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```
% Repeat Problem 3, except use the problem matrices and measurement data
% that are defined by the Matlab script kf example02b.m. Notice that the R
% and Q values are different for this problem and that there is a different
% measurement time history.

% Run your Kalman filter two additional times using the two alternate Q
% values that are mentioned in the comments in the file kf example02b.m.
% It is uncertain which is the correct Q value.

clear; clc;
close all;

kf_example02b

% Qk_a = Qk
% Qk_b = first alternate Qk
% Qk_c = second alternate Qk
Gk = Gammak;
```

KALMAN FILTER

Decide which is the best value in the following way: Calculate err(v(k))

for k = 1, 2, ..., 50 for each of your runs. Compute the average of these 50 values. This average times 50, i.e., {err_v(1) + err_v(2) + ... + err_v(50)}, will be a sample of chi-square distribution of degree 50 if the filter model is correct. Develop upper and lower limits between which the average {err_v(1) + err_v(2) + ... + err_v(50)}/50 must lie 99% of the time if the Kalman filter model is correct, and test your averages for each of the three candidate Q values. Which is the most reasonable? Look at the state estimate differences between the best filter and the other two filters.

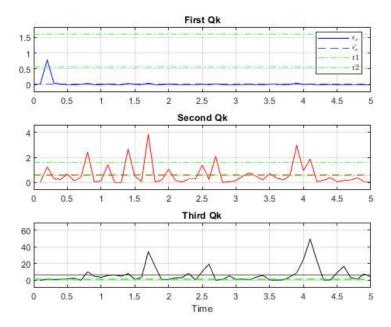
```
% err_nu_k = 1/2 * [ nu_k ]' * inv(S_k) * nu_k
for i = 1:length(nu a)
    err_nu_a(i) = nu_a(i) * inv(S_a(i)) * nu_a(i);
    err_nu_b(i) = nu_b(i) * inv(S_b(i)) * nu_b(i);
    err_nu_c(i) = nu_c(i) * inv(S_c(i)) * nu_c(i);
err_nu_a_mean = mean(err_nu_a);
err_nu_b_mean = mean(err_nu_b);
err_nu_c_mean = mean(err_nu_c);
N = 50;
Nx = 2;
Nz = 1;
% NEED STATISTICS TOOLBOX
a = .01;
r1 = chi2inv(a/2, N*Nz) / N;
r2 = chi2inv(1 - a/2, N * Nz) / N;
figure
```

```
subplot(3,1,1)
        plot( thist, err_nu_a, 'b' ); hold on; grid on;
        yline(err_nu_a_mean, 'b--');
        yline(r1, 'g-.');
        yline(r2, 'g-.');
        bigger_ylim
        title('First Qk');
        \label{legend('$\epsilon_{\nu}}', '$\bar{\epsilon_{\nu}}$', 'r1', 'r2', 'interpreter', 'latex')
        plot( thist, err_nu_b, 'r');
        yline(err_nu_b_mean, 'r--');
        yline(r1, 'g-.');
yline(r2, 'g-.');
        bigger_ylim
        title('Second Qk');
    subplot(3,1,3)
        plot( thist, err_nu_c, 'k');
        yline(err_nu_c_mean, 'k');
        yline(r1, 'g--');
yline(r2, 'g--');
        title('Third Qk')
        bigger_ylim
    xlabel('Time')
if err_nu_a_mean > r1 && err_nu_a_mean < r2</pre>
    sprintf('Qk = %g is reasonable', Qk_a)
else
    sprintf('Qk = %g is NOT reasonable', Qk_a)
end
if err_nu_b_mean > r1 && err_nu_b_mean < r2</pre>
    sprintf('Qk = %g is reasonable', Qk_b)
    sprintf('Qk = %g is NOT reasonable', Qk_b)
end
if err_nu_c_mean > r1 && err_nu_c_mean < r2</pre>
    sprintf('Qk = %g is reasonable', Qk_c)
    sprintf('Qk = %g is NOT reasonable', Qk_c)
% possibly use Nz? Otherwise Q is not within bounds
```

```
ans =
  'Qk = 40 is NOT reasonable'

ans =
  'Qk = 0.4 is reasonable'

ans =
  'Qk = 0.004 is NOT reasonable'
```



Compute the RMS value of the difference time history

for each state vector element. Do the averaging over the last 40 points.

```
x1_rms_ab = rms( xhat_a(11:end,1) - xhat_b(11:end,1) );
x2_rms_ab = rms( xhat_a(11:end,2) - xhat_b(11:end,2) );

x1_rms_bc = rms( xhat_b(11:end,1) - xhat_c(11:end,1) );
x2_rms_bc = rms( xhat_b(11:end,2) - xhat_c(11:end,2) );

% Are these differences significant compared to the computed state
% estimation error standard deviations for the best filter?
x1_rms_ab / P11_b(end)
x2_rms_ab / P22_b(end)

x1_rms_bc / P11_b(end)
x2_rms_bc / P22_b(end)

disp('The second filter is the best filter. The first filter is closer to the second filter than the third filter is to the second filter');
```

```
ans = 68.0758727260931

ans = 15.8780431653489

ans = 1211.25919352871

ans = 86.8008602066569
```

The second filter is the best filter. The first filter is closer to the second filter than the third filter is to the second filter

subfunctions KALMAN FILTER

```
xhat = xhat0;
P = P0;
% Initialize saved output arrays
xbar_arr = [xhat'];
Pbar_cell = {P};
xhat_arr = [xhat'];
P_cell = {P};
Pxx_arr = [P(1,1)];
Pzz\_arr = [P(2,2)];
nu_arr = [];
S_arr = [];
% Propagate and filter through all measurements
for k = 0: length(zhist)-1
    % propagate state and covar
    xbar = Fk * xhat;
                                              % a priori state est
    Pbar = Fk * P * Fk' + Gk * Qk * Gk'; % a priori covar est
    % update
    nu = zhist(k+1) - Hk * xbar;
                                              % innovation
    S = Hk * Pbar * Hk' + Rk;
                                               % innovation covariance
    W = Pbar * Hk' * inv(S);
                                               % Kalman gain
    xhat = xbar + W * nu;
                                              % a posteriori state est
    P = Pbar - W * S * W';
                                              % a posteriori covar est
    % next step
    k = k + 1;
    \% save states and covariances
    xbar_arr = [xbar_arr; xbar'];
    Pbar_cell = {Pbar_cell; Pbar};
    xhat_arr = [xhat_arr; xhat'];
    P_{cell} = \{P_{cell}; P\};
    Pxx_arr = [Pxx_arr; P(1,1)];
Pzz_arr = [Pzz_arr; P(2,2)];
nu_arr = [nu_arr; nu];
              = [S_arr; S];
    S_arr
end
```

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