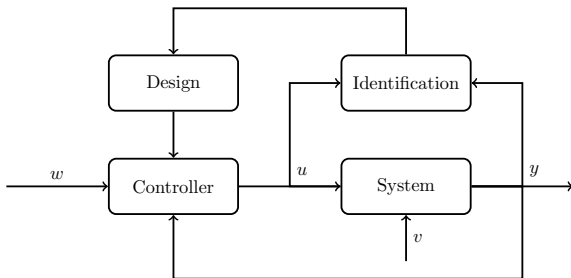


- 1 System theory
- 2 Stochastics
- 3 State estimation 1
- 4 State estimation 2
- 5 Optimal control 1
- 6 System identification 1 + adaptive control 1
- 7 External models + prediction
- 8 Optimal control 2
- 9 Optimal control 3
- 10 System identification 2
- 11 System identification 3 + model validation**
- 12 System identification 4 + adaptive control 2
- 13 Adaptive control 3



- Time-varying estimation
- Design of experiments
- Model validation

Stochastic Adaptive Control - Follow-up

Follow-up from last time



Questions?

Heuristics for time-varying systems

ARX model

$$A(q^{-1})y_t = B(t, q^{-1})u_t + e_t, \quad (1)$$

$$b_1(t) = b_{1,0} + b_{1,1}t \quad (2)$$

Treat time-varying coefficient as two coefficients with their own inputs

$$y_t = \phi^T \theta + e_t \quad (3)$$

$$\theta^T = [a_1 \quad a_2 \quad \cdots \quad a_{n_a} \quad b_{1,0} \quad b_{1,1} \quad b_2 \quad \cdots \quad b_{n_b}] \quad (4)$$

$$\phi^T = [-y_{t-1} \quad -y_{t-2} \quad \cdots \quad -y_{t-n_a} \quad u_{t-1} \quad tu_{t-1} \quad u_{t-2} \quad \cdots \quad u_{t-n_b}] \quad (5)$$

For deterministic time varying systems, rearrange the parameters

$$y_t = \phi_t^T \theta_t + e_t \quad (6)$$

$$\theta_t = \alpha + f(t)\beta = \begin{bmatrix} I & f(t) \end{bmatrix} \begin{bmatrix} \alpha \\ \beta \end{bmatrix} \quad (7)$$

$$y_t = \begin{bmatrix} \phi_t^T & \phi_t^T f(t) \end{bmatrix} \begin{bmatrix} \alpha \\ \beta \end{bmatrix} + e_t \quad (8)$$

For piece-wise linear parameters, rearrange the parameters

$$y_t = \phi_t^T \theta_t + e_t \quad (9)$$

$$\theta_t = \alpha_i + (t - T_i)\beta_i, \quad T_i \leq t \leq T_{i+1} \quad (10)$$

$$y_t = \begin{bmatrix} \phi_t^T & \phi_t^T (t - T_i) \end{bmatrix} \begin{bmatrix} \alpha_i \\ \beta_i \end{bmatrix} + e_t \quad (11)$$

System with general time-varying parameters

$$\theta_{t+1} = f(t, \theta_t, v_t) \quad (12)$$

- The methods discussed so far cannot estimate the time-varying dynamics and were not designed to do it
- In practice, the problem is that the correction factor diminishes over time

$$P_t \rightarrow 0 \quad (13)$$

Reset the covariance after some time, t_i

$$P_{t_i} = P_i > P_{t_i-1}, \quad \hat{\theta}_{t_i} = \hat{\theta}_{t_i-1} \quad (14)$$

The appropriate restarting time depends on the application

For instance, restart at fixed intervals

$$t_i = Ni \quad (15)$$

This can be useful for periodic systems

Another method: Keep the correction term large

For instance, keep the correction term κ constant

$$\hat{\theta}_t = \hat{\theta}_{t-1} + \kappa \epsilon_t \quad (16)$$

$$\tilde{\theta}_t = (I - \kappa \phi_t^T) \tilde{\theta}_{t-1} - \kappa \epsilon_t \quad (17)$$

Alternatively, keep the variance constant

$$P_t = P \quad (18)$$

$$\hat{\theta}_t = \hat{\theta}_{t-1} + \kappa \epsilon_t \quad (19)$$

$$\kappa_t = \frac{P \phi_t}{1 + \phi_t^T P \phi_t} \quad (20)$$

Time-varying systems - Forgetting methods: Exponential Forgetfulness

Another approach: Forget a little bit all the time (exponential forgetfulness)

$$J_t = \frac{1}{2} \sum_{i=1}^t \lambda^{t-i} \epsilon_i^2 = \lambda J_{t-1} + \frac{1}{2} \epsilon_t^2 \quad (21)$$

The recursion is similar to the previous methods

$$\hat{\theta}_t = \hat{\theta}_{t-1} + P_t \phi_t \epsilon_t \quad (22)$$

$$\epsilon_t = y_t - \phi_t^T \hat{\theta}_{t-1} \quad (23)$$

$$P_t^{-1} = \lambda P_{t-1}^{-1} + \phi_t \phi_t^T \quad (24)$$

The forgetting factor λ can be expressed in terms of a horizon, N_∞

$$\lambda = 1 - \frac{1}{N_\infty} \quad (25)$$

Model

$$y_t = \frac{-1}{4}y_{t-1} + \frac{1}{2}y_{t-2} + u_{t-1} + e_t \quad (26)$$

New measurement ($t = 10$)

$$y_t = 1.47, \quad u_{t-1} = 2 \quad (27)$$

Covariates, parameter estimate and covariance

$$\phi_t = \begin{bmatrix} 2.42 \\ 2.57 \\ 2 \end{bmatrix}, \quad \hat{\theta}_{t-1} = \begin{bmatrix} -0.2505 \\ 0.4960 \\ 0.9991 \end{bmatrix}, \quad P_{t-1} = \begin{bmatrix} 0.1355 & -0.0431 & -0.1057 \\ -0.0431 & 0.0620 & -0.0088 \\ -0.1057 & -0.0088 & 0.1242 \end{bmatrix}$$

Forgetting factor

$$\lambda = 0.95 \quad (28)$$

Residual

$$\epsilon_t = y_t - \phi_t^T \hat{\theta}_{t-1} = 1.47 - \begin{bmatrix} 2.42 & 2.57 & 2 \end{bmatrix} \begin{bmatrix} -0.2505 \\ 0.4960 \\ 0.9991 \end{bmatrix} = -1.1967 \quad (29)$$

Covariance

$$P_t^{-1} = \lambda P_{t-1}^{-1} + \phi_t \phi_t^T = 0.95 \begin{bmatrix} 259.9763 & 214.2840 & 236.4348 \\ 214.2840 & 192.9152 & 196.0344 \\ 236.4348 & 196.0344 & 223.1583 \end{bmatrix} \quad (30)$$

$$+ \begin{bmatrix} 2.42 \\ 2.57 \\ 2 \end{bmatrix} \begin{bmatrix} 2.42 & 2.57 & 2 \end{bmatrix} = \begin{bmatrix} 252.8339 & 209.7892 & 229.4530 \\ 209.7892 & 189.8743 & 191.3727 \\ 229.4530 & 191.3727 & 216.0004 \end{bmatrix} \quad (31)$$

Parameter estimate

$$\hat{\theta}_t = \hat{\theta}_{t-1} + P_t \phi_t \epsilon_t = \begin{bmatrix} -0.2505 \\ 0.4960 \\ 0.9991 \end{bmatrix} \quad (32)$$

$$- 1.1967 \begin{bmatrix} 0.1426 & -0.0456 & -0.1111 \\ -0.0456 & 0.0638 & -0.0081 \\ -0.1111 & -0.0081 & 0.1298 \end{bmatrix} \begin{bmatrix} 2.42 \\ 2.57 \\ 2 \end{bmatrix} = \begin{bmatrix} -0.2574 \\ 0.4512 \\ 1.0350 \end{bmatrix} \quad (33)$$

Model

$$y_t = \frac{-1}{4}y_{t-1} + \frac{1}{2}y_{t-2} + u_{t-1} + e_t \quad (34)$$

New measurement ($t = 10$)

$$y_t = 1.47, \quad u_{t-1} = 2 \quad (35)$$

Covariates, parameter estimate and covariance

$$\phi_t = \begin{bmatrix} 2.42 \\ 2.57 \\ 2 \end{bmatrix}, \quad \hat{\theta}_{t-1} = \begin{bmatrix} -0.2505 \\ 0.4960 \\ 0.9991 \end{bmatrix}, \quad P_{t-1} = \begin{bmatrix} 0.1355 & -0.0431 & -0.1057 \\ -0.0431 & 0.0620 & -0.0088 \\ -0.1057 & -0.0088 & 0.1242 \end{bmatrix}$$

Forgetting factor

$$\lambda = 0.90 \quad (36)$$

Solve the exercise in 15 min.

Residual

$$\epsilon_t = y_t - \phi_t^T \hat{\theta}_{t-1} = 1.47 - \begin{bmatrix} 2.42 & 2.57 & 2 \end{bmatrix} \begin{bmatrix} -0.2505 \\ 0.4960 \\ 0.9991 \end{bmatrix} = -1.1967 \quad (37)$$

Covariance

$$P_t^{-1} = \lambda P_{t-1}^{-1} + \phi_t \phi_t^T = 0.90 \begin{bmatrix} 259.9763 & 214.2840 & 236.4348 \\ 214.2840 & 192.9152 & 196.0344 \\ 236.4348 & 196.0344 & 223.1583 \end{bmatrix} \quad (38)$$

$$+ \begin{bmatrix} 2.42 \\ 2.57 \\ 2 \end{bmatrix} \begin{bmatrix} 2.42 & 2.57 & 2 \end{bmatrix} = \begin{bmatrix} 239.8351 & 199.0750 & 217.6313 \\ 199.0750 & 180.2286 & 181.5709 \\ 217.6313 & 181.5709 & 204.8424 \end{bmatrix} \quad (39)$$

Parameter estimate

$$\hat{\theta}_t = \hat{\theta}_{t-1} + P_t \phi_t \epsilon_t = \begin{bmatrix} -0.2505 \\ 0.4960 \\ 0.9991 \end{bmatrix} \quad (40)$$

$$- 1.1967 \begin{bmatrix} 0.1505 & -0.0481 & -0.1172 \\ -0.0481 & 0.0672 & -0.0085 \\ -0.1172 & -0.0085 & 0.1369 \end{bmatrix} \begin{bmatrix} 2.42 \\ 2.57 \\ 2 \end{bmatrix} = \begin{bmatrix} -0.2577 \\ 0.4488 \\ 1.0369 \end{bmatrix} \quad (41)$$

Time-varying systems - Fortescue's Method

Improve with a time-varying forgetting factor depending on the prediction error, ϵ_t

$$\lambda_t = 1 - \frac{1}{N_0} \frac{\epsilon_t^2}{\sigma^2 s_t} \quad (42)$$

N_0 is the approx. horizon over which the parameter is roughly constant

Recursion

$$\epsilon_t = y_t - \phi_t^T \hat{\theta}_{t-1} \quad (43)$$

$$s_t = 1 + \phi_t^T P_{t-1} \phi_t \quad (44)$$

$$K_t = \frac{P_{t-1} \phi_t}{\lambda_t + s_t} \quad (45)$$

$$\hat{\theta}_t = \hat{\theta}_{t-1} + K_t \epsilon_t \quad (46)$$

$$P_t = (I - K_t \phi_t^T) P_{t-1} \frac{1}{\lambda_t} \quad (47)$$

If the variance is unknown, we can introduce an estimate

$$\lambda_t = 1 - \frac{1}{N_0} \frac{\epsilon_t^2}{r_t s_t} \quad (48)$$

$$r_t = r_{t-1} + \frac{1}{t} \left(\frac{\epsilon_t^2}{s_t} - r_{t-1} \right), \quad r_0 = \epsilon_0^2 \quad (49)$$

Introduce model of parameters

$$\theta_{t+1} = \theta_t + v_t, \quad v_t \sim N(0, R_1 \sigma^2) \quad (50)$$

$$y_t = \phi_t^T \theta_t + e_t, \quad e_t \sim N(0, \sigma^2) \quad (51)$$

Estimate parameters using the Kalman filter

Data update

$$\hat{\theta}_{t|t} = \hat{\theta}_{t|t-1} + P_{t|t-1} \phi_t (y_t - \phi_t^T \hat{\theta}_{t|t-1}) \quad (52)$$

$$P_{t|t}^{-1} = P_{t|t-1}^{-1} + \phi_t \phi_t^T \quad (53)$$

Time update

$$\hat{\theta}_{t+1|t} = \hat{\theta}_{t|t} \quad (54)$$

$$P_{t+1|t} = P_{t|t} + R_1 \quad (55)$$

Experiment design

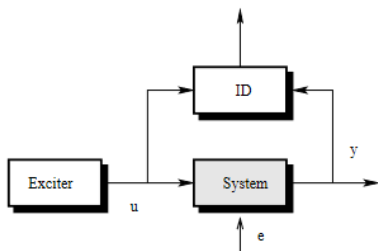
When attempting to identify a system, consider the following:

- ① What are the outputs?
- ② What are the inputs?
- ③ What are the disturbances?

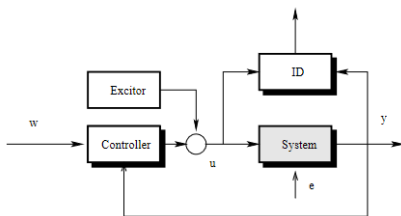
Also consider some practical aspects of the system

- ① What are we allowed to do?
- ② What type of model are we interested in?

Open Loop



Closed Loop



For any system \mathcal{S} , we can construct a set of models \mathcal{M} to describe it

$$\mathcal{S} : y = G_0(q)u + H_0(q)e \quad (56)$$

$$\mathcal{M} = \{G(q, \theta), H(q, \theta) | \theta \in \mathcal{D}\} \quad (57)$$

Ideally, the system should be included in the set of possible models

$$\mathcal{S} \in \mathcal{M} \quad (58)$$

Given two models in \mathcal{M}

$$\mathcal{M}_1 : y = G_1(q)u + H_1(q)e_1 \quad (59)$$

$$\mathcal{M}_2 : y = G_2(q)u + H_2(q)e_2 \quad (60)$$

we want to be able to determine which that describes the system better

Therefore, we need to perform an *informative* (open-loop) experiment

Informative Experiments

We want to determine an input signal resulting in data that is *sufficiently informative* to distinguish between models in \mathcal{M}

For two models identified using data that is sufficiently informative, the expectation

$$\overline{\mathbb{E}}[\Delta\epsilon^2] = \lim_{N \rightarrow \infty} \frac{1}{N} \sum_{t=1}^N \mathbb{E}[\Delta\epsilon_t^2] = \int_{-\pi}^{\pi} \phi_1(w) + \phi_2(w) dw = 0 \quad (61)$$

only holds if

$$\phi_2(w) = \left| \frac{H_0 \Delta H}{H_1 H_2} \right|^2 \sigma^2 = 0 \quad \Rightarrow \quad \Delta H(e^{jw}) \equiv 0 \quad (62)$$

$$\phi_1(w) = \left| \frac{1}{H_1} \right|^2 \left| \Delta G + \frac{G_0 - G_2}{H_2} \Delta H \right|^2 \Phi_u(w) = 0 \quad (63)$$

$$\Rightarrow |\Delta G(e^{jw})|^2 \Phi_u(w) \equiv 0 \Rightarrow \Delta G(e^{jw}) \equiv 0 \quad (64)$$

Consequently, the input should have a spectrum $\Phi_u(w)$ for which the above expectation only becomes zero for identical models in \mathcal{M} .

A quasi-stationary signal with spectrum $\Phi_u(w)$ is persistently excited of order n ($pe(n)$) if, for all filters in the form

$$M(q^{-1}) = m_0 + m_1q^{-1} + \dots + m_{n-1}q^{-(n-1)} \quad (65)$$

the relation

$$\Phi_z(w) = |M(e^{jw})|^2\Phi_u(w) = 0, \quad z_t = M(q^{-1})u_t \quad (66)$$

implies that for all w

$$M(e^{jw}) = 0 \quad (67)$$

$M(q^{-1})$ has n parameters and $n - 1$ zeros; implying that $M(q)M(q^{-1})$ has at most $n - 1$ different zeros

Equivalently, the spectrum, $\Phi_u(w)$, has to be non-zero at at least n different points in the interval $w \in [-\pi, \pi]$

The reason is that a signal which is pe(n) cannot be filtered to zero by an MA filter of order $n - 1$, but n or higher might do it

$$u_t = \text{const} \neq 0, \text{ signal is pe}(1) \quad (68)$$

$$M_1(q^{-1}) = 1 - q^{-1} : M_1(q^{-1})u_t = u_t - u_{t-1} = 0 \quad (69)$$

$$M_0(q^{-1}) = 1 : M_0(q^{-1})u_t = u_t \neq 0 \quad (70)$$

or looking at the spectrum: it is always zero

$$\Phi_u = \tilde{d}\delta(w) \quad (71)$$

$$\Phi_{M_1 u} = 2(1 - \cos(w))\tilde{d}\delta(w) = 0 \quad (72)$$

Transfer function

$$G = q^{-k} \frac{B(q)}{A(q)} = q^{-k} \frac{b_0 + b_1 q^{-1} + \dots + b_{n_b} q^{-n_b}}{1 + a_1 q^{-1} + \dots + a_{n_a} q^{-n_a}} \quad (73)$$

The signal u_t has to be $pe(n_b + n_a + 1)$

$$\Delta G = \frac{B_1}{A_1} - \frac{B_2}{A_2} = \frac{B_1 A_2 - B_2 A_1}{A_1 A_2} = 0 \quad \Rightarrow \quad |B_1 A_2 - B_2 A_1|^2 \Phi_u(w) = 0 \quad (74)$$

where it can be seen that the effective part of ΔG has the order $n_b + n_a$

Crest factor (for zero-mean signals)

$$C_r^2 = \frac{\max_t u_t^2}{\lim_{N \rightarrow \infty} \frac{1}{N} \sum_{t=1}^N u_t^2} \quad (75)$$

The crest factor should be as low as possible (the minimum is 1)

For binary signals, $u_t = \pm \bar{u}$, the crest factor is minimum, $C_r^2 = 1$

Consequently, binary signals are useful for linear systems, but cannot, in general, handle nonlinear systems

$$y_t = \frac{B(q)}{A(q)} f(u_t) \quad (76)$$

$$f(u_t) = \alpha \cos(\pm \bar{u}) = \alpha \cos(\bar{u}) \quad (77)$$

Single harmonic signal

$$u_t = A \sin(\omega t), \quad (78)$$

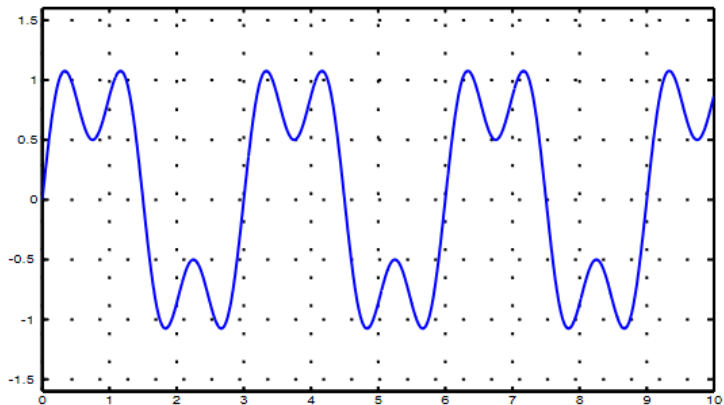
- Two non-zero frequency components in its spectrum (at $\pm\omega$)
- It is $\text{pe}(2)$
- Its crest factor is $C_r^2 = 2$

Sum of sines

$$u_t = \sum_{k=1}^n A_k \sin(\omega_k t + \phi_k) \quad (79)$$

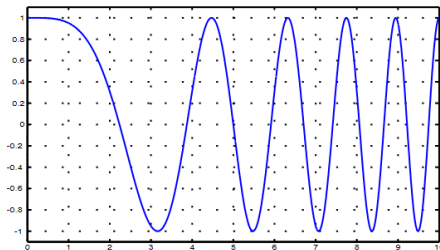
- Two components for each ω_k , so the signal is $\text{pe}(2n)$
- If $\omega_k = 0$ or $\omega_k = \frac{\pi}{T_s}$, the order goes down by 1 to $\text{pe}(2n - 1)$ (by 2 if both)
- The crest factor is, in the worst case, $C_r^2 = 2n$, and lowest if the sinusoids are maximally out of phase

Sum of 2 harmonics, with maximum phase difference (180°)



Single sine function: The chirp signal

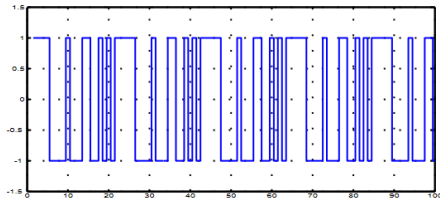
$$u_t = A \sin((w_0 + \alpha t)t), \quad C_r^2 = \sqrt{2} \quad (80)$$



PRBS signal

$$z_t = \text{mod}(B(q^{-1})z_{t-1}, 2) \quad (81)$$

B is order m and the signal has the maximum length $M = 2^m - 1$



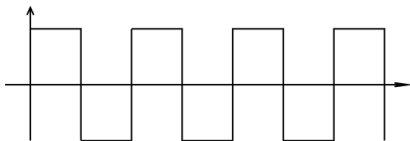
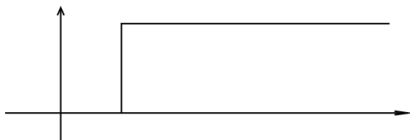
- PRBS signals are deterministic, but have properties similar to those of white noise
- A PRBS signal is $\text{pe}(M - 1)$ and $C_r^2 = 1$

Alternative: Apply random Gaussian signals that are filtered/colored white noise signals

$$u_t = H_u(q)\check{e}_t, \quad \check{e}_t \sim \mathcal{F}_{iid}(0, \sigma_u^2)(white) \quad (82)$$

- In practice, we would have to use a truncated Gaussian to keep the control bounded, e.g., within $\pm 3\sigma$ ($\approx 99\%$ coverage), resulting in $C_r^2 = 3$
- Random binary signals can be generated by taking the sign of a suitable Random Gaussian signal

Step and square wave signals are also commonly used



For a step at time M and a square (both between d_0 and d_1)

$$C_r^2 = \frac{d_1^2}{\lim_{N \rightarrow \infty} \frac{Md_0^2 + (N-M)d_1^2}{N}} = \frac{d_1^2}{d_1^2 + \lim_{N \rightarrow \infty} \frac{M}{N}d_0^2} = 1, \quad C_r^2 = \frac{d_1^2}{\frac{1}{2}d_1^2 + \frac{1}{2}d_0^2}$$

The pulse can also be represented as an infinite harmonic sum

ARMAX model

$$A(q^{-1})y_t = B(q^{-1})u_t + C(q^{-1})e_t, \quad e_t \sim N(0, 0.05) \quad (83)$$

Polynomials

$$A(q^{-1}) = 1 - 1.2q^{-1} + 0.8q^{-2}, \quad (84)$$

$$B(q^{-1}) = -0.5q^{-1} + 0.2q^{-2}, \quad (85)$$

$$C(q^{-1}) = 1 + 0.3q^{-1} \quad (86)$$

Define system

```
>> A = [1, -1.2, 0.8]; D = 1;
>> B = [0, -0.5, 0.2]; F = 1;
>> C = [1, 0.3]; R = 0.05;
>> Ts = 1;
>> M = idpoly(A, B, C, D, F, R, Ts);
```

Simulate system

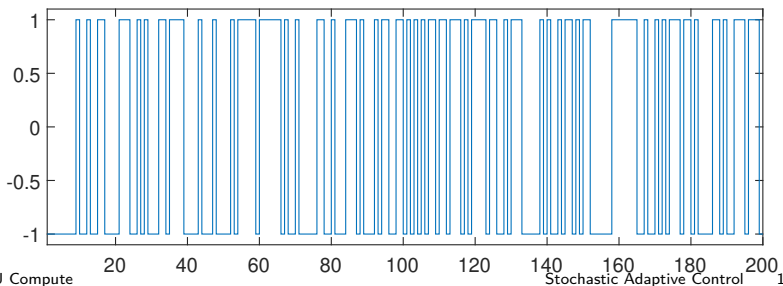
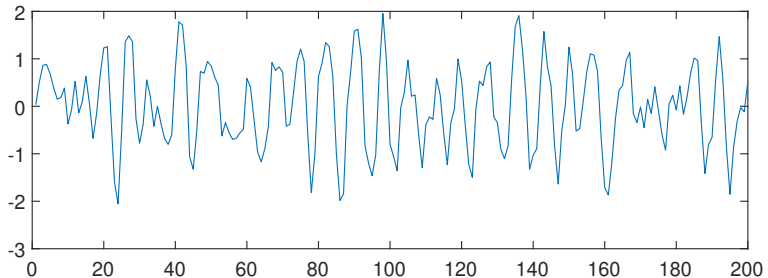
```
>> N = 200;
>> u = iddata([], idinput(N, 'prbs'));
>> e = iddata([], sqrt(R)*randn(N, 1));
>> y = sim(M, [u, e]);
>> simdata = [y.y, u.u];
```

Visualize simulation

```
>> figure(1);
>> subplot(211);
>> plot(y.y);
>> subplot(212);
>> stairs(u.u([1:end, end]));
>> xlim([1, numel(u.u)]);
>> ylim([-1.1, 1.1]);
```

Stochastic Adaptive Control - Time-variant Estimation

Informative experiments - Example



ARMAX model

$$A(q^{-1})y_t = B(q^{-1})u_t + C(q^{-1})e_t, \quad e_t \sim N(0, 0.05) \quad (87)$$

Polynomials

$$A(q^{-1}) = 1 + 1.3q^{-1} - 0.7q^{-2}, \quad (88)$$

$$B(q^{-1}) = -0.3q^{-1} - 0.1q^{-2}, \quad (89)$$

$$C(q^{-1}) = 1 + 0.5q^{-1} \quad (90)$$

Define system

```
>> A = [1, 0.8, -0.1];   D = 1;
>> B = [0, -0.3, -0.2]; F = 1;
>> C = [1, 0.5];       R = 0.05;
>> Ts = 1;
>> M = idpoly(A, B, C, D, F, R, Ts);
```

Simulate system

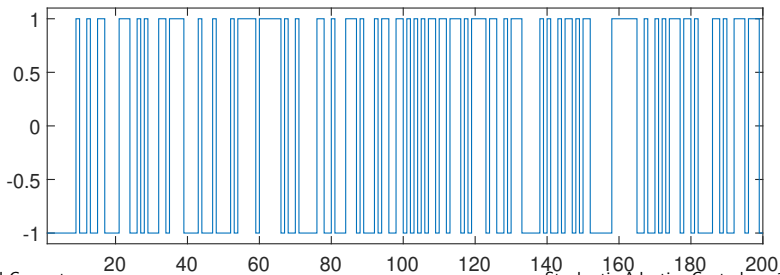
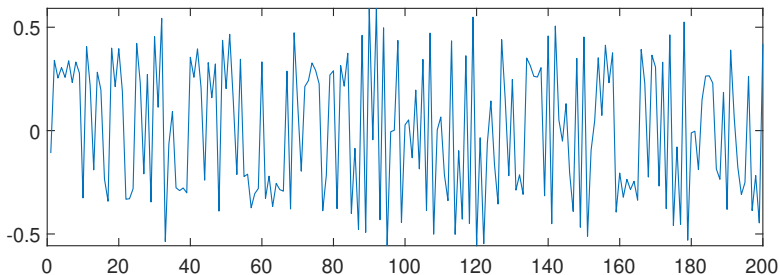
```
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```

Visualize simulation

```
>> figure(1);
>> subplot(211);
>> plot(y.y);
>> subplot(212);
>> stairs(u.u([1:end, end]));
>> xlim([1, numel(u.u)]);
>> ylim([-1.1, 1.1]);
```


Stochastic Adaptive Control - Time-variant Estimation

Informative experiments - Exercise

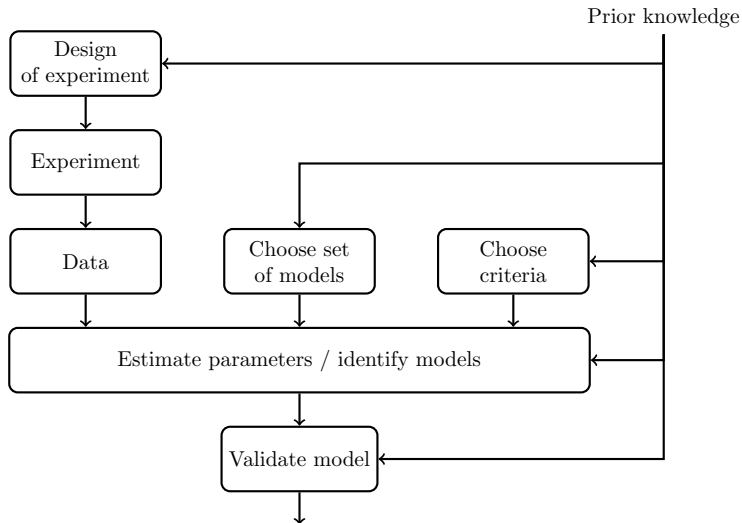


Model validation

We now know how to estimate a model, but how do we check if it is correctly estimated when we don't know the true parameters?

Essentially, we are asking the following two questions

- ① Is our model too simple?
- ② Is our model too complex?



Three available quantities for validation

- 1 the estimated parameters
- 2 the uncertainty (the variance)
- 3 the undescribed model parts (the residuals)

The last is the source of measurement deviations

$$\text{measurement}(y) = \text{model}(\theta, u) + \text{residual}(\epsilon) \quad (91)$$

Question: Does our model have too many parameters?

Unbiased estimate

$$\hat{\theta} \sim \mathcal{F}(\theta, P) \quad (92)$$

θ_i is *significant* if it, with reasonable certainty, is different from zero

Use a *marginal parameter test* to validate that a parameter is significant

For sufficiently many measurements, the distribution approaches a normal distribution

$$\hat{\theta} \sim N(\theta, P) \quad (93)$$

If the following holds, θ_i is, with $(1 - \alpha)\%$ confidence not insignificant

$$|\hat{\theta}_i| > f_{1-\frac{\alpha}{2}} \sqrt{P_{i,i}} \quad (94)$$

f_x is the x th quantile of the standard normal distribution. This approach requires that the variance, P , is known

If the variance, P , was estimated, use the t-distribution

$$z_i = \frac{\hat{\theta}_i}{\sqrt{P_{i,i}}} \sim t(M - d_p) \quad (95)$$

d_p is the number of parameters and M is the number of measurements

If the following holds, θ_i is, with $(1 - \alpha)\%$ confidence not insignificant

$$|\hat{\theta}_i| > f_{1-\frac{\alpha}{2}}^t(M - d_p) \sqrt{P_{i,i}} \quad (96)$$

f_x^t is the x th quantile of the t-distribution. Again, if $M \gg d_p$, this will approach the normal distribution

Model Validation - Multiple Insignificant Parameters?

More than one parameter might be insignificant, but we cannot tell whether its some or all

But we can test whether all parameters in a subset θ_b are significant

$$\hat{\theta} = \begin{bmatrix} \hat{\theta}_a \\ \hat{\theta}_b \end{bmatrix} \sim N \left(\begin{bmatrix} \theta_a \\ \theta_b \end{bmatrix}, \begin{bmatrix} P_a & P_{ab} \\ P_{ab}^T & P_b \end{bmatrix} \right) \quad (97)$$

Test statistic for the hypothesis of insignificant parameters ($\theta_b = 0$)

$$z_b = \hat{\theta}_b^T P_b^{-1} \hat{\theta}_b \sim F(d_b, M - d_p) \quad (98)$$

If the following holds, all parameters in θ_b are, with $(1 - \alpha)\%$ confidence significant

$$z_b > f_{1-\alpha}^F(d_b, M - d_p) \quad (99)$$

d_b is the size of the subset and f_x^F is the x th quantile of the F-distribution. For large M , we can apply a $\chi^2(d_b)$ instead of the F-distribution

Model Reduction

Distribution of parameter estimates

$$\begin{bmatrix} \theta_a \\ \theta_b \end{bmatrix} \sim N \left(\begin{bmatrix} \hat{\theta}_a \\ \hat{\theta}_b \end{bmatrix}, \begin{bmatrix} P_a & P_{ab} \\ P_{ab}^T & P_b \end{bmatrix} \right) \quad (100)$$

If a subset of the parameters, $\hat{\theta}_b$, is insignificant, we can reduce the model using the projection theorem

$$\theta_a | \theta_b \sim N(\hat{\theta}_a, \bar{P}_a) \quad (101)$$

$$\hat{\theta}_a = \hat{\theta}_a - P_{ab}^T P_b^{-1} \hat{\theta}_b \quad (102)$$

$$\bar{P}_a = P_a - P_{ab} P_b^{-1} P_{ab}^T \quad (103)$$

When have we used the projection theorem before and what for?

Think about it for yourself for one minute and then discuss with the person next to you for one minute.

Insignificant: singular analysis of the variance matrix P

Most estimations methods involve solving linear equations in the form

$$H\hat{\theta} = g \quad (104)$$

H is a measure of the data set related to the variance, $H^{-1} = P$

$$P = \left(\sum_{i=0}^N \psi_i \psi_i^T \right)^{-1} \sigma^2 \quad (105)$$

- If a model is overparameterized, then (in the ideal case) H will be singular
- In the less ideal case, H is invertible, but has eigenvalues that are significantly smaller than the rest

$$\text{eig}(H)_i \ll \text{eig}(H)_j \quad \Leftrightarrow \quad \text{eig}(P)_i \gg \text{eig}(P)_j \quad (106)$$

- This requires that the system is sufficiently excited – insufficiently excited systems will result in similar issues

Another way to evaluate if a model is overparameterized is to consider the condition number of its variance.

$$\text{cond}(P) = \frac{|\lambda_{\max}|}{|\lambda_{\min}|}, \quad \lambda = \text{eig}(P) \quad (107)$$

where λ_{\min} and λ_{\max} are the smallest and largest eigenvalues of P

If $\text{cond}(P)$ is large, it indicates overparameterization

Example:

$$\text{Model 1: } \text{cond}(P_1) = 1000 \quad (108)$$

$$\text{Model 2: } \text{cond}(P_2) = 40 \quad (109)$$

Model 1 appears to be too complex, while model 2 is more balanced

If the model is overparameterized, some zeros and poles might be close to each other

$$y_t = H_{yu}(q)u_t + H_{ye}(q)e_t \quad (110)$$

Use linearization to approximate uncertainty in zero and poles

$$\hat{p}_i = f_i(\hat{\theta}) \simeq f_i(\theta) + \frac{\partial f_i}{\partial \theta} \tilde{\theta}, \quad \tilde{\theta} \sim N(0, P) \quad (111)$$

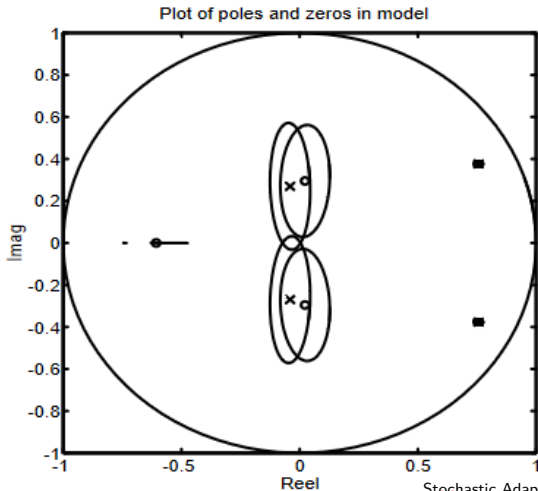
$$\hat{p}_i \sim N \left(p_i, \frac{\partial f_i}{\partial \theta} P \left(\frac{\partial f_i}{\partial \theta} \right)^T \right) \quad (112)$$

If the confidence intervals of a pole and a zero overlap, it is a strong indication that they cancel each other out

Hint: Use Matlab's `zppplot`

$$(1 - 1.5q^{-1} + 0.7q^{-2})y_t = (1 - 0.5q^{-1})u_t + e_t \quad (113)$$

$$(1 - a_1q^{-1} + \dots + a_4q^{-4})y_t = (b_0 + \dots + b_3q^{-3})u_t + e_t \quad (114)$$



Residual Analysis

Question: Is the model too simple?

Residuals

$$\text{measurement}(y) = \text{model}(\theta, u) + \text{residual}(\epsilon) \quad (115)$$

For a perfect model, the residuals would have the following properties

- 1 $\epsilon_t \sim \mathcal{F}(0, \sigma^2)$.
- 2 ϵ_t has a symmetric distribution
- 3 ϵ_t is white
- 4 ϵ_t is uncorrelated with current and prior inputs

Equivalently (in terms of co-variance functions)

$$r_\epsilon(k) = \mathbb{E}[\epsilon_{t+k}\epsilon_t] = \begin{cases} \sigma^2 & k = 0, \\ 0 & \text{otherwise,} \end{cases} \quad r_{\epsilon_t, u_t}(k) = \mathbb{E}[\epsilon_{t+k}u_t] = 0 \quad (116)$$

Important: use one data set for *estimation* and another for the *validation* (cross-validation)

Residual Analysis - mean and variance test

Simple approach: Test whether the distribution of the residuals has the right mean and variance

If the below holds, the residuals are not zero mean

$$|\bar{\epsilon}| > f_{1-\frac{\alpha}{2}}^t(M-1) \sqrt{\frac{S^2}{M}} \quad (117)$$

$$\bar{\epsilon} = \frac{1}{M} \sum_{i=1}^M \epsilon_i, \quad S^2 = \frac{1}{M-1} \sum_{i=1}^M (\epsilon_i - \bar{\epsilon})^2 \quad (118)$$

If either of the below hold, the variance is time-varying

$$\frac{S_1^2}{S_2^2} < f_{\alpha/2}^F(M_1, M_2) \text{ or } \frac{S_1^2}{S_2^2} > f_{1-\alpha/2}^F(M_1, M_2) \quad (119)$$

$$S_i^2 = \frac{1}{M_i} \sum_{j=1}^{M_i} \epsilon_{i+j}^2 \quad (120)$$

Note: The intervals must be non-overlapping

Test for whiteness: The number of sign changes, z , should follow (M is the number of data points)

$$z \sim N\left(\frac{M-1}{2}, \frac{M-1}{4}\right) \quad (121)$$

We reject the hypothesis if either of the below holds

$$z < \frac{M-1}{2} - \sqrt{\frac{M-1}{4}} f_{1-\frac{\alpha}{2}}^N \quad \text{or} \quad z > \frac{M-1}{2} + \sqrt{\frac{M-1}{4}} f_{1-\frac{\alpha}{2}}^N \quad (122)$$

That is, the hypothesis is rejected if the test statistic is outside the confidence interval

Alternative test for whiteness: The auto-covariance must be in the form

$$r_{\epsilon}(k) = \mathbb{E}[\epsilon_{t+k}\epsilon_t] = \begin{cases} \sigma^2 & k = 0 \\ 0 & \text{otherwise} \end{cases} \quad (123)$$

Estimates of auto-covariance and auto-correlation

$$\hat{r}_{\epsilon}(k) = \frac{1}{M} \sum_{t=1}^{M-k} \epsilon_{t+k}\epsilon_t, \quad \hat{\rho}_{\epsilon}(k) = \frac{\hat{r}_{\epsilon}(k)}{\hat{r}_{\epsilon}(0)} \quad (124)$$

Test the covariance at each time step

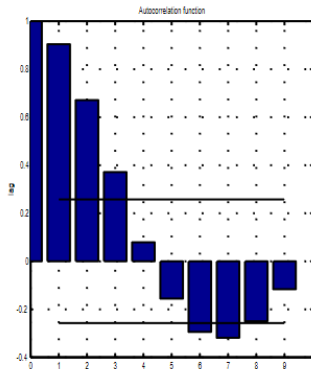
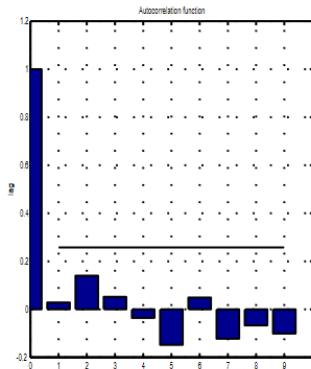
$$H_0 : \sqrt{M}\hat{\rho}_{\epsilon}(k) \sim N(0, 1), \text{ reject if } |\hat{\rho}_{\epsilon}(k)| > \frac{f_{1-\frac{\alpha}{2}}^N}{\sqrt{M}} \quad (125)$$

Test if the covariance is zero for $k \neq 0$

$$H_0 : z = M \sum_{i=1}^m \hat{\rho}_{\epsilon}^2(i) \sim \chi^2(m), \text{ reject if } z > f_{1-\alpha}^{\chi^2}(m) \quad (126)$$

Stochastic Adaptive Control - Time-variant Estimation

Residual Analysis - test of autocorrelation



Test the cross-covariance

$$r_{\epsilon,u}(k) = \mathbb{E}[\epsilon_{t+k}u_t] = 0 \quad (127)$$

Cross-covariance and cross-correlation

$$\hat{r}_{\epsilon,u}(k) = \frac{1}{M} \sum_{i=1}^{M-k} \epsilon_{t+k}u_t, \quad \hat{\rho}_{\epsilon,u}(k) = \frac{\hat{r}_{\epsilon,u}(k)}{\sqrt{\hat{r}_{\epsilon}(0)\hat{r}_u(0)}} \quad (128)$$

Marginal test of the cross-covariance

$$H_0 : \sqrt{M}\hat{\rho}_{\epsilon,u}(k) \sim N(0, 1), \text{ reject if } |\hat{\rho}_{\epsilon,u}(k)| > \frac{f_{1-\frac{\alpha}{2}}^N}{\sqrt{M}} \quad (129)$$

Check if the covariance is zero for $k \neq 0$

$$H_0 : z = M \sum_{i=1}^m \hat{\rho}_{\epsilon,u}^2(i) \sim \chi^2(m), \text{ reject if } z > f_{1-\alpha}^{\chi^2}(m) \quad (130)$$

Alternative test for whiteness: Consider the Fourier transformed residuals

$$X(w_k) = \frac{1}{M} \sum_{t=1}^M \epsilon_t e^{jw_k t} \quad (131)$$

Hint: Matlab's `fft` can be used to compute $X(w_k)$

Estimated spectral density (periodogram)

$$\hat{\phi}(w_k) = |X(w_k)|^2 \quad (132)$$

Hint: Matlab's `etfe` can be used to compute $\hat{\phi}(w_k)$

If x_t is white noise

$$\mathbb{E}[\hat{\phi}(w_k)] = 2\sigma^2 \quad (133)$$

Question: Can we validate a model using a single data set?

Coefficient of determination

$$R^2 = \frac{J_0 - J(\hat{\theta})}{J_0} \quad (134)$$

$$J_0 = \frac{1}{2} \sum_{i=1}^M (y_i - \bar{y})^2, \quad J(\hat{\theta}) = \frac{1}{2} \sum_{i=1}^M \epsilon_i^2 \quad (135)$$

$J(\hat{\theta})$ is the loss-function and a perfect model results in $R^2 = 1$. Lower values of R^2 indicate worse models

Alternative loss functions

$$W(\hat{\theta}) = \sum_{i=1}^M \epsilon_i^2, \quad W_M(\hat{\theta}) = \frac{1}{M} \sum_{i=1}^M \epsilon_i^2 \quad (136)$$

The loss functions are monotonically decreasing with model complexity

Objective: Compare two model classes, \mathcal{M}_1 and \mathcal{M}_2 using the F-test

Hypothesis: $\mathcal{M}_{\text{true}} \subset \mathcal{M}_1 \subset \mathcal{M}_2$ where $d_2 \geq d_1$ are the number of model parameters. Consequently, the loss-function $J_i = J_i(\hat{\theta})$ does not decrease significantly by increasing the model size if $\mathcal{M}_1 \subset \mathcal{M}_2$

Test statistic

$$H_0 : z = \frac{J_1 - J_2}{J_2} \frac{M - d_2}{d_2 - d_1} \sim F(d_2 - d_1, M - d_2) \quad (137)$$

Reject hypothesis if

$$z > f_{1-\alpha}^F(d_2 - d_1, M - d_2) \quad (138)$$

Residual Analysis - model comparison: Information Criteria

Information criteria

- ① Akaike's Information Criterion (AIC); tends towards higher complexity

$$AIC = \left(1 + \frac{2d}{M}\right) W_M \quad (139)$$

- ② Bayesian Information Criterion (BIC);

$$BIC = \left(1 + \frac{\log(M)d}{M}\right) W_M \quad (140)$$

- ③ Akaike's Final Prediction Error (FPE) Criterion; expresses the variance of the prediction error, also $FPE \rightarrow AIC, M \gg d$

$$FPE = \frac{M+d}{M-d} W_M = \left(1 + \frac{2d}{M-d}\right) W_M \quad (141)$$

If two models have the same d , choose the one with the lowest loss function

Information criteria

$$AIC = \left(1 + \frac{2d}{M}\right) W_M \quad (142)$$

$$BIC = \left(1 + \frac{\log(M)d}{M}\right) W_M \quad (143)$$

$$FPE = \frac{M+d}{M-d} W_M = \left(1 + \frac{2d}{M-d}\right) W_M \quad (144)$$

What happens to the criteria as $M \rightarrow \infty$?

Think about it for yourself for one minute and then discuss with the person next to you for one minute.

Model validation - Example

ARMAX model

$$A(q^{-1})y_t = B(q^{-1})u_t + C(q^{-1})e_t, \quad e_t \sim N(0, 0.05) \quad (145)$$

Polynomials

$$A(q^{-1}) = 1 - 1.2q^{-1} + 0.8q^{-2}, \quad (146)$$

$$B(q^{-1}) = -0.5q^{-1} + 0.2q^{-2}, \quad (147)$$

$$C(q^{-1}) = 1 + 0.3q^{-1} \quad (148)$$

Define system

```
>> A = [1, -1.2, 0.8]; D = 1;
>> B = [0, -0.5, 0.2]; F = 1;
>> C = [1, 0.3]; R = 0.05;
>> Ts = 1;
>> M = idpoly(A, B, C, D, F, R, Ts);
```

Simulate system

```
>> N = 200;
>> u = iddata([], idinput(N, 'prbs'));
>> e = iddata([], sqrt(R)*randn(N, 1));
>> y = sim(M, [u, e]);
>> simdata = [y.y, u.u];
```

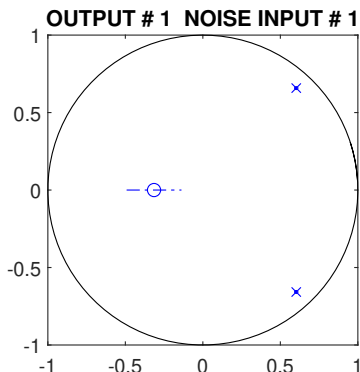
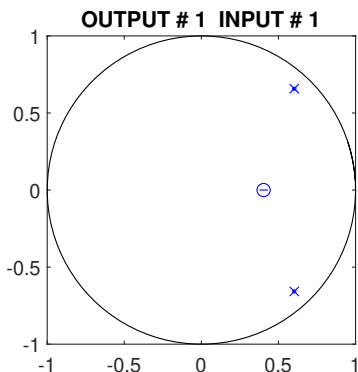
Model validation - Example

Estimate parameters

```
>> th = armax(simdata, [2, 2, 1, 1])
```

Zero-pole cancellation

```
>> figure(1)
>> subplot(121)
>> zpplot(th2zp(th, 1), norminv(0.995)) % 99% Gaussian CI
>> subplot(122)
>> zpplot(th2zp(th, 0), norminv(0.995)) % 99% Gaussian CI
```



Eigenvalue analysis

```
>> [TH,P] = th2par(th);  
>> log10(eig(P))
```

```
ans =
```

```
-5.1145  
-4.8291  
-4.7172  
-3.8705  
-2.3304
```

Insignificant parameters

```
>> TH + sqrt(diag(P))*[-1,1]*norminv(0.995)
```

```
ans =
```

```
-1.2319    -1.1860  
 0.7790     0.8178  
-0.5121    -0.4936  
 0.1862     0.2139  
 0.1555     0.5077
```

Model validation - Exercise

ARMAX model

$$A(q^{-1})y_t = B(q^{-1})u_t + C(q^{-1})e_t, \quad e_t \sim N(0, 0.05) \quad (149)$$

Polynomials

$$A(q^{-1}) = 1 - 1.2q^{-1} + 0.8q^{-2}, \quad (150)$$

$$B(q^{-1}) = -0.5q^{-1} + 0.2q^{-2}, \quad (151)$$

$$C(q^{-1}) = 1 + 0.3q^{-1} \quad (152)$$

Estimate the parameters in a model with

$$n_a = 3, \quad (153)$$

$$n_b = 3, \quad (154)$$

$$n_c = 2, \quad (155)$$

$$k = 1 \quad (156)$$

Solve the exercise in 15 min.

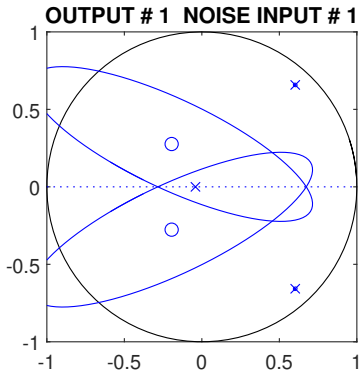
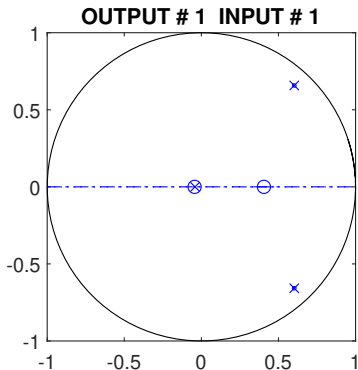
Model validation - Exercise

Estimate parameters

```
>> th = armax(simdata, [3, 3, 2, 1])
```

Zero-pole cancellation

```
>> figure(1)
>> subplot(121)
>> zpplot(th2zp(th, 1), norminv(0.995)) % 99% Gaussian CI
>> subplot(122)
>> zpplot(th2zp(th, 0), norminv(0.995)) % 99% Gaussian CI
```



Eigenvalue analysis

```
>> [TH,P] = th2par(th);  
>> log10(eig(P))  
  
ans =  
  
-5.2588  
-5.0938  
-4.8826  
-4.7356  
-4.2485  
-2.5110  
-2.2328  
0.3504
```

Insignificant parameters

```
>> TH + sqrt(diag(P))*[-1,1]*norminv(0.995)  
  
ans =  
  
-2.9920    0.6649  
-1.4583    2.9499  
-1.4232    1.4888  
-0.5054   -0.4884  
-0.7291    1.0882  
-0.3571    0.3749  
-1.4256    2.2050  
-0.4790    0.7077
```

Questions?