rightarrow$ a (co)sine at this frequency will also be appropriate

Let us thus refine the optimal spectrum with the multisine parametrization of $\Phi_u(\omega)$

This yields the optimal signal:

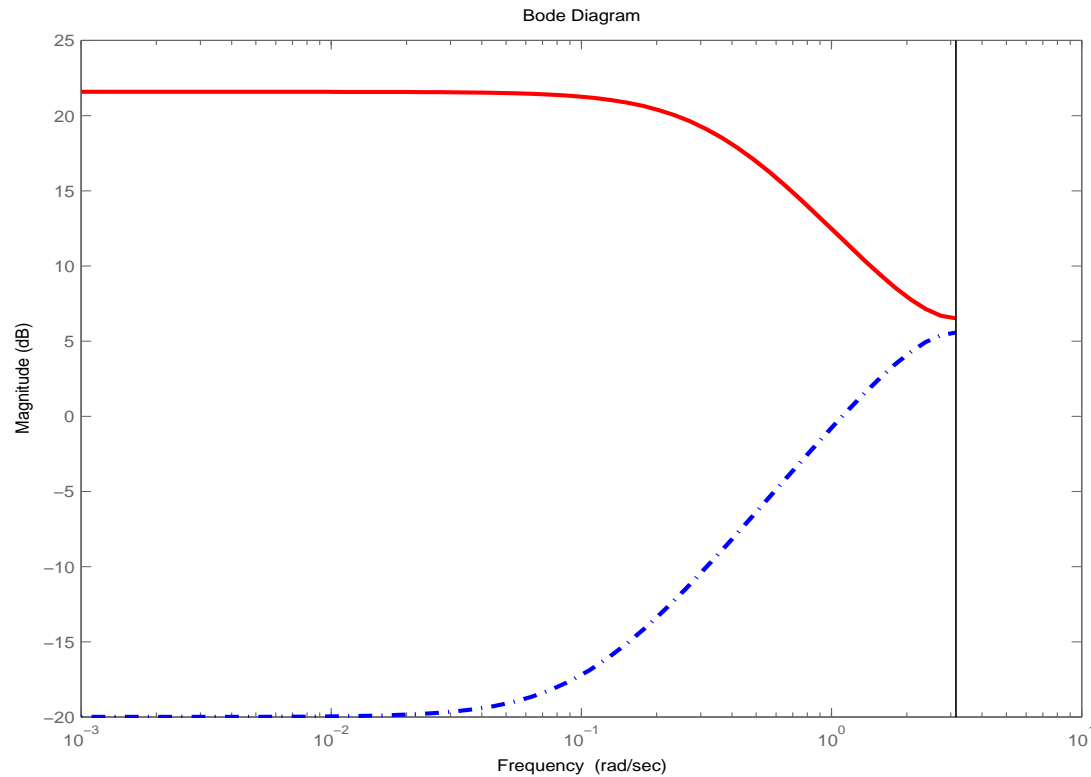
$$u(t) = 0.57 \sin(0.15 t)$$

which corresponds to an even **smaller** $\mathcal{P}_{u,opt} = \frac{(0.57)^2}{2} = 0.16$

Remark. If we choose a sine at $\omega = 0.01$, we would need an amplitude $A = 3$ corresponding to a power $\mathcal{P}_u = 4.5$

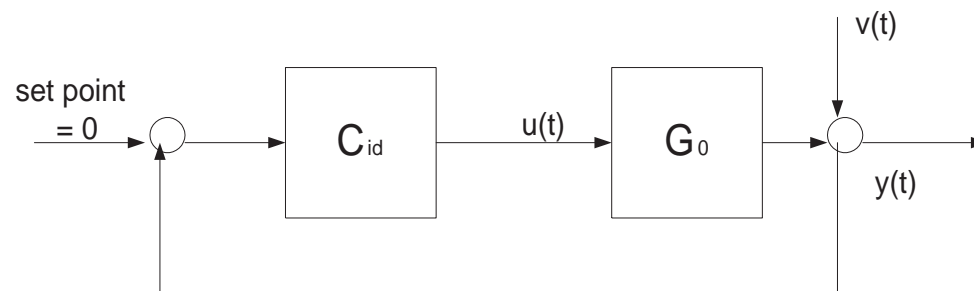
Why concentrating the power around $\omega = 0.15$?

This is the frequency range with the pole of G_0 (red) and the zero of H_0 (blue) !!!!



6. Illustration 2: least intrusive identification experiment for control

We consider a real-life system G_0 (with two resonances) operated in closed loop whose objective is to reject the disturbance $v(t) = H_0(z)e(t)$:



This closed-loop delivers a product $y(t)$

Objective: (re)-identify a model of G_0 (and H_0) to update C_{id} by a new robust controller designed with the model

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$$\rightarrow \text{a constraint } \alpha \sqrt{\Lambda_G^* \frac{P_\theta}{N} \Lambda_G} < r_{adm}(\omega) \quad \forall \omega$$

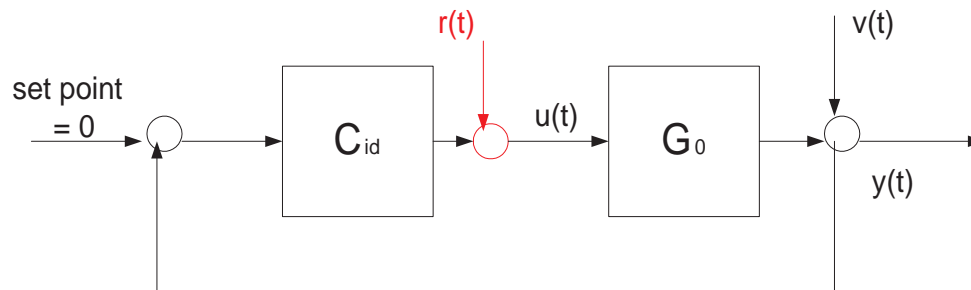
Extra requirement: in order to be able to continue to produce a good $y(t)$, we would like to achieve this objective with the least intrusive identification experiment

How can we evaluate the effects of an identification experiment on the production quality??

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Normal operation vs. Identification experiment operation

$$\left\{ \begin{array}{l} y(t) = S_{id}v(t) \\ u(t) = -C_{id}S_{id}v(t) \end{array} \right. \quad vs. \quad \left\{ \begin{array}{l} y(t) = \overbrace{G_0 S_{id} r(t)}^{y_r(t)} + S_{id}v(t) \\ u(t) = \overbrace{S_{id} r(t)}^{u_r(t)} - C_{id}S_{id}v(t) \end{array} \right.$$



The perturbations y_r and u_r are present during the experiment duration N

Their presence reduce the production quality

For fixed N , we can measure the **performance degradation** due to the application of a signal with power spectrum $\Phi_r(\omega)$ by:

$$\begin{aligned}\mathcal{J}_r &= \alpha_y \mathcal{P}_{y_r} + \alpha_u \mathcal{P}_{u_r} \\ &= \alpha_y \left(\frac{1}{2\pi} \int_{-\pi}^{\pi} \Phi_{y_r}(\omega) d\omega \right) + \alpha_u \left(\frac{1}{2\pi} \int_{-\pi}^{\pi} \Phi_{u_r}(\omega) d\omega \right)\end{aligned}$$

e.g. $\alpha_u = \alpha_y = 1$

⇒ **Least intrusive identification experiment for control**

$$\arg \min_{\Phi_r(\omega)} \overbrace{\mathcal{P}_{u_r} + \mathcal{P}_{y_r}}^{\mathcal{J}_r}$$

under the constraint that $N P_\theta^{-1} \geq R_{adm}(\omega) \quad \forall \omega$

Reminder: $N P_\theta^{-1} > R_{adm}(\omega)$ is equivalent to:

$$\alpha \sqrt{\Lambda_G^* \frac{P_\theta}{N} \Lambda_G} < r_{adm}(\omega)$$

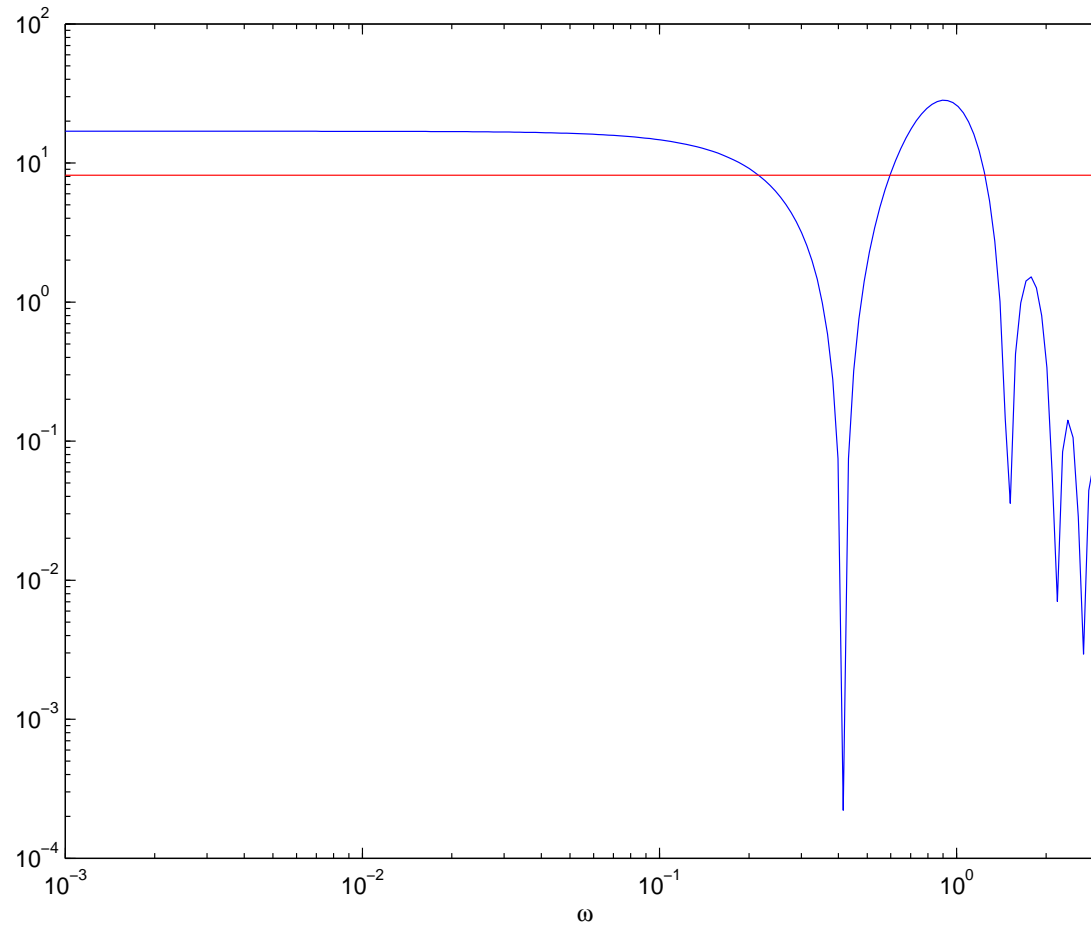
Here, $r_{adm}(\omega)$ has been chosen based on robust control considerations and the duration of the experiment is fixed to $N = 500$.

$$\mathit{arg} \min_{\Phi_r(\omega)} \mathcal{P}_{u_r} + \mathcal{P}_{y_r} \quad \text{subject to } N P_\theta^{-1} \geq R_{adm}(\omega) \quad \forall \omega$$

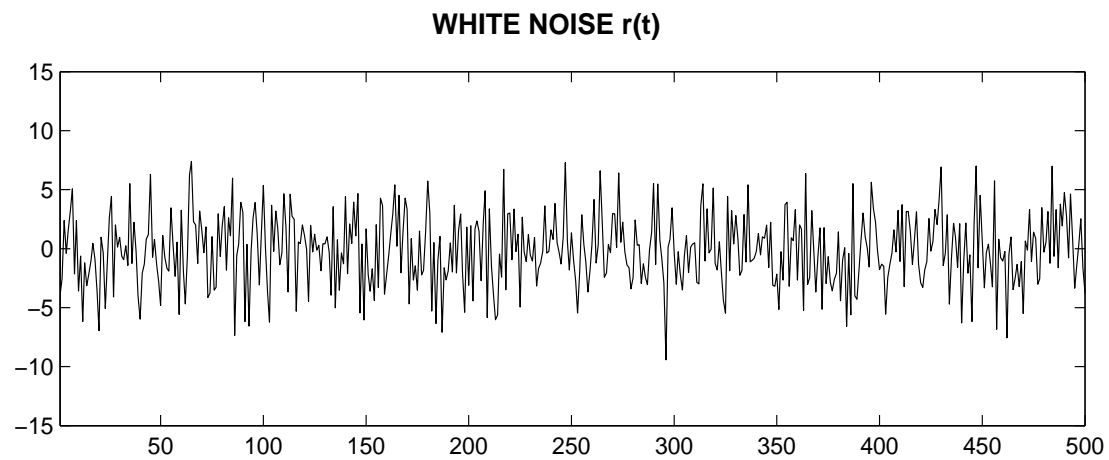
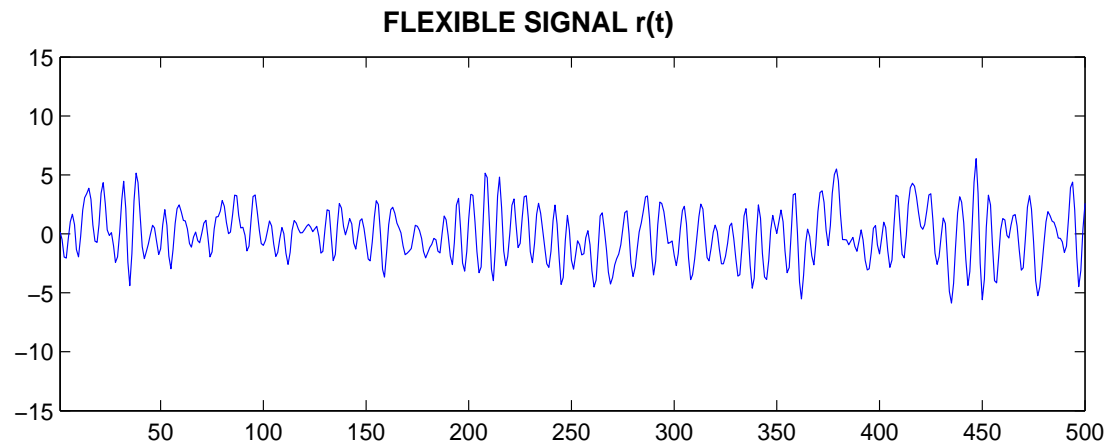
We consider two cases for the optimization on $\Phi_r(\omega)$ (filtered white noise parametrization):

- **the first one constrains $r(t)$ to be a white noise ($m = 0$)**
- **the second one allows $r(t)$ to have more flexible $\Phi_r(\omega)$ ($m = 10$)**

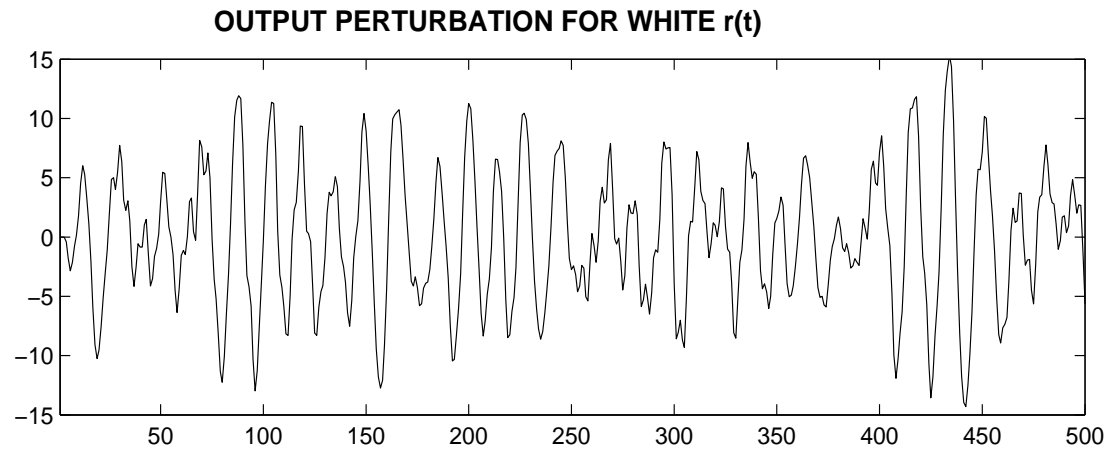
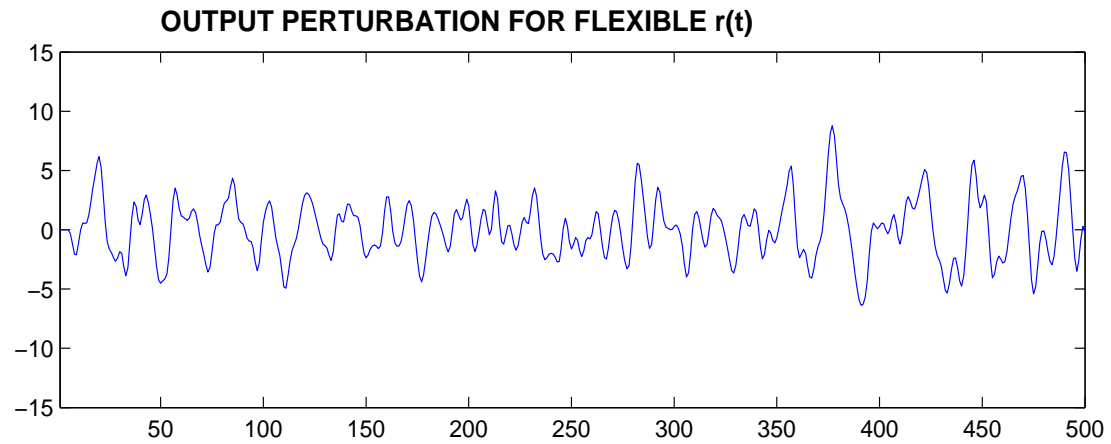
The two optimizations deliver the following spectra: white noise (red) and flexible spectrum (blue):



Let us generate a realization of length $N = 500$ of those two spectra...



This leads to the following perturbations $y_r(t)$



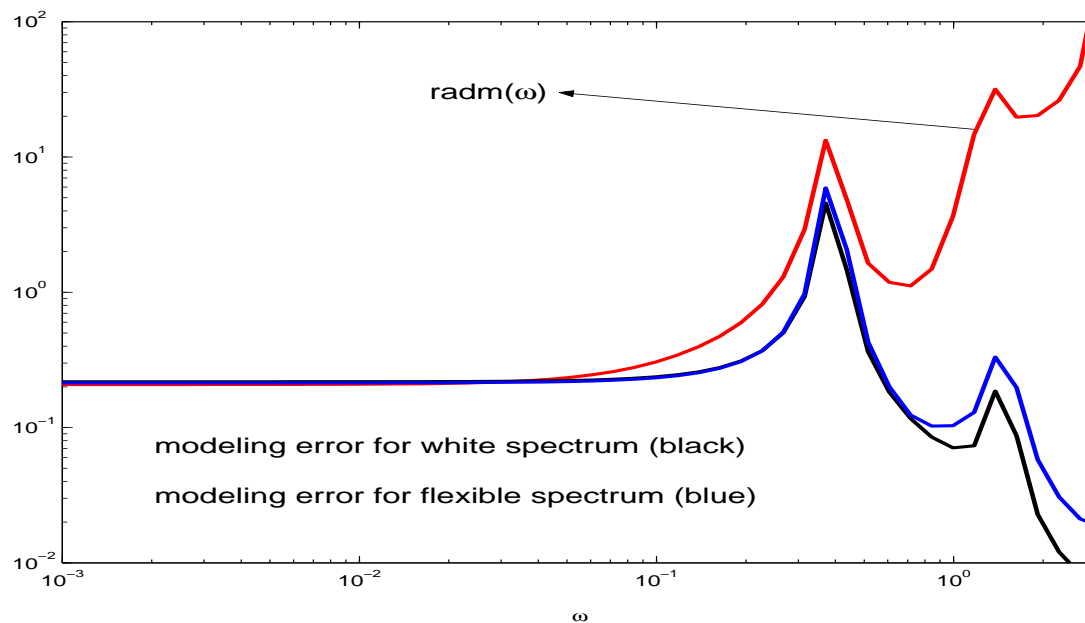
$\Phi_r(\omega)$	required \mathcal{J}_r for $N = 500$
white	22.5
flexible	9.9

Conclusions:

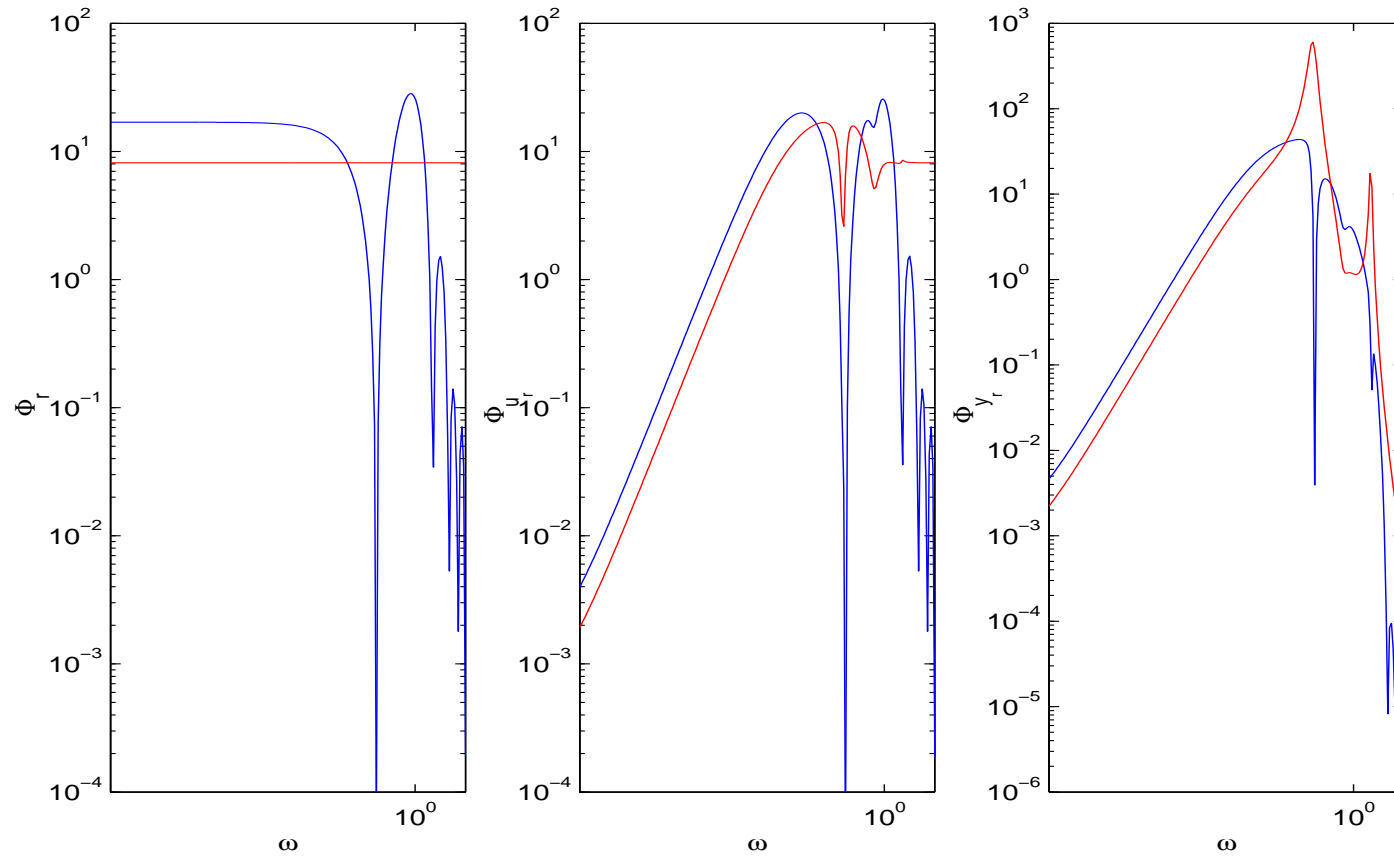
- both experimental conditions deliver a model accurate enough for robust control (see next slide)
- $\mathcal{J}_r = \mathcal{P}_{u_r} + \mathcal{P}_{y_r}$ is two times larger with the RBS than with the frequency-dependent signal (i.e. 22.5 vs. 9.9)

That both the flexible and the white spectra deliver a model accurate enough for robust control can be seen by verifying in both cases that

$$\underbrace{\alpha \sqrt{\Lambda_G^* \frac{P_\theta}{N} \Lambda_G}}_{\text{modeling error}} < r_{adm}(\omega) \quad \forall \omega$$



Explanation of the better \mathcal{J}_r for the flexible signal:



$\Phi_r(\omega)$, $\Phi_{u_r}(\omega)$ and $\Phi_{y_r}(\omega)$ for the white $r(t)$ (red) and for the flexible $r(t)$ (blue)

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