

LECTURE 2

* Kalman Filter

The recursive LS formulation from last time was:

$$\begin{aligned}
 P_i^{-1} &= P_{i-1}^{-1} + A_i^T R_i^{-1} A_i \\
 K_i &= P_i A_i^T R_i^{-1} \\
 \hat{X}_i &= \hat{X}_{i-1} + K_i (Y_i - A_i \hat{X}_{i-1})
 \end{aligned}$$

→ The fundamental assumption is that X is constant and that its estimation \hat{X} must be updated as new data points arrive.

Let's assume now that X can actually evolve in time:

$$\begin{aligned}
 X_i &= F_{i-1} X_{i-1} + \epsilon_{i-1} & X_i &= F_{i-1} X_{i-1} + G_{i-1} u_{i-1} + \epsilon_{i-1} \\
 Y_i &= A_i X_i + \eta_i & Y_i &= A_i X_i + \eta_i
 \end{aligned}$$

or

$\epsilon_i \triangleq$ noise in the model
 $\eta_i \triangleq$ noise in the measurements (\Rightarrow it's the " ϵ " in lecture 1)
 $u_i \triangleq$ control vector (if any)

Let's also assume:

$$\begin{aligned}
 \epsilon_i &\sim N(0, R_i) \quad \swarrow \text{Gaussian} \\
 \eta_i &\sim N(0, R_i) \quad \swarrow \text{noise}
 \end{aligned}$$

and F_i, G_i, A_i known for every i

Because of ϵ_i, η_i , X_i is a random variable \Rightarrow Our estimation will be on the mean and we can write:

$$\begin{aligned}
 E(X_i) &= E(F_{i-1} X_{i-1} + G_{i-1} u_{i-1} + \epsilon_{i-1}) \\
 &= F_{i-1} E(X_{i-1}) + G_{i-1} u_{i-1}
 \end{aligned}$$

\Rightarrow The mean of X_i is propagated by the formula:

②

$$\bar{x}_i = F_{i-1} \bar{x}_{i-1} + G_{i-1} u_{i-1} \quad \text{where } \bar{x}_i \triangleq E(X_i) \quad (1)$$

$$E[(X_i - \bar{x}_i)(X_i - \bar{x}_i)^T] =$$

$$= E\left[\left(F_{i-1} X_{i-1} + G_{i-1} u_{i-1} + \varepsilon_{i-1} - F_{i-1} \bar{x}_{i-1} - G_{i-1} u_{i-1}\right) \left(\dots\right)^T\right]$$

$$= E\left[\left(F_{i-1} (X_{i-1} - \bar{x}_{i-1}) + \varepsilon_{i-1}\right) \left(F_{i-1} (X_{i-1} - \bar{x}_{i-1}) + \varepsilon_{i-1}\right)^T\right]$$

$$= F_{i-1} E\left[(X_{i-1} - \bar{x}_{i-1})(X_{i-1} - \bar{x}_{i-1})^T\right] F_{i-1}^T + E\left[\varepsilon_{i-1} \varepsilon_{i-1}^T\right] \xrightarrow{\triangleq Q_{i-1}}$$

$$+ E\left[\varepsilon_{i-1} \left(F_{i-1} (X_{i-1} - \bar{x}_{i-1})\right)^T\right] + E\left[(X_{i-1} - \bar{x}_{i-1})^T F_{i-1}^T \varepsilon_{i-1}\right]$$

\downarrow
 $= 0$

\downarrow
 $= 0$

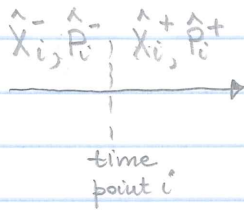
By definition,
the noise is
uncorrelated with
the variable X_i

Because $E[(X_i - \bar{x}_i)(X_i - \bar{x}_i)^T] = P_i$ - covariance of the estimation,

we can write:

$$P_i = F_{i-1} P_{i-1} F_{i-1}^T + Q_{i-1} \quad (2)$$

Note: In eq. (1) and (2), we have just used the evolution model of X_i , not the estimations $\hat{X}_i, \hat{P}_i \Rightarrow$ When it comes about estimation and we are at time i , two options are viable:



a) We use data collected up to i , EXCLUDED the data collected at time $i \Rightarrow$ Let's use the notation: \hat{X}_i^- , \hat{P}_i^-

b) We use data collected up to i , INCLUDED the data collected at time $i \Rightarrow$ Let's use the notation: \hat{X}_i^+ , \hat{P}_i^+

Therefore, we can have a two-tiered approach:

- If we are waiting for the data to be collected at time i , we can enhance our estimation from time $i-1$ by using the model:

$$(*) \quad \begin{aligned} \hat{X}_i^- &= F_{i-1} \hat{X}_{i-1}^+ + G_{i-1} u_{i-1} \\ \hat{P}_i^- &= F_{i-1} \hat{P}_{i-1}^+ F_{i-1}^T + Q_{i-1} \end{aligned}$$

- If we have collected the data at time i , we can refine the estimation by using the recursive Least square formula:

$$\hat{P}_i^+ = \left[(\hat{P}_i^-)^{-1} + A_i^T R_i^{-1} A_i \right]^{-1}$$

$$(**) \quad K_i = \hat{P}_i^+ A_i^T R_i^{-1}$$

$$\hat{X}_i^+ = \hat{X}_i^- + K_i (Y_i - A_i \hat{X}_i^-)$$

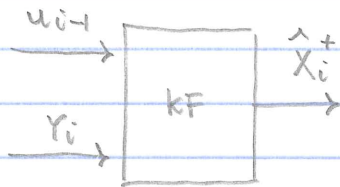
(*) is a prediction step and (**) is a correction step \Rightarrow The combination is the Kalman filter

Note a few things:

$$\hat{X}_i^+ = F_{i-1} \hat{X}_{i-1}^+ + G_{i-1} u_{i-1} + K_i Y_i - K_i A_i F_{i-1} \hat{X}_{i-1}^+ - K_i A_i G_{i-1} u_{i-1}$$

④

$$= (F_{i-1} - K_i A_i F_{i-1}) \hat{X}_{i-1}^+ + (G_{i-1} - K_i A_i G_{i-1}) u_{i-1} + K_i Y_i$$



\Rightarrow kF is actually a time-varying filter that operates on the input variables u and Y

In (*), we provide an estimation for X in the absence of new data \Rightarrow If no further data were provided, our estimation would evolve simply based on the model \Rightarrow Uncertainty would grow in time because of the term Q_{i-1}
 \Rightarrow Estimation worsens between measures

In (**), a large \hat{P}_i^- is overshadowed by $A_i^T R_i^{-1} A_i$, since it contributes as $(\hat{P}_i^-)^{-1} \Rightarrow$ A small uncertainty on the measurements (i.e., R_i small) will likely make \hat{P}_i^+ small, i.e.:

$$P_i^+ \cong (A_i^T R_i^{-1} A_i)^{-1}$$

However, in $K_i = P_i^+ A_i^T R_i^{-1}$, a small P_i^+ is bad for updating the estimation $\hat{X}_i^+ \Rightarrow$ That's why the presence of R_i^{-1} compensates for P_i^+ and makes K_i sizable enough to use Y_i in the update of \hat{X}_i^+

Moreover, note: R_i^{-1} large $\Rightarrow A_i^T R_i^{-1} A_i$ - small $\Rightarrow P_i^+ \cong P_i^-$ and K_i - small \Rightarrow The uncertainty of the estimation is let evolve as in the absence of new measurements and new measurements are trusted very little when it comes about updating the estimate of X

Finally note: while a small R_i is generally useful (i.e., we can trust the measurements), a small Q_i may eventually lead to undesired conditions:

Q_i - small $\Rightarrow P_i^-$ small $\Rightarrow P_i^+ \cong P_i^-$ small $\Rightarrow K_i$ small, i.e., if we

trust the model too much, the measurements will be eventually ignored. \square

* Kalman Filter for Nonlinear Models

$$\left. \begin{aligned} X_i &= F_{i-1} X_{i-1} + G_{i-1} u_{i-1} + \varepsilon_{i-1} \\ Y_i &= A_i X_i + \eta_i \end{aligned} \right\} \begin{array}{l} \text{If the model is accurate, then the} \\ \text{error } (Y_i - A_i \hat{X}_i^-) \rightarrow \text{Gaussian RV} \end{array}$$



The KF is able to extract all the available information buried in noisy data

What if the model is nonlinear?

$$\begin{aligned} X_i &= f_i(X_{i-1}, u_{i-1}) + \varepsilon_i & \varepsilon_i &\sim N(0, Q_i) \\ Y_i &= g_i(X_i) + \eta_i & \eta_i &\sim N(0, R_i) \end{aligned}$$

EKF) One approach focuses on the variations of X_i from its expected value μ_i and uses the first order Taylor series expansion:

$$f_i(X_{i-1}, u_{i-1}) \cong f_i(\mu_{i-1}, u_{i-1}) + \frac{df_i}{dX}(\mu_{i-1}, u_{i-1})(X_{i-1} - \mu_{i-1})$$

$$g_i(X_i) \cong g_i(\mu_i) + \frac{dg_i}{dX}(\mu_i)(X_i - \mu_i)$$



$$\mu_i = E(X_i) = E(f_i(X_{i-1}, u_{i-1})) \cong f_i(\mu_{i-1}, u_{i-1})$$

$$\text{Define: } Z_i \triangleq X_i - \mu_i; \quad V_i \triangleq Y_i - g_i(\mu_i); \quad F_{i-1} \triangleq \frac{df_i}{dX}(\mu_{i-1}, u_{i-1}); \quad A_i \triangleq \frac{dg_i}{dX}(\mu_i)$$

and consider the linearized model:
$$\begin{cases} Z_i = F_{i-1} Z_{i-1} + \varepsilon_{i-1} \\ V_i = A_i Z_i + \eta_i \end{cases} \quad (1)$$

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Now, in order to estimate X_i , one can observe:

Our estimation focuses on the expected value of X_i (so it was in the linear case)

Denoted with \hat{X}_{i-1}^+ the best estimate of μ_{i-1} after collecting the data at time $i-1$, we can write:

$$\hat{X}_i^- = f_i(\hat{X}_{i-1}^+, u_{i-1})$$

$$F_{i-1} = \frac{df_i}{dx}(\hat{X}_{i-1}^+, u_{i-1})$$

Also note: $P_i \triangleq E((X_i - \mu_i)(X_i - \mu_i)^T) = E(\underbrace{z_i z_i^T}_{\text{def of } z_i}) \Rightarrow$ We can apply the KF

to (1) and estimate P_i :

$$\hat{P}_i^- = F_{i-1} \hat{P}_{i-1}^+ F_{i-1}^T + Q_{i-1}$$

$$\hat{X}_i^- = f_i(\hat{X}_{i-1}^+, u_{i-1})$$

(E1)

where: \hat{P}_{i-1}^+ is the best estimate of P_{i-1} after collecting the data at time $i-1$

$$F_{i-1} = \frac{df_i}{dx}(\hat{X}_{i-1}^+, u_{i-1})$$

Similarly note:

$$\hat{z}_i^+ = \hat{z}_i^- + k_i (y_i - A_i \hat{z}_i^-) \Rightarrow \hat{X}_i^+ = \hat{X}_i^- + k_i (y_i - \underbrace{g_i(\mu_i)}_{\cong g_i(\hat{X}_i^-)}) - A_i (\hat{X}_i^- - \mu_i)$$

Taylor series expansion

$$\Rightarrow \hat{X}_i^+ = \hat{X}_i^- + k_i (y_i - g_i(\hat{X}_i^-))$$

Remember: we need to update the estimate of P_i in order to compute k_i :

$$\hat{P}_i^+ = E((\hat{X}_i^+ - \mu_i)(\hat{X}_i^+ - \mu_i)^T) = E(\hat{z}_i^+ \hat{z}_i^{+T}) \Rightarrow$$
 By using KF on (1):

$$\hat{P}_i^+ = \left[(\hat{P}_i^-)^{-1} + A_i^T R_i^{-1} A_i \right]^{-1}$$

$$K_i = \hat{P}_i^+ A_i^T R_i^{-1}$$

$$\hat{X}_i^+ = \hat{X}_i^- + K_i (Y_i - g_i(\hat{X}_i^-))$$

(E2) where:

- \hat{X}_i^- is the best estimate of X_i before collecting the data at time i
- $A_i = \frac{dg_i}{dX}(\hat{X}_i^-)$

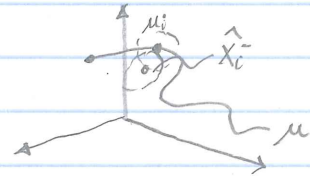
The combination (E1)-(E2) is called "Extended Kalman Filter" (EKF)

Problems:

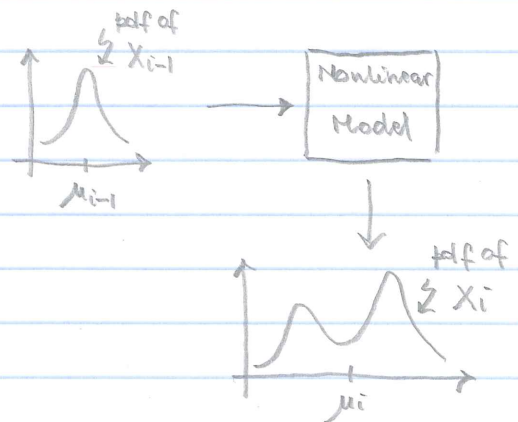
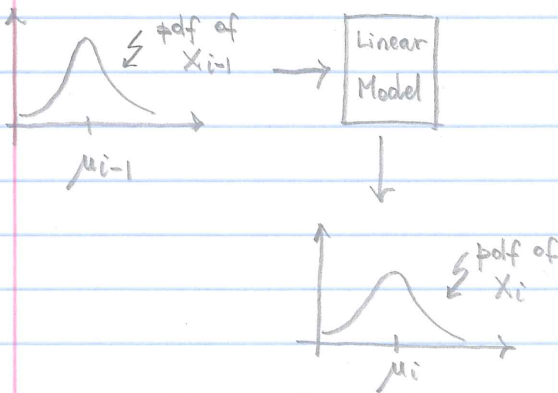
- F_{i-1} and A_i are obtained by linearizing around the estimated \hat{X}_{i-1}^+ and \hat{X}_i^- , respectively, not the actual mean values μ_{i-1} , μ_i , respectively \Rightarrow If the estimates are bad, the filter may drift away



EKF can give acceptable results as long as the nonlinearity is weak (i.e., the linearization introduces a small error)

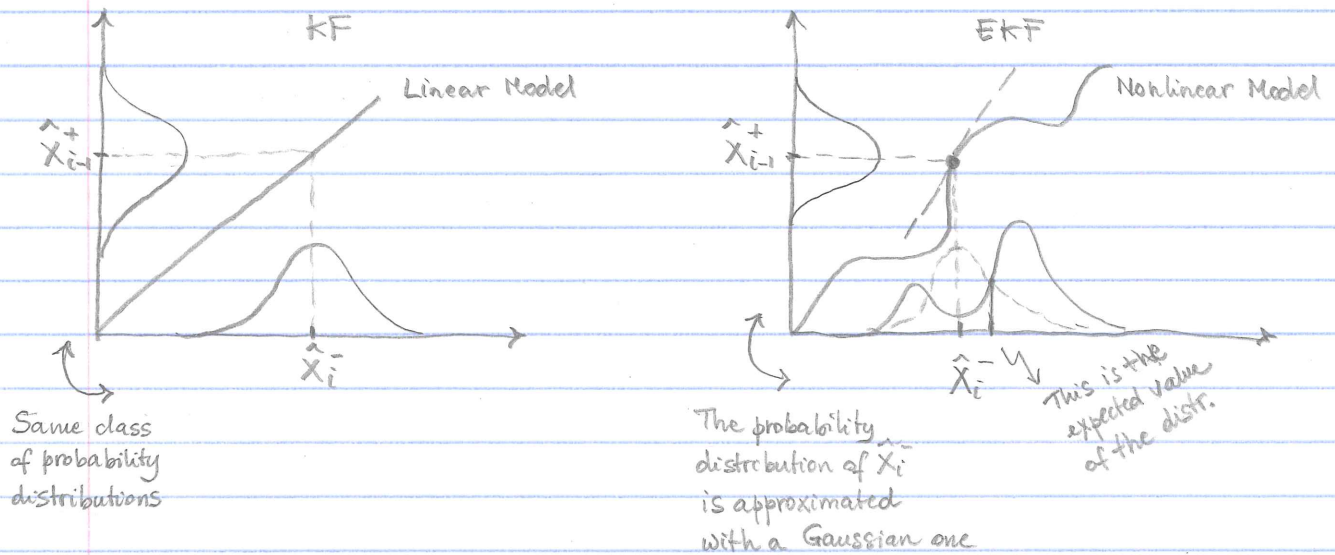


- F_{i-1} and A_i are jacobian matrices \Rightarrow Their calculation can be challenging and numerical issues can rise
- By linearizing, we assume that the distribution of X propagates from time $i-1$ to time i according to a linear model \Rightarrow It is not always the case:



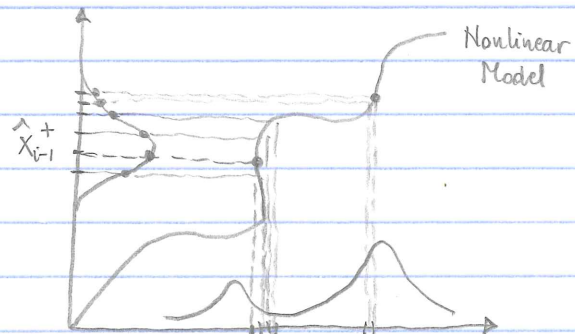
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Now, when we propagate our estimation by using KF (linear case) and EKF (nonlinear case), we have:



With EKF, \hat{X}_i^- may be a very bad estimation of $\mu_i \Rightarrow$ We need a different approach that aims to sample the distributions of \hat{X}_{i-1}^+ , \hat{X}_i^- rather than just assume they are Gaussian.

UKF) First, let's assume that \hat{X}_{i-1}^+ is Gaussian and let's propagate many values drawn from this distribution through the nonlinear model



Second, as the number of values (a.k.a. "particles") increases, a better estimation of the distribution of \hat{X}_i^- can be given

We don't need to follow a Monte Carlo approach, though. It can be proved that - if \hat{X}_{i-1}^+ is Gaussian and n is the size of X - we only need $n+1$ particles

to estimate mean and variance of distribution determined by the nonlinear transformation of \hat{X}_{i-1}^+ \Rightarrow We have:

\hat{X}_{i-1}^+ is the mean of the input distribution
 \hat{P}_{i-1}^+ is the covariance matrix of the input distribution \Rightarrow It is a $n \times n$ matrix with columns $p_{i-1,1}^+, p_{i-1,2}^+, \dots, p_{i-1,n}^+$

\Rightarrow The $n+1$ points must be:

$\sigma_0 = \hat{X}_{i-1}^+$
 $\sigma_j = \sigma_0 + \sqrt{n} p_{i-1,j}^+ \quad j=1,2,\dots,n$
 $\sigma_{n+j} = \sigma_0 - \sqrt{n} p_{i-1,j}^+ \quad j=1,2,\dots,n$

\Rightarrow The transformation of these $2n+1$ points is:

$\gamma_0 = f_i(\sigma_0, u_{i-1})$
 $\gamma_j = f_i(\sigma_j, u_{i-1}) \quad j=1,2,\dots,2n$

\Rightarrow Sample mean and covariance of this distribution can be used for prediction:

$$\hat{X}_i^- = \frac{1}{2n+1} \sum_{j=0}^{2n} \gamma_j$$

$$\hat{P}_i^- = \text{cov}(\{\gamma_j\}_{j=0,2n}) + Q_{i-1}$$

$$= \frac{1}{2n+1} \sum_{j=0}^{2n} (\gamma_j - \hat{X}_i^-)(\gamma_j - \hat{X}_i^-)^T + Q_{i-1}$$

Analogously, one can compute:

$\xi_0 = g_i(\sigma_0)$
 $\xi_j = g_i(\sigma_j) \quad j=1,2,\dots,2n$

\Rightarrow One can estimate sample mean and covariance:

$$\tilde{y}_i = \frac{1}{2n+1} \sum_{j=0}^{2n} \xi_j$$

$$P_{yy} = \frac{1}{2n+1} \sum_{j=0}^{2n} (\xi_j - \tilde{y}_i)(\xi_j - \tilde{y}_i)^T + R_i$$

(10)

Note that we add Q_{i-1} and R_i to \hat{P}_i^- and P_{yy} , respectively, because we want to compute an estimation of the covariances of samples of X and Y , respectively, but the samples have been chosen in a deterministic way

With $\{\chi_j\}$ and $\{\xi_j\}$, one can also compute:

$$P_{xy} = \frac{1}{2n+1} \sum_{j=0}^{2n} (\chi_j - \hat{X}_i^-) (\xi_j - \tilde{Y}_i)^T$$

If X and Y have size n and m , respectively, this is a $n \times m$ matrix

It can be shown that the value:

$$\hat{X}_i^+ = \hat{X}_i^- + k_i (Y_i - \tilde{Y}_i) \quad \text{with } k_i = P_{xy} P_{yy}^{-1}$$

is the unbiased estimation of X_i that minimizes the sum of the variances of the individual components of $X_i \Rightarrow$ Therefore, the following filter is derived:

Sampling Step:

$$\sigma_0 = \hat{X}_{i-1}^+; \quad \sigma_j = \hat{X}_{i-1}^+ + \sqrt{n [\hat{P}_{i-1}^+]_j}; \quad \sigma_{n+j} = \hat{X}_{i-1}^+ - \sqrt{n [\hat{P}_{i-1}^+]_j} \quad j=1, 2, \dots, n$$

(U1) with $[\hat{P}_{i-1}^+]_j \triangleq p_{i-1,j}^+$ j -th column of \hat{P}_{i-1}^+

$$\chi_j = f_i(\sigma_j, u_{i-1}) \quad j=0, 1, \dots, 2n$$

$$\xi_j = g_i(\sigma_j) \quad j=0, 1, \dots, 2n$$

Prediction Step:

$$(u2) \quad \hat{X}_i^- = \frac{1}{2n+1} \sum_{j=0}^{2n} \chi_j$$

$$\hat{P}_i^- = \frac{1}{2n+1} \sum_{j=0}^{2n} (\chi_j - \hat{X}_i^-) (\chi_j - \hat{X}_i^-)^T + Q_{i-1}$$

Correction Step:

$$(U3) \quad \tilde{y}_i = \frac{1}{2n+1} \sum_{j=0}^{2n} \xi_j \quad P_{yy} = \frac{1}{2n+1} \sum_{j=0}^{2n} (\xi_j - \tilde{y}_i)(\xi_j - \tilde{y}_i)^T + R_i$$

$$P_{xy} = \frac{1}{2n+1} \sum_{j=0}^{2n} (\xi_j - \hat{x}_i^-)(\xi_j - \tilde{y}_i)^T$$

$$\hat{P}_i^+ = \hat{P}_i^- - P_{xy} P_{yy}^{-1} P_{xy}^T$$

$$\hat{x}_i^+ = \hat{x}_i^- + P_{xy} P_{yy}^{-1} (y_i - \tilde{y}_i)$$

The combination of steps (U1) - (U3) is the "Unscented Kalman Filter" (UKF) \square

Slides present examples of application of Kalman filter to neural decoding of kinematic variables. \square

