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# Practice 4. Optimal Estimation

# **Activity 1. State estimation (Kalman filter)**

Let us consider a water tank that can be modeled through the following equation

$$\frac{\mathrm{dV}(t)}{\mathrm{dt}} = q_{\mathrm{in}}(t) - q_{\mathrm{out}}(t)$$

where V(t) is the tank volume,  $q_{\rm in}(t)$  is the output flow by pump and  $q_{\rm out}(t)$  is the output flow controllerd by a valve. The input flow provided by the pump is controlled through the voltage  $v_h(t)$  according to  $q_{\rm in}(t) = k_b v_b(t)$ . On the other hand, the output flow  $q_{\rm out}(t)$  through valve is given by  $q_{\rm out}(t) = k_v V(t)$ .

Considering that the system parameters are  $k_b=2$  and  $k_v=0.1$ , we want to the problem of states estimation using optimization. In particular, the tank volume V should be estimated using the input voltage  $v_b$  and the output flow measurement  $q_{\rm out}$  when there is a noncontrolled system noise (disturbance)  $w_k$  (Gaussian with zero mean and variance equal to Q=0.1) and a non-controlled measurement noise  $k_v$  (Gaussian with zero mean and variance equal to R=0.1). The value of N=100 input voltages ( $v_b(k)$ ) and output flows ( $q_{\rm out}(k)$ ) are acquired with a sampling time equal to T=1s. They can be generated as indicated in the guide of the lab. Use as initial conditions of the state estimation filter:  $\hat{V}_1=10$  and  $P_1=0.2$ .

# 1.a - Problem formulation and Data generation

```
% Problem formulation
T=1; % sample time
kv=0.1; % parameter kv
kb=2; % parameter kb
A=exp(-kv*T); % parameter a (using exact discretization)
B=(1-exp(-kv*T))*kb/kv; % parameter b (using exact discretization)
C=kv; % parameter c
D=0; % parameter d
```

For calculation of Pareto plot we change the variance *Q* and then *R* using a for loop. Here we also define initial conditions.

```
p = 0;
for e = 0.1:0.1:2
p=p+1; % index for pareto plot
```

```
Q=0.1; % variance of system noise (disturbance)
R=0.1; % variance of measurement noise
x1=10;P1=0.2; % Initial conditions from estimation
```

We generate data with 100 data points but only one time because we don't want to work with different data all the time. The data are generated using the following functions:

$$x(k+1) = A \cdot x(k) + B \cdot u(k) + w(k)$$
$$y(k) = C \cdot x(k) + D \cdot u(k) + v(k)$$

# Kalman filter via numerical optimization

From saved data we extract our input u and output y.

Then we formulate our optimization of the prediction problem using the following minimization function.

$$\begin{aligned} & \text{min} & & J = \underbrace{\sum_{k=1}^{N} w(k)^T \cdot Q^{-1} \cdot w(k)}_{k=1} + \underbrace{\sum_{k=1}^{N} v(k)^T \cdot R^{-1} \cdot v(k)}_{k=1} + \underbrace{w(1)^T \cdot P_1^{-1} \cdot w(1)}_{1} \\ & \text{s.t.} & & w(k) = \hat{x}(k+1) - \left(A \cdot \hat{x}(k) + B \cdot u(k)\right) \quad k = 2, ..., N \\ & & v(k) = y(k) - \left(C \cdot \hat{x}(k) + D \cdot u(k)\right) \quad k = 1, ..., N \\ & & w(1) = x(1) - \hat{x}(1) \end{aligned}$$

min 
$$J = \sum_{k=1}^{N} (\hat{x}(k+1) - A \cdot \hat{x}(k) - B \cdot u(k))^{T} \cdot Q^{-1} \cdot (\hat{x}(k+1) - A \cdot \hat{x}(k) - B \cdot u(k)) + \sum_{k=1}^{N} (y(k) - C \cdot \hat{x}(k) - D \cdot u(k))^{T} \cdot R^{-1} \cdot (y(k) - C \cdot \hat{x}(k) - D \cdot u(k)) + (x_{1} - \hat{x}(1))^{T} \cdot P_{1}^{-1} \cdot (x_{1} - \hat{x}(1))$$

To introduce this problem to yalmip which will solve it numerically, we initialize variable x and define our objective function and constraints.

```
x = sdpvar(repmat(1,1,N+1),repmat(1,1,N+1));
objective = (x1-x{1})'*P1^(-1)*(x1-x{1}); % disturbance
constraints = [];
for k = 1:N
    objective = objective+...
    (x{k+1}-A*x{k}-B*u(k))'*Q^(-1)*(x{k+1}-A*x{k}-B*u(k))...
    +(y(k)-C*x{k})'*R^(-1)*(y(k)-C*x{k});
end
```

And finally we solve the problem using *quadprog* solver and *yalmip*.

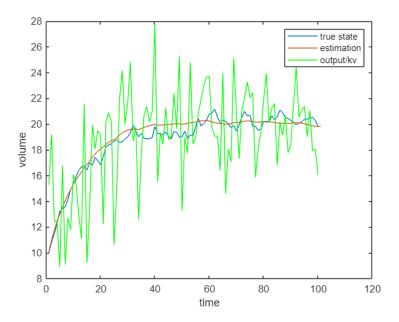
```
options = sdpsettings('solver', 'quadprog');
optimize(constraints, objective, options);
```

We extract the value of estimated state

```
x_est = zeros(N+1,1);
for k = 1:N+1
    x_est(k,1) = value(x{k});
end
```

And plot the result but only for Q = 0.1, other values are only used for calculation of pareto plot.

```
if Q == 0.1 && R == 0.1
    x_est_a = x_est;
    figure
    plot(data(:,2)); hold on
    plot(x_est)
    plot(y/kv,'g')
    legend('true state','estimation','output/kv')
    xlabel('time')
    ylabel('volume')
end
```



To calculate Pareto plot we use the following equations

$$\begin{aligned} & \text{min} & & J = \sum_{k=1}^{N} w(k)^T \cdot Q^{-1} \cdot w(k) \\ & \text{s.t.} & & w(k) = \hat{x}(k+1) - \left(A \cdot \hat{x}(k) + B \cdot u(k)\right) \quad k = 2, ..., N \\ & & v(k) = y(k) - \left(C \cdot \hat{x}(k) + D \cdot u(k)\right) \quad k = 1, ..., N \\ & & w(1) = x(1) - \hat{x}(1) \end{aligned}$$

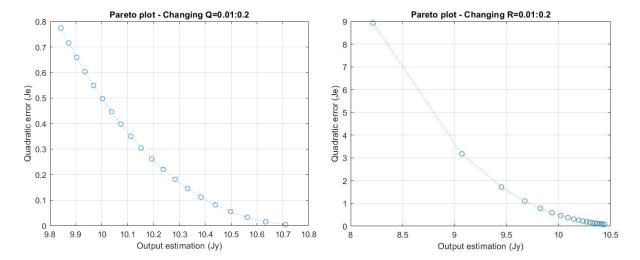
Where w(k) is the system noise and v(k) is the measurement noise regarding the state.

```
Je=0;
Jy=0;
for k=1:N
    w(k)=x_est(k+1)-(A*x_est(k)+B*u(k));    % System noise (disturbance)
    v(k)=y(k)-(C*x_est(k)+D*u(k));    % Measurement noise
    Je = Je + w(k)'*w(k);    % Quadratic error
    Jy = Jy + v(k)'*v(k);    % Output estimation
end

all_Je(p) = Je;
all_Jy(p) = Jy;

end % end of for loop for pareto plot
figure
plot(all_Jy,all_Je,'o:')
title('Pareto plot - Changing Q=0.01:0.2')
```

```
xlabel('Output estimation (Jy)')
ylabel('Quadratic error (Je)')
grid on
```



On both of these Pareto plots we can see that with increasing system error we have better output estimation and vice versa. That's why we need to find an optimal value of Q and R that will fit our requirements.

### 1.b - Kalman filter with 98% confidence interval

To calculate the the analytical solution we use Kalman filter. The equations for Kalman filter are the following

$$\begin{array}{lll} P(1) & & & & \\ \hat{x}(1) & & & & \\ y(k) & k = 1 = ...N & & & \\ u(k) & k = 1 = ...N & & & \\ \hline \end{array}$$
 
$$\begin{array}{lll} L(k) = A \cdot P(k) \cdot C^T \cdot \left[ R + C \cdot P(k) \cdot C^T \right]^{-1} & & \\ \hat{x}(k+1) = A \cdot \hat{x}(k) + B \cdot u(k) + L(k) \cdot \left( y(k) - C \cdot \hat{x}(k) \right) & & \\ P(k+1) = Q + \left( A - L(k) \cdot C \right) \cdot P(k) \cdot A^T & & \\ \end{array}$$

And then the confidence interval is defined by upper and lower bound with the following equations.

$$lb(k) = x(k) - \alpha \cdot \sqrt{P(k)}$$

$$ub(k) = x(k) + \alpha \cdot \sqrt{P(k)}$$

Where alpha defines our confidence interval and for 98%  $\alpha = 3$ .

```
L=zeros(N,1); % initialization of L matrix
% For 98% -> alpha=3; for 95% -> alpha=2
alpha=3; % select according confidence interval
p = 0;
for e = 0.1:0.1:2
p=p+1; % index for pareto plot
Q=0.1*e; % controll system noise variance
         % measurement noise variance
R=0.1;
x(1)=x1; % initial condition for x;
P(1)=P1; % initial condition for P;
lb(1)=x(1)-sqrt(P(1))*alpha*0.98;
ub(1)=x(1)+sqrt(P(1))*alpha*0.98;
for k=1:N  % codifying Kalman filter in Matlab
   L(k)=A*P(k)*C'/(R+C*P(k)*C');
   x(k+1)=A*x(k)+B*u(k)+L(k)*(y(k)-C*x(k));
   P(k+1)=Q+(A-L(k)*C)*P(k)*A';
   lb(k+1)=x(k+1)-alpha*sqrt(P(k+1)); % lower bound
   ub(k+1)=x(k+1)+alpha*sqrt(P(k+1)); % upper bound
end
```

Displaying the graphs with Q and R equaled to 0.1

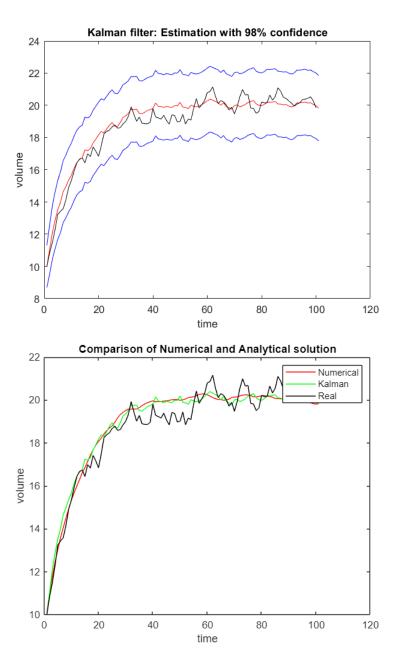
```
if Q == 0.1 && R == 0.1
   x_kal = x;
   figure
   plot(x,'r'),hold on; % plot state estimation
   plot(lb, 'b'); % plot lower bounds interval
    plot(ub, 'b'); % plot upper bound interval
   plot(data(:,2),'k')
   title('Kalman filter: Estimation with 98% confidence')
   xlabel('time')
   ylabel('volume')
   figure
    plot(x_est_a,'r'), hold on
   plot(x, 'g')
    plot(data(:,2),'k')
   legend('Numerical','Kalman','Real')
   title('Comparison of Numerical and Analytical solution')
```

```
xlabel('time')
ylabel('volume')
end
```

Calculating the Pareto plot for Kalman filter.

```
Je=0;
Jy=0;
for k=1:N
    w(k)=x(k+1)-(A*x(k)+B*u(k));  % System noise (disturbance)
    v(k)=y(k)-(C*x(k)+D*u(k));  % Measurement noise
    Je = Je + w(k)'*w(k);  % Quadratic error
    Jy = Jy + v(k)'*v(k);  % Output estimation
end

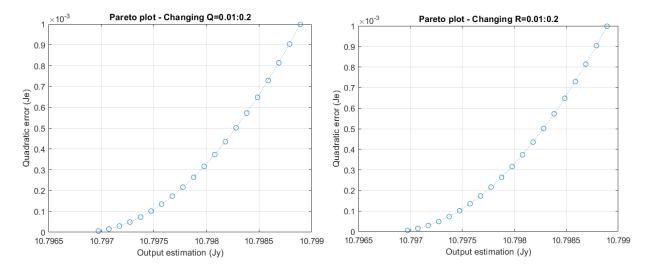
all_Je_A(p) = Je;
all_Jy_A(p) = Jy;
end % end of for loop for pareto plot
```



On the first graph we can see red line which is estimated state by Kalman and blue lines represent upper and lower bound.

Comparing the analytical Kalman solution and numerical using quadprog, we can see that they are quite similar. The Kalman filter looks a bit smoother than a state estimated numerically

```
figure
plot(all_Jy_A,all_Je_A,'o:')
title('Pareto plot - Changing Q=0.01:0.2')
xlabel('Output estimation (Jy)')
ylabel('Quadratic error (Je)')
```



On the Pareto plots for the Kalman filter we can see different correlation between quadratic error and output estimation. Higher the quadratic error is, higher the output estimation error is and vice versa.

# 1.c - Steady-State Kalman

To calculate the steady-state Kalman we needed to solve the following equation to find P.

$$L(k) = A \cdot P(k) \cdot C^{T} \cdot \left[ R + C \cdot P(k) \cdot C^{T} \right]^{-1}$$
$$P(k+1) = Q + \left( A - L(k) \cdot C \right) \cdot P(k) \cdot A^{T}$$

$$L = A \cdot P \cdot C^{T} \cdot \left[ R + C \cdot P \cdot C^{T} \right]^{-1}$$
$$P = Q + \left( A - L \cdot C \right) \cdot P \cdot A^{T}$$

$$P = Q + \left(A - A \cdot P \cdot C^T \cdot \left[R + C \cdot P \cdot C^T\right]^{-1} \cdot C\right) \cdot P \cdot A^T$$

We substituted L in P with its equation and found 2 possible solutions.

```
syms z
z=vpa(solve(z==Q+(A-A*z*C'*inv(R+C*z*C')*C)*z*A',z,'Real',true))

z =
    (-2.1729057180862011395865087288967
    0.46021324886601872041966785430985)
```

We are gonna pick the one that is positive, because that is the correct one for us.

```
PSS=z(2)

PSS = 0.46021324886601872041966785430985
```

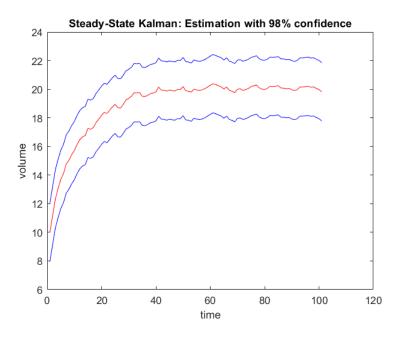
And with the PSS we are able to calculate our steady-state gain LSS.

```
LSS=A*PSS*C*(R+C*PSS*C)^-1

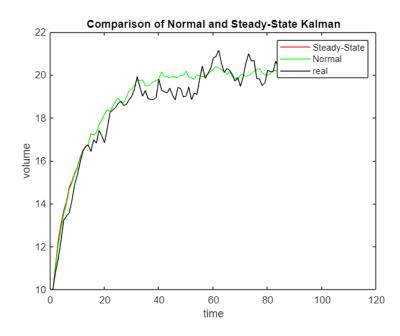
LSS = 0.39809720695226966940828200508746
```

Now we can calculate all the states using constants PSS and LSS.

```
% Kalman filter with confidence interval using steady state
xss=zeros(N+1,1); % initialization of states
lbss=zeros(N+1,1); % lower bounds of states
ubss=zeros(N+1,1); % upper bounds of states
alpha=3; % choose the desired confidence
xss(1)=x1; % initial condition for x;
lbss(1)=x(1)-alpha*sqrt(PSS);
ubss(1)=x(1)+alpha*sqrt(PSS);
for k=1:N % codifying steady state Kalman filter
    xss(k+1)=A*xss(k)+B*u(k)+LSS*(y(k)-C*xss(k));
    lbss(k+1)=xss(k+1)-alpha*sqrt(PSS);
    ubss(k+1)=xss(k+1)+alpha*sqrt(PSS);
end
figure
plot(xss,'r'); % plot state estimation
hold on;
plot(lbss,'b'); % plot lower bounds interval
plot(ubss, 'b'); % plot upper bound interval
title('Steady-State Kalman: Estimation with 98% confidence')
xlabel('time')
ylabel('volume')
```



```
figure
plot(xss,'r'), hold on;
plot(x_kal,'g');
plot(data(:,2),'k');
legend('Steady-State','Normal','real')
title('Comparison of Normal and Steady-State Kalman')
xlabel('time')
ylabel('volume')
```



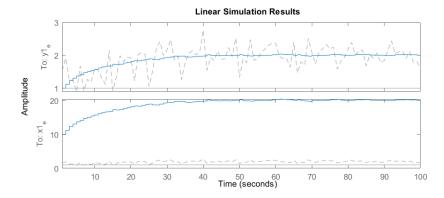
When comparing the Normal Kalman with changing P and L and Steady-State Kalman we can see that they are very similar, almost the same even if the the Steady-State Kalman uses constant PSS and LSS. We can also see that the biggest difference is in the beggining when the P and L are still changing but after some time they become constant.

Steady state Kalman with MATLAB function

```
% MATLAB Kalman function
SYS=ss(A,[B 1],C,[0 0],-1);
[KEST,L,P] = kalman(SYS,Q,R,'delayed')
KEST =
  A =
         x1 e
   x1_e 0.865
  B =
           u1
                   у1
   x1_e 1.903 0.3981
        x1_e
   y1_e 0.1
   x1_e
       u1 y1
   v1 e 0 0
   x1_e 0 0
Input groups:
                Channels
      Name
                1
    KnownInput
    Measurement
                    2
Output groups:
                  Channels
        Name
    OutputEstimate 1
    StateEstimate
Sample time: unspecified
Discrete-time state-space model.
L = 0.3981
P = 0.4602
```

Using the kalman MATLAB function we get the same results of PSS and gain LSS as in our steady-state Kalman.

```
figure('Position',[0 0 1200 400])
lsim(KEST,[u,y],1:100,x1)
```

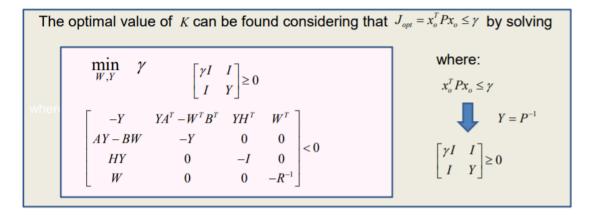


# 1.d - Steady state Kalman filter using LQR LMIs and Duality

Firsrt we initialize our system variables.

```
T=1; % sample time
kv=0.1; % parameter kv
kb=2; % parameter kb
a=exp(-kv*T); % parameter a (using exact discretization)
b=(1-exp(-kv*T))*kb/kv; % parameter b (using exact discretization)
c=kv; % parameter c
d=0; % parameter d
Q=0.1; % deviation of system noise (disturbance)
R=0.1; % deviation of measurement noise
```

We define our matrices, objective function and constraints. The optimal value of LSS can be found solving the ricatti equation for the LQR that is formulated as LMI and constrained as on the following image.



```
%LMI solution
A=a'; % Duality principle for observer design
B=c'; % Duality principle for observer design
Y = sdpvar(1,1);
```

```
W = sdpvar(1,1);
gamma=sdpvar(1,1);
H=sqrt(Q);
F = [Y >= 0];
F = [F, [gamma*eye(1) eye(1);eye(1) Y]>=0];
F = [F, [-Y]]
                    Y*A'-W'*B'
                                    Y*H'
         A*Y-B*W
                        -Y
                                     0
                                                 0;
            H*Y
                         0
                                    -1
                                                 0;
                         0
                                     0
                                           -R^(-1)]
             W
                                                     <= 0];
% Solving the problem
options = sdpsettings('solver', 'sedumi');
optimize(F,gamma,options)
SeDuMi 1.3.5 by AdvOL, 2005-2008 and Jos F. Sturm, 1998-2003.
Alg = 2: xz-corrector, theta = 0.250, beta = 0.500
eqs m = 3, order n = 8, dim = 22, blocks = 3
nnz(A) = 9 + 0, nnz(ADA) = 7, nnz(L) = 5
 it:
         b*v
                   gap
                          delta rate t/tP* t/tD*
                                                      feas cg cg prec
  0:
                1.66E+01 0.000
  1 : -1.04E+00 1.39E+00 0.000 0.0838 0.9900 0.9900
                                                     1.38 1 1 2.1E+00
  2 : -6.47E-01 7.36E-01 0.000 0.5297 0.9000 0.9000
                                                    1.56 1 1 1.5E+00
  3 : -4.77E-01 3.25E-01 0.000 0.4412 0.9000 0.9000
                                                     1.31 1 1 6.8F-01
  4 : -4.49E-01 1.02E-01 0.000 0.3133 0.9000 0.9000
                                                     1.20 1 1 1.9E-01
  5 : -4.54E-01 2.90E-02 0.000 0.2854 0.9000 0.9000
                                                     1.09 1 1 5.1E-02
      -4.57E-01 8.72E-03 0.000 0.3003 0.9000 0.9000
                                                     1.03 1 1 1.5E-02
  7 : -4.59E-01 2.62E-03 0.000 0.3002 0.9000 0.9000
                                                     1.01 1 1 4.4E-03
                                                     1.00 1 1 1.2E-03
  8 : -4.60E-01 7.29E-04 0.000 0.2784 0.9000 0.9000
  9 : -4.60E-01 1.79E-04 0.000 0.2460 0.9000 0.9000
                                                     1.00 1 1 3.0E-04
 10 : -4.60E-01 3.81E-05 0.000 0.2126 0.9000 0.9000
                                                     1.00 1 1 6.4E-05
 11 : -4.60E-01 6.98E-06 0.000 0.1832 0.9000 0.9000
                                                     1.00 1 1 1.2E-05
 12 : -4.60E-01 6.44E-07 0.000 0.0923 0.9900 0.9900
                                                     1.00 1 1 1.1E-06
 13 : -4.60E-01 1.40E-07 0.000 0.2167 0.9000 0.9000
                                                     1.00 1 1 2.3E-07
 14 : -4.60E-01 2.65E-08 0.000 0.1899 0.9000 0.9000
                                                    1.00 1 1 4.5E-08
 15 : -4.60E-01 2.18E-09 0.000 0.0821 0.9900 0.9900
                                                    1.00 1 1 3.7E-09
 16 : -4.60E-01 1.73E-10 0.000 0.0793 0.9900 0.9900
                                                    1.00 2 2 2.9E-10
iter seconds digits
                         c*x
                                          b*y
        0.1 9.4 -4.6021324851e-01 -4.6021324871e-01
 |Ax-b| = 4.6e-11, [Ay-c]_+ = 5.2E-11, |x| = 2.2e+00, |y| = 2.4e+00
Detailed timing (sec)
   Pre
                IPM
                            Post
2.500E-02
             8.200E-02
                         3.007E-03
Max-norms: ||b||=1, ||c|| = 10,
Cholesky |add|=0, |skip|=0, ||L.L||=1725.6.
ans = struct with fields:
    yalmipversion: '20210331'
    matlabversion: '9.11.0.1769968 (R2021b)'
       yalmiptime: 0.0536
       solvertime: 0.1134
             info: 'Successfully solved (SeDuMi-1.3)'
          problem: 0
```

```
LSS_lqr = (value(W)*inv(value(Y)))' % Duality principle for observer design

LSS_lqr = 0.3981
```

We can see that this gain calculated using LMI LQR method is the same as the gain calculated by Steady-State Kalman and MATLAB Kalman function.

# 1.e - Moving Horizon Estimation (MHE) via numerical optimization

In MHE we are considering limited data points for estimation in our case N=10.clc, clear

```
% formulation of the optimization problem
N=10; % determine the number of points
x_hat = sdpvar(repmat(1,1,N+1),repmat(1,1,N+1));
x_0=sdpvar(1,1);
yd=sdpvar(N,1);
ud=sdpvar(N,1);
```

We set our matrices and our objective function using the same minimization function as before.

```
A=a;
B=b;
C=c;
Q=0.1;
R=0.1;

objective = (x_0-x_hat{1})'*P1^(-1)*(x_0-x_hat{1});
constraints = [];
for k = 1:N
    objective = objective+...
    (x_hat{k+1}-A*x_hat{k}-B*ud(k))'*Q^(-1)*(x_hat{k+1}-A*x_hat{k}-B*ud(k))...
    +(yd(k)-C*x_hat{k})'*R^(-1)*(yd(k)-C*x_hat{k});
end
```

We create our estimator of the state which we will use to perdict each state by only looking at N points.

```
options = sdpsettings('solver','quadprog');
estimator=optimizer(constraints,objective,options, {yd,ud,x_0},x_hat{N+1});
```

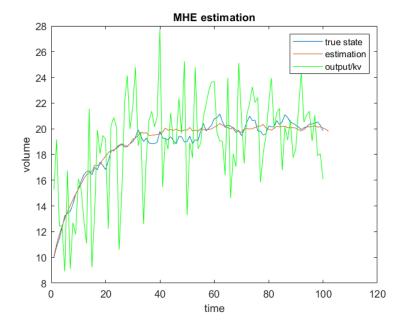
We define first 11 states using the estimation from exercise 1.a in order to calculate other states, because we have to look at the state that is k-N, and the first state we can calculate with MHE method and with control horizon of 10 points is x\_est\_mhe(12).

```
M=100; % Estimation time
```

```
x_est_mhe=zeros(M+1,1); % initialization of states
%initialize the first N state estimation using the result of exercise (a)
x_est_mhe(1:11)=x_est(1:11);
% run the MHE in real time
for k=N+1:M+1
x_est_mhe(k+1)=estimator(y(k-N:k-1),u(k-N:k-1),x_est_mhe(k-N));
end
```

Displaying the results.

```
figure
plot(data(:,2));hold on
plot(x_est_mhe)
plot(y/kv,'g')
title('MHE estimation')
legend('true state','estimation','output/kv')
xlabel('time')
ylabel('volume')
```



MHE is also quite precise state estimation even if we do not consider all the known values to estimate the state.

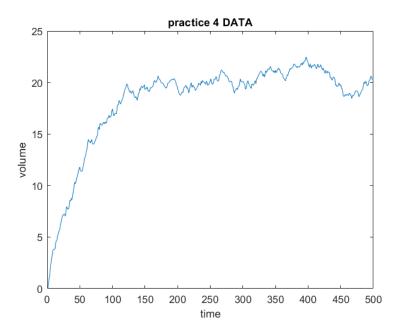
# **Activity 2: Parameter Estimation**

In order to estimate the parameters of the pump  $k_b$  and valve  $k_v$ , input and output data is collected using a sampling time T=0.2s generated as indicated in the guide of the lab. Considering that the model (1) is discretized using the Euler model.

We first generate and save the data.

```
% Data generation
rng(5) % Inizialize random seed (for repetitibility)
T=0.2; % Sample time
kv=0.1;kb=2; % parameters for data generation
A=-kv;B=kb;sys=ss(A,B,1,0); % continuous system sys:dx=A\cdot x+B\cdot u
y=1\cdot x+0\cdot u
% y=x=V=volume (Output = State variable)
% u=v b (Input variable)
dsys=c2d(sys,T); % it converts to discrete system
[nd,dd]=tfdata(dsys,'v'); % it returns numerator and denominator
% as vectors
b=nd(2); a=-dd(2); % x(k)=a\cdot x(k-1)+b\cdot u(k)
N=500; % Number of input/output patterns
y=zeros(N,1); % it inizializes vector y (ouputs)
u=ones(N,1); % it inizializes vector u (inputs)
sigma=0.2; % Deviation of the white noise
% Generating and saving data
for k=2:N % We obtain the data adding WHITE NOISE
y(k)=a*y(k-1)+b*u(k-1)+sigma*randn(1);
end
data=[u y];
save data data % Save data for future calculations
```

Displaying our generated data.



# 2.a - Estimating $k_b$ and $k_v$ using quadprog solver with and without constraints

Without contstraints:

Minimum found that satisfies the constraints.

Optimization completed because the objective function is non-decreasing in feasible directions, to within the value of the optimality tolerance, and constraints are satisfied to within the value of the constraint tolerance.

<stopping criteria details>

# theta\_est = value(theta)

```
theta_est = 2 \times 1
0.9818
```

```
theta_est_2=(C'*C)\C'*Y
                           % another way to obtain theta est
  theta_est_2 = 2 \times 1
     0.9818
      0.3710
 kv_est = (1-theta_est(1))/T %Estimation of kv (Euler approximation)
  kv est = 0.0908
 kb_est = theta_est(2)/T %Estimation of kb (Euler approximation)
  kb est = 1.8548
With constraints:
 Y=y(2:N);
                        % column matrix Y
 C=[y(1:N-1) u(1:N-1)]; % matrix C
 %Numerical Optimization
 J = @(x)(Y-C*x)'*(Y-C*x);
                           % performance index
 theta_est_con=fminsearch(J,[0;0]); % parameters estimators from numerical
 theta = sdpvar(2,1);
 objective= (Y-C*theta)'*(Y-C*theta); % performance index
 constraints=[Y-C*theta<=0.56];</pre>
 options = sdpsettings('solver', 'quadprog');
 optimize(constraints,objective,options); % numerical optimization
  Minimum found that satisfies the constraints.
  Optimization completed because the objective function is non-decreasing in
  feasible directions, to within the value of the optimality tolerance,
  and constraints are satisfied to within the value of the constraint tolerance.
  <stopping criteria details>
 theta_est_con = value(theta);
 kv_est_con = (1-theta_est_con(1))/T % Estimation of kv (Euler approximation)
  kv est con = 0.0908
 kb_est_con = 1.8548
```

We obtained the same value for non-constrained and constrained problem where  $k_{\nu}=0.0908$  and  $k_b=1.8548$ . We used constraint  $Y-C\cdot\theta\leq0.56$ . When we lover the value 0.56 the estimated parameters change.

### 2.b - Non recursive method

theta\_est\_3 =  $2 \times 1$ 

Here we estimate our parameters and their confidence intervals. Confidence intervals are defined by the following equations.

$$\begin{split} & \text{Model} \qquad y(k+1) = a \cdot y(k) + b \cdot u(k) + \epsilon \qquad \epsilon \to N(0,\sigma) \\ & \hat{\theta}(N) = \left[ C(N)^T \cdot C(N) \right]^{-1} \cdot C(N)^T \cdot y(N) \qquad \hat{\theta}(N) = \begin{pmatrix} \hat{a} \\ \hat{b} \end{pmatrix} \qquad \text{estimators of parameters a and b} \\ & s^2 = \frac{J(\hat{\theta})}{N-2} \qquad \qquad \text{estimator of parameter } \sigma^2 \\ & s^2 \cdot \left[ C(N)^T \cdot C(N) \right]^{-1} \qquad \qquad \text{Covariance matrix of estimators } \hat{\theta}(N) = \begin{pmatrix} \hat{a} \\ \hat{b} \end{pmatrix} \\ & \begin{pmatrix} \hat{a} - \sqrt{cov_{11}} \cdot z_{\frac{\alpha}{2}} \\ \hat{a} \end{pmatrix}, \hat{a} + \sqrt{cov_{11}} \cdot z_{\frac{\alpha}{2}} \end{pmatrix} \qquad \qquad \text{Confidence interval for a} \\ & \begin{pmatrix} \hat{b} - \sqrt{cov_{22}} \cdot z_{\frac{\alpha}{2}} \\ \hat{b} \end{pmatrix}, \hat{b} + \sqrt{cov_{22}} \cdot z_{\frac{\alpha}{2}} \end{pmatrix} \qquad \qquad \text{Confidence interval for b} \\ & \text{exact discret: } a = e^{AT} \qquad b = \begin{pmatrix} T \\ 0 \end{pmatrix} e^{A\lambda} d\lambda \end{pmatrix} \cdot B = \frac{B}{A} \cdot \left( e^{AT} - 1 \right) \quad \text{Euler approx } a = 1 + AT \qquad b = BT \end{split}$$

Where we are going to consider 98% confidence interval so alpha=3.

theta\_est\_3 = (C'\*C)\C'\*Y %non-recursive least squares formula

```
%Confidence intervals
sigma2=J(theta_est)/(N-2); % estimation of noise variance
sigma2b = sigma2; % saving variance for 2.d
Cov_mat=sigma2*(C'*C)'; % covariance matrix of the parameters estimators
std_error=sqrt(diag(Cov_mat)); % standard errors

kv_est=(1-theta_est(1))/T; % point 1 estimation of kv using Euler approximation
lbkv=(1-theta_est(1)-std_error(1)*3)/T; % lower bound for 98% confidence
interval for kv
ubkv=(1-theta_est(1)+std_error(1)*3)/T; % upper bound for 98% confidence
interval for kv
% Exact determination of confidence interval for kb_est is too difficult
kb_est=theta_est(2)/T; % point estimation of kb using Euler approximation
```

```
lbkb=(theta_est(2)-std_error(2)*3)/T; % lower bound for 98% confidence interval
for kb
ubkb=(theta_est(2)+std_error(2)*3)/T; % upper bound for 98% confidence interval
for kb
```

```
Calculation using arx.
 sys_arx = arx([y,u],[1,1,0])
  Warning: For transient data (step or impulse experiment), make sure that the change in input
  signal does not happen too early relative to the order of the desired model. You can achieve
  this by prepending sufficient number of zeros (equilibrium values) to the input and output
  signals. For example, a step input must be represented as [zeros(nx,1); ones(N,1)] rather
  than ones(N,1), such that nx > model order.
  sys_arx =
  Discrete-time ARX model: A(z)y(t) = B(z)u(t) + e(t)
    A(z) = 1 - 0.9818 z^{-1}
    B(z) = 0.371
  Sample time: 1 seconds
  Parameterization:
     Polynomial orders: na=1 nb=1 nk=0
     Number of free coefficients: 2
     Use "polydata", "getpvec", "getcov" for parameters and their uncertainties.
  Status:
  Estimated using ARX on time domain data.
  Fit to estimation data: 95.43% (prediction focus)
  FPE: 0.04074, MSE: 0.04025
 [theta est ARX, theta est ARX sd] = getpvec(sys arx) %theta and standard
deviation of theta
  theta_est_ARX = 2 \times 1
     -0.9818
       0.3710
  theta_est_ARX_sd = 2 \times 1
       0.0021
       0.0384
 theta est ARX = abs(theta est ARX);
 % Parameter nominal values and confidence interval
 kv_est_ARX = (1 - theta_est_ARX(1)) / T;
                                                                    % point 1
estimation of kv using Euler approximation
 lbkv_ARX = (1 - theta_est_ARX(1) - theta_est_ARX_sd(1)*3) / T; % lower bound
for 98% confidence interval for kv
 ubkv_ARX = (1 - theta_est_ARX(1) + theta_est_ARX_sd(1)*3) / T; % upper bound
for 98% confidence interval for kv
 % Exact determination of confidence interval for kb_est is too difficult
```

We create a table to compare the results.

```
Comparison = ["kv";"lbkv";"ubkv";"kb";"lbkb";"ubkb"];
Non_recursive = [kv_est;lbkv;ubkv;kb_est;lbkb;ubkb];
ARX = [kv_est_ARX;lbkv_ARX;ubkv_ARX;kb_est_4;lbkb_ARX;ubkb_ARX];
results = table(Comparison,Non_recursive,ARX)
```

results =  $6 \times 3$  table

	Comparison	Non_recursive	ARX
1	"kv"	0.0908	0.0908
2	"lbkv"	-1.2598e+03	0.0600
3	"ubkv"	1.2600e+03	0.1215
4	"kb"	1.8548	1.8548
5	"lbkb"	-65.5036	1.2795
6	"ubkb"	69.2132	2.4301

We can see in the table that the values of  $k_{\nu}$  and  $k_b$  are the same but the confidence intervals is different, smaller for the arx model which means its more precise or confident.

## 2.c - Recursive method

Recursive method is designed using the following equations.

```
 \begin{pmatrix} y(2) \\ y(3) \end{pmatrix} = \begin{pmatrix} y(1) & u(1) \\ y(2) & u(2) \end{pmatrix} \cdot \begin{pmatrix} \hat{a} \\ \hat{b} \end{pmatrix}   P(3) = \begin{bmatrix} C(3)^T \cdot C(3) \end{bmatrix}^{-1}   \hat{\theta}(3) = P(3) \cdot C^T(3) \cdot Y(3)
```

```
c(k+1) = (y(k) \quad u(k))
P(k+1) = P(k) - \frac{P(k) \cdot c^{T}(k+1) \cdot c(k+1) \cdot P(k)}{1 + c(k+1) \cdot P(k) \cdot c^{T}(k+1)}
K(k+1) = \frac{P(k) \cdot c^{T}(k+1)}{1 + c(k+1) \cdot P(k) \cdot c^{T}(k+1)}
\hat{\theta}(k+1) = \hat{\theta}(k) + K(k+1) \cdot \left[ y(k+1) - c(k+1) \cdot \hat{\theta}(k) \right]
```

```
Y=[y(2);y(3)];
                        %Y(3)
C=[y(1:2) u(1:2)];
                        %C(3)
P=inv(C'*C);
                       %P(3)
theta_est_rec=P*C'*Y; %theta(3)
for i=3:N-1
c=[y(i) u(i)];
D=1+c*P*c'; %Denominator
 K=P*c'/D;
theta est rec=theta est rec+K*(y(i+1)-c*theta est rec);
 P=P-P*c'*c*P/D;
end
kv_est_rec=-log(theta_est_rec(1))/T % point estimation of kv
kv est rec = 0.0916
```

kb\_est\_rec=kv\_est\_rec\*theta\_est\_rec(2)/(1-exp(-kv\_est\_rec\*T)) % kb

kb est rec = 1.8718

```
clear lbkb ubkb lbkv ubkv kb_est kv_est;
lbkb=zeros(N,1);
ubkb=zeros(N,1);
lbkv=zeros(N,1);
ubkv=zeros(N,1);
kb_est=zeros(N,1);
kv_est=zeros(N,1);
for i=3:N-1
```

```
c=[y(i) u(i)];
    D=1+c*P*c'; %Denominator
    K=P*c'/D;
    theta_est_rec=theta_est_rec+K*(y(i+1)-c*theta_est_rec);
    P=P-P*c'*c*P/D;
    sigma2=J(theta_est_rec)/(i-2);  % estimation of noise variance
    Cov mat=sigma2*P;
                                   % covariance matrix
    std_error=sqrt(diag(Cov_mat)); % standard errors
                                                    % point estimation of kv
    kv est(i)=(1-theta est rec(1))/T;
    lbkv(i)=(1-theta_est_rec(1)-std_error(1)*3)/T; % lower bound for 98%
    ubkv(i)=(1-theta_est_rec(1)+std_error(1)*3)/T; % upper bound for 98%
    kb est(i)=theta est rec(2)/T;
                                                    % point estimation of kb
    lbkb(i)=(theta_est_rec(2)-std_error(2)*3)/T;
                                                    % lower bound for 98%
    ubkb(i)=(theta_est_rec(2)+std_error(2)*3)/T;
                                                    % upper bound for 98%
end
```

# Calculation using recursive arx

```
load data2 data; % Read data
u = data(:,1); % it extracts inputs
y = data(:,2); % it extracts outputs
N = size(data,1); % determine the number of points
T = 0.2; % Sample time
lbkb_ARX1=zeros(N,1);
ubkb_ARX1=zeros(N,1);
lbkv_ARX1=zeros(N,1);
ubkv_ARX1=zeros(N,1);
kv_est_ARX=zeros(N,1);
kb est ARX=zeros(N,1);
std_error_ARX = [];
estimator = recursiveARX([1 1 0]);
for k = 1:N
    [theta1,theta2,y_hat] = estimator(y(k),u(k));
    cov = estimator.ParameterCovariance; % covariance matrix
    theta ARX = [-theta1(2) theta2];
    std_error_ARX = diag(sqrt(cov));
    kv_est_ARX(k) =(1-theta_ARX(1))/T;
                                                            % Estimation of kv
```

```
kv est ARX1 = 0.0881
```

```
kb_est_ARX1=kv_est_ARX1*theta_ARX(2)/(1-exp(-kv_est_ARX1*T)) % kb
```

```
kb_est_ARX1 = 1.8056
```

Table for comparison.

```
Comparison = ["kv";"lbkv";"ubkv";"kb";"lbkb";"ubkb"];
recursive = [kv_est_rec;lbkv(N-2);ubkv(N-2);kb_est_rec;lbkb(N-2);ubkb(N-2)];
recARX = [kv_est_ARX1;lbkv_ARX1(N-2);ubkv_ARX1(N-2);kb_est_ARX1;lbkb_ARX1(N-2);ubkb_ARX1(N-2)];
results = table(Comparison, recursive, recARX)
```

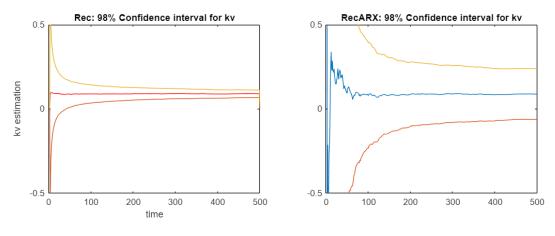
results =  $6 \times 3$  table

	Comparison	recursive	recARX
1	"kv"	0.0916	0.0881
2	"lbkv"	0.0672	-0.0629
3	"ubkv"	0.1115	0.2378
4	"kb"	1.8718	1.8056
5	"lbkb"	1.4119	-1.0180
6	"ubkb"	2.2413	4.5994

Comparing the recursive method and arx recursive method, we can see that the recursive ARX is now less precise then just recursive method, and the bound for recursive ARX are bigger.

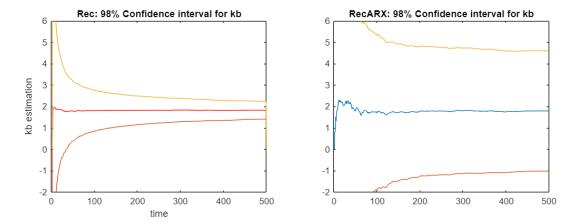
```
figure('Position',[0 0 1200 400])
subplot(1,2,1)
plot(kv_est,'r'), hold on;
plot(lbkv);
plot(ubkv);
axis([0 N -0.5 0.5]);
title('Rec: 98% Confidence interval for kv');
```

```
xlabel('time');
ylabel('kv estimation');
subplot(1,2,2)
plot(kv_est_ARX), hold on
plot(lbkv_ARX1)
plot(ubkv_ARX1)
title('RecARX: 98% Confidence interval for kv')
axis([0 N -0.5 0.5]);
```



Comparison of  $k_{\nu}$  estimated using recursive method and recursiveARX.

```
figure('Position',[0 0 1200 400])
subplot(1,2,1)
plot(kb_est,'r'), hold on;
plot(lbkb);
plot(ubkb);
axis([0 N -2 6]);
title('Rec: 98% Confidence interval for kb');
xlabel('time');
ylabel('kb estimation');
subplot(1,2,2)
plot(kb_est_ARX), hold on
plot(lbkb_ARX1)
plot(ubkb_ARX1)
title('RecARX: 98% Confidence interval for kb')
axis([0 N -2 6]);
```



Comparison of  $k_b$  estimated using recursive method and recursive ARX.

# 2.d - Using Kalman filter as an estimator

sigma2 = 0.0404

We designed our Kalman filter using the equations in the slides.

The parameters are considered as the states  $x(k)=\theta(k)$ 

```
Then, regressor model y(k)=c^T(k)\theta in state space form:  \theta(k+1)=\theta(k)   y(k)=c^T(k)\theta+v(k)
```

v(k) measurement noise with variance R

```
Considering state space form of y(k)=c^T(k)\theta
A=I
B=0
C=c^T
and no disturbances: Q=0

Then, Kalman filter formulas for parameter estimation are:
L(k) = P(k)c(k) \left[R + c^T(k)P(k)c(k)\right]^{-1}
\hat{\theta}(k+1) = \hat{\theta}(k) + L(k) \left[y(k) - c^T(k)\hat{\theta}(k)\right]
P(k+1) = \left[I - L(k)c^T(k)\right]P(k)
```

Considering that the output y(k) is y(k+1).

```
for i=3:N-1
  c=[y(i) u(i)]';
  L=P*c/(sigma2+c'*P*c);
  theta_est_kal=theta_est_kal+L*(y(i+1)-transpose(c)*theta_est_kal);
  P=(eye(2)-L*transpose(c))*P;
  Cov mat=sigma2*P;
                          % covariance matrix
  kv_est_kal(i)=(1-theta_est_kal(1))/T;  % point estimation of kv
  lbkv kal(i)=(1-theta est kal(1)-std error(1)*3)/T; % lower bound for 98%
  ubkv_kal(i)=(1-theta_est_kal(1)+std_error(1)*3)/T; % upper bound for 98%
  kb est kal(i)=theta est kal(2)/T;
                            % point estimation of kb
  kv_est_Kal1=-log(theta_est_kal(1))/T
                          % point estimation of kv
```

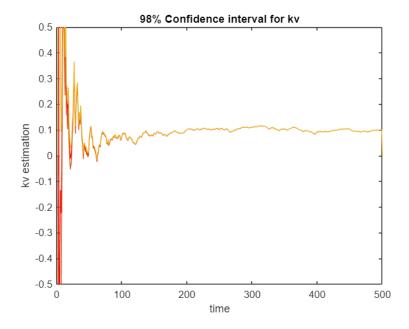
 $kv_est_Kal1 = 0.1006$ 

```
kb_est_Kal1=kv_est_Kal1*theta_est_kal(2)/(1-exp(-kv_est_Kal1*T)) % kb
```

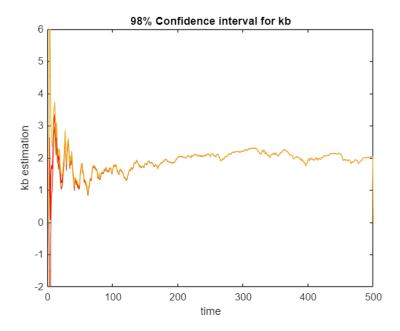
```
kb_est_Kal1 = 2.0596
```

The results using Kalman are very precise and similar to the initial parameters of  $k_{\nu} = 0.1$  and  $k_b = 2$ . This method looks to be the best for parameter estimation.

```
figure
plot(kv_est_kal,'r');
hold on;
plot(lbkv_kal);
plot(ubkv_kal);
axis([0 N -0.5 0.5]);
title('98% Confidence interval for kv');
xlabel('time');
ylabel('kv estimation');
```



```
figure
plot(kb_est_kal,'r');
hold on;
plot(lbkb_kal);
plot(ubkb_kal);
axis([0 N -2 6]);
title('98% Confidence interval for kb');
xlabel('time');
ylabel('kb estimation');
```



On both graphs we can see that the confidence interval is really small which means that the estimated value is very close to the real value.