

# OPTIMAL FILTERING

## LECTURE 6

- Smoothing
  - Using an extended state vector
  - General smoothing formulas
- Introduction to non-linear filtering



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**Reading instructions:** Kailath, Sect. 10.1-10.4.

Arulampalam, S., Maskell, S., Gordon, N. and Clapp, T. (2002)  
“A tutorial on particle filters for on-line nonlinear/non-Gaussian  
Bayesian tracking”. IEEE Trans. SP, 50(2):174–188.

F. Gustafsson, “Particle filter theory and practice with  
positioning applications,” IEEE A&E Syst. Mag. Part II:  
Tutorials, vol. 25, no. 7, pp. 53–82, 2010.

## SMOOTHING

Often, an estimation delay is acceptable, or off-line processing can be  
done  $\implies$  More (future) data can be used



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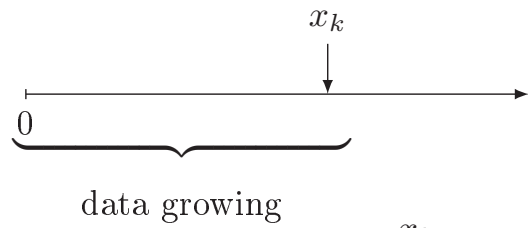
 More accurate estimates

 Increased computational complexity

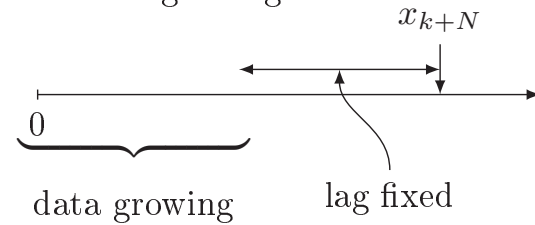
How to extend Kalman?

# DIFFERENT FORMS OF SMOOTHING

**Filtering**  $\hat{x}_{k|k}$

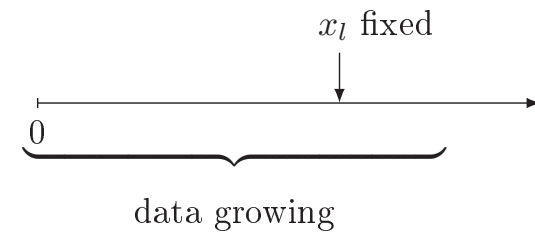


**Prediction**  $\hat{x}_{k+N|k}$



**Fixed point smoothing**

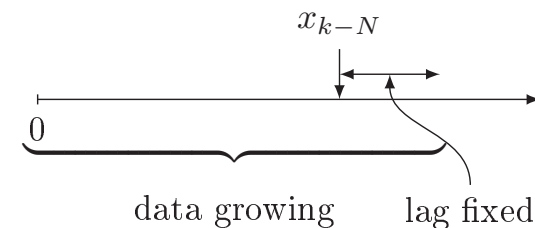
$\hat{x}_{l|k}$ ,  $l$  fixed,  $k$  growing



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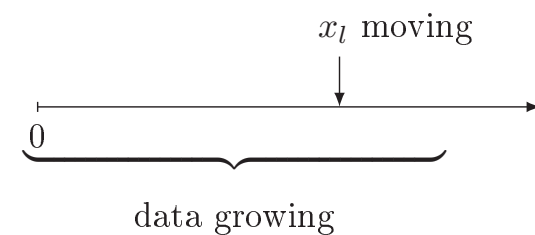
# DIFFERENT FORMS OF SMOOTHING, CONT.

**Fixed lag smoothing**  $\hat{x}_{k-N|k}$ ,



**Fixed interval smoothing**

$\hat{x}_{l|k}$ ,  $k = M$ ,  $0 \leq l \leq M$



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## APPLICATIONS



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**Fixed point smoothing:** Initial condition estimation

**Fixed lag smoothing:** on-line processing with delay

**Fixed interval smoothing:** off-line processing

...

## MULTIPLE-STEP AHEAD PREDICTOR



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$$\hat{x}_{k+N|k} = F_{k+N-1}F_{k+N-2} \cdots F_k \hat{x}_{k|k}$$

Show!

## FIXED POINT SMOOTHING

**Problem:** For all  $k > j$ ,  $j$  fixed, find  $\hat{x}_{j|k}$  and

$$P_k^{aa} = E[(x_j - \hat{x}_{j|k})(x_j - \hat{x}_{j|k})^*].$$

**Solution:** (assuming  $S_k = 0$ )

$$K_k^a = P_k^a H_k^* (H_k P_{k|k-1} H_k^* + R_k)^{-1}$$

$$\hat{x}_{j|k} = \hat{x}_{j|k-1} + K_k^a (y_k - H_k \hat{x}_{k|k-1})$$

$$P_{k+1}^a = P_k^a F_k^* - K_k^a H_k P_{k|k-1} F_k^*$$

$$P_{k+1}^{aa} = P_k^{aa} - K_k^a H_k P_k^a$$

**Initial conditions:**  $\hat{x}_{j|j-1} = \hat{x}_{j|j-1}$ ,  $P_j^{aa} = P_j^a = P_{j|j-1}$ .

**Improvement:**  $P_k^{aa} - P_{j|j-1} = \sum_{i=j}^{k-1} K_i^a H_i P_i^a$



## FIXED POINT SMOOTHING, COMMENTS

- What happens when  $R \rightarrow \infty$ ?
- Complexity?
- Most improvement obtained after 2-3 time constants of the system.
- Time varying even for time invariant systems!



## FIXED LAG SMOOTHING

**Problem:**  $N$  fixed,  $k$  growing. Find  $\hat{x}_{k-N|k}$  and

$$P_{k-N|k} = E[(x_{k-N} - \hat{x}_{k-N|k})(x_{k-N} - \hat{x}_{k-N|k})^*].$$

**Solution:** For  $i = 1, 2, \dots, N$

$$K_k^{(i+1)} = P_{k|k-1}^{(i)} H_k^* (H_k P_{k|k-1} H_k^* + R_k)^{-1}$$

$$\hat{x}_{k-i|k} = \hat{x}_{k-i|k-1} + K_k^{(i+1)} e_k$$

$$P_{k+1|k}^{(i+1)} = P_{k|k-1}^{(i)} (F_k - K_k H_k)^*$$

$$P_{k-i|k} = P_{k-i|k-1} - P_{k|k-1}^{(i)} H_k^* \left( K_k^{(i+1)} \right)^*$$

**Initialization:**  $\hat{x}_{0|-1} = 0$ ,  $\hat{x}_{-i|-1} = 0$ ,  $P_{k|k-1}^{(0)} = P_{k|k-1}$ .

- Time invariant for time invariant systems!
- Complexity?



## FIXED INTERVAL SMOOTHING

**Problem:**  $M$  fixed,  $0 \leq k \leq M$ , find  $\hat{x}_{k|M}$  and

$$P_{k|M} = E[(x_k - \hat{x}_{k|M})(x_k - \hat{x}_{k|M})^*].$$

**Solution 1:** Fixed lag smoother with fixed lag  $N = M$  and  $k = M$ .

**Solution 2:** Ordinary Kalman,  $0 \leq k \leq M$ , save  $\hat{x}_{j|j-1}$ ,  $\hat{x}_{j|j}$ ,  $P_{j|j-1}$ ,  $P_{j|j}$ , followed by a backward recursion  $j = M, M-1, \dots, 1$ :

$$\hat{x}_{j-1|M} = \hat{x}_{j-1|j-1} + P_{j-1|j-1} F_{j-1}^* P_{j|j-1}^{-1} (\hat{x}_{j|M} - \hat{x}_{j|j-1})$$

$$A_{j-1} = P_{j-1|j-1} F_{j-1}^* P_{j|j-1}^{-1}$$

$$P_{j-1|M} = P_{j-1|j-1} + A_{j-1} (P_{j|M} - P_{j|j-1}) A_{j-1}^*$$

- 2<sup>nd</sup> approach requires less work
- Normally, fixed lag smoother with sufficiently large lag is adequate.



# GENERAL SMOOTHING FORMULAS

“Bryson-Frazier” formulas (1963)  $0 \leq k \leq N$

$$F_{p,k} = F_k - K_k H_k$$

$$\lambda_{k|N} = F_{p,k}^* \lambda_{k+1|N} + H_k^* R_{e_k}^{-1} e_k, \quad \lambda_{N+1|N} = 0$$

$$\hat{x}_{k|N} = \hat{x}_{k|k-1} + P_{k|k-1} \lambda_{k|N}$$

$$\Lambda_{k|N} = F_{p,k}^* \Lambda_{k+1|N} F_{p,k} + H_k^* R_{e_k}^{-1} H_k, \quad \Lambda_{N+1|N} = 0$$

$$P_{k|N} = P_{k|k-1} - P_{k|k-1} \Lambda_{k|N} P_{k|k-1}$$



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## Two-pass algorithm

- (1) Forward  $\hat{x}_{k|k-1}, P_{k|k-1}$
- (2) Backward  $\lambda_{k|N}, \Lambda_{k|N}$
- (3) Combine to obtain  $\hat{x}_{k|N}, P_{k|N}$

**Filtered version** (using  $\hat{x}_{k|k}$  and  $P_{k|k}$ ), see Kailath page 374.

# ALTERNATIVE FORM

**Rauch-Tung-Striebel (RTS) Formulas** (1965)

$$\hat{x}_{k|N} = \hat{x}_{k|k} + P_{k|k-1} F_{p,k}^* P_{k+1|k}^{-1} (\hat{x}_{k+1|N} - \hat{x}_{k+1|k})$$

$$P_{k|N} = P_{k|k} + P_{k|k-1} F_{p,k}^* P_{k+1|k}^{-1} (P_{k+1|N} - P_{k+1|k}) P_{k+1|k}^{-1} F_{p,k} P_{k|k-1}$$

**Also two-pass algorithm**



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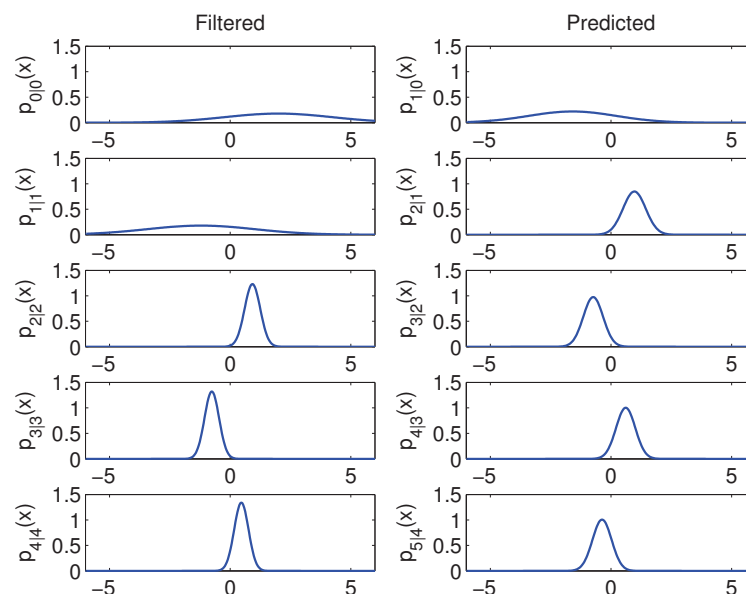
# NON-LINEAR FILTERING

- Assumption of linearity + Gaussianity often a rough approximation of reality!
- A single estimate of  $x_k$  is not always that useful.
- Full posterior distribution of  $x_k$  given observations contains more information, in general.
- Alternative view of Kalman:
  - Recursive calculation of posterior distributions  $p(x_k|\{y_1, \dots, y_{k-1}\})$  and  $p(x_k|\{y_1, \dots, y_{k-1}\})$
  - $p(x_k|\{y_1, \dots, y_{k-1}\}) \in \mathcal{CN}(\hat{x}_{k|k-1}, P_{k|k-1})$
  - $p(x_k|\{y_1, \dots, y_k\}) \in \mathcal{CN}(\hat{x}_{k|k}, P_{k|k})$
  - Parametric representation:  $\hat{x}_{k|m}$  and  $P_{k|m}$  specifies the Gaussian distribution fully



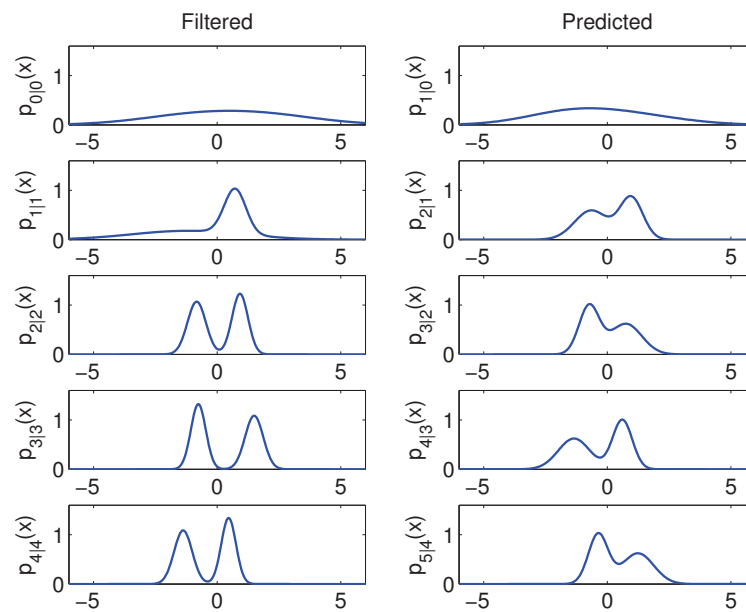
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## KALMAN FILTER – ALTERNATIVE VIEW



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## NON-GAUSSIAN EXAMPLE



## GENERAL MODEL

$$x_{k+1} = f(x_k, w_k)$$

$$y_k = h(x_k) + v_k$$

More general:

$$f(x_{k+1}, x_k, w_k) = 0$$

$$h(y_k, x_k, v_k) = 0$$

Even more general:

$$x_{k+1} \sim p(x_{k+1} | x_k)$$

$$y_k \sim p(y_k | x_k)$$



## GENERAL BAYESIAN SOLUTION

$$p(x_k|y_{1:k}) = \frac{p(y_k|x_k)p(x_k|y_{1:k-1})}{p(y_k|y_{1:k-1})}$$

$$p(y_k|y_{1:k-1}) = \int p(y_k|x_k)p(x_k|y_{1:k-1}) dx_k$$

$$p(x_{k+1}|y_{1:k}) = \int p(x_{k+1}|x_k)p(x_k|y_{1:k}) dx_k$$



Problems:

- How represent the distributions?
- How calculate the integrals?

## ALTERNATIVE REPRESENTATIONS OF DISTRIBUTIONS

- Parametric. Only possible for Gaussian + a few more  $\implies$  alternative derivation of Kalman
- Approximate by Gaussian, update mean and variance. Many possibilities.
- Discrete representation
  - Fixed grid
  - Stochastic grid (Monte Carlo, ...)
- Gaussian mixture model – linear combination of a number of Gaussian.



## IMPORTANCE SAMPLING

- Standard Monte Carlo:

$$\mathbb{E}[g(x)] = \int g(x)p(x) dx \approx \frac{1}{N} \sum_{n=1}^N g(x_n)$$

where  $x_n$  i.i.d. drawn from distribution  $p(x)$ .

- Importance Sampling: Introduce suitable proposal density  $q(x)$ :

$$\mathbb{E}[g(x)] = \int g(x)p(x) dx = \int g(x) \underbrace{\frac{p(x)}{q(x)}}_{\text{weight}} \underbrace{q(x)}_{\text{proposal dens.}} \approx \sum_{n=1}^N g(x_n)w_n$$

where  $x_n$  drawn from the proposal density  $q(x)$  and  $w_n = \frac{p(x_n)}{Nq(x_n)}$ .

## IMPORTANCE SAMPLING CONT.

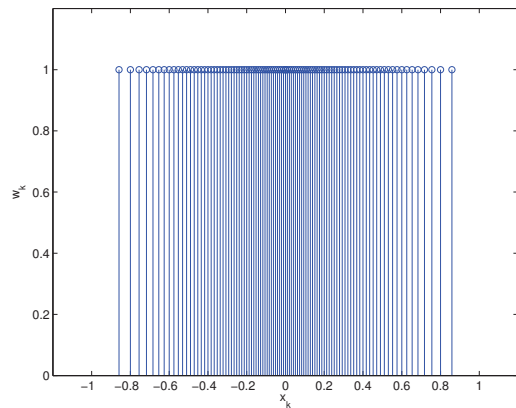
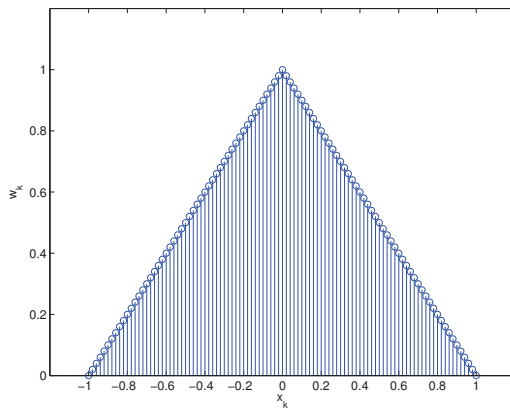
Equivalently  $p(x)$  is approximated by

$$p(x) \approx \sum_{n=1}^N w_n \delta(x - x_n)$$

$\{x_n, w_n\}$  called **particle distribution**.

Sometimes  $p(x)$  contain an unknown scaling  $\implies$  weights  $w_n$  are scaled. Easy solution is to normalize so  $\sum w_n = 1$ .

# ILLUSTRATION IMPORTANCE SAMPLING



**Case 1:** Uniformly distributed particles

**Case 2:** Equally weighted particles

Note: Both represent the same distribution !