

TTK4250

# Week 3

From the Kalman filter to stochastic processes

Edmund Førland Brekke

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## Recap from last week

### An estimator is a random variable

- ... because it depends on the (random) data.
- We can talk about its distribution, expectation and covariance.
- An estimator is unbiased if  $E[\mathbf{x} - \hat{\mathbf{x}}] = \mathbf{0}$ .

### LMMSE estimation

$$\hat{\mathbf{x}} = E[\mathbf{x}] + \text{Cov}(\mathbf{x}, \mathbf{z})\text{Cov}(\mathbf{z})^{-1}(\mathbf{z} - E[\mathbf{z}])$$

is the estimator of the form  $\hat{\mathbf{x}} = \mathbf{A}\mathbf{z} + \mathbf{b}$  that minimizes

$$\text{MSE}(\hat{\mathbf{x}}) = E \left[ \|\mathbf{x} - \hat{\mathbf{x}}\|_2^2 \right].$$

### The multivariate Gaussian

$$p(\mathbf{x}, \mathbf{y}) = \mathcal{N} \left( \begin{bmatrix} \mathbf{x} \\ \mathbf{y} \end{bmatrix}; \begin{bmatrix} \mathbf{a} \\ \mathbf{b} \end{bmatrix}, \begin{bmatrix} \mathbf{P}_{xx} & \mathbf{P}_{xy} \\ \mathbf{P}_{xy}^T & \mathbf{P}_{yy} \end{bmatrix} \right)$$

- Quadratic forms.
- Moment parametrization vs canonical parametrization.

### Marginalization and conditioning

In moment parametrization, conditioning is given by

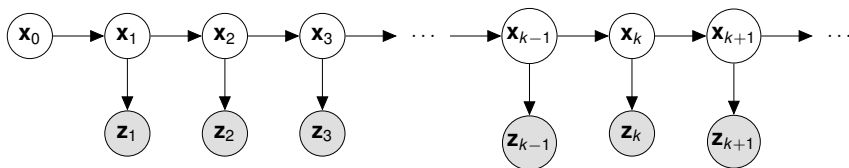
$$\boldsymbol{\mu}_{x|y} = \mathbf{a} + \mathbf{P}_{xy}\mathbf{P}_{yy}^{-1}(\mathbf{y} - \mathbf{b})$$

$$\mathbf{P}_{x|y} = \mathbf{P}_{xx} - \mathbf{P}_{xy}\mathbf{P}_{yy}^{-1}\mathbf{P}_{xy}^T.$$

This leads to the Product Identity.

# Recursive Bayesian estimation: Model and key concepts

We study systems whose structure fits the **graphical model** below:



- The horizontal arrows represent a **process model** of the form  $p(\mathbf{x}_k | \mathbf{x}_{k-1})$
- The vertical arrows represent a **measurement model** of the form  $p(\mathbf{z}_k | \mathbf{x}_k)$ .

This structure reflects the following **Markov assumptions**

$$p(\mathbf{x}_k | \mathbf{x}_1, \dots, \mathbf{x}_{k-2}, \mathbf{x}_{k-1}, \mathbf{z}_1, \dots, \mathbf{z}_{k-2}, \mathbf{z}_{k-1}) = p(\mathbf{x}_k | \mathbf{x}_{k-1})$$
$$p(\mathbf{z}_k | \mathbf{x}_1, \dots, \mathbf{x}_{k-2}, \mathbf{x}_{k-1}, \mathbf{x}_k, \mathbf{z}_1, \dots, \mathbf{z}_{k-2}, \mathbf{z}_{k-1}) = p(\mathbf{z}_k | \mathbf{x}_k)$$

## Recursive Bayesian estimation: The Bayes filter

In the Bayesian philosophy we want a pdf as our solution. This pdf may or may not be given by parameters such as expectation, covariance etc.

What do we know about  $\mathbf{x}_k$  after observing  $\mathbf{z}_{1:k} = (\mathbf{z}_1, \mathbf{z}_2, \dots, \mathbf{z}_k)$ ?

- The total probability theorem yields the predicted density

$$p(\mathbf{x}_k | \mathbf{z}_{1:k-1}) = \int p(\mathbf{x}_k | \mathbf{x}_{k-1}) p(\mathbf{x}_{k-1} | \mathbf{z}_{1:k-1}) d\mathbf{x}_{k-1}.$$

- Bayes' rule yields the posterior density

$$p(\mathbf{x}_k | \mathbf{z}_{1:k}) = \frac{p(\mathbf{z}_k | \mathbf{x}_k) p(\mathbf{x}_k | \mathbf{z}_{1:k-1})}{p(\mathbf{z}_k | \mathbf{z}_{1:k-1})} \propto p(\mathbf{z}_k | \mathbf{x}_k) p(\mathbf{x}_k | \mathbf{z}_{1:k-1}).$$

**Remark:** Violations of the Markov assumptions can be handled by replacing the Markov chain by a higher order Markov chain that models the temporal correlations. We must then extend the state vector with corresponding states.

# Linearity, Gaussianity and the Kalman filter

“Everything should be made as simple as possible, but not simpler.”

- In general, we cannot find a closed-form solution to the Bayes filter.
- If the posterior can be described with reasonable accuracy by a few parameters (e.g., expectation and covariance), then we should look for a compact representation.

Closed-form solution to the Bayes filter = Kalman filter

When does a closed-form solution to the Bayes filter exist?

- When the initial density is Gaussian  $\mathcal{N}(\mathbf{x}_0; \hat{\mathbf{x}}_0, \mathbf{P}_0)$
- ... and the Markov model is Gaussian-linear  $\mathcal{N}(\mathbf{x}_k; \mathbf{F}\mathbf{x}_{k-1}, \mathbf{Q})$
- ... and the likelihood is Gaussian-linear  $\mathcal{N}(\mathbf{z}_k; \mathbf{H}\mathbf{x}_k, \mathbf{R})$
- ... and standard independence assumptions apply.

# The prediction step of the Kalman filter

The predicted density is given by

$$\begin{aligned} p(\mathbf{x}_k | \mathbf{z}_{1:k-1}) &= \int p(\mathbf{x}_k | \mathbf{x}_{k-1}) p(\mathbf{x}_{k-1} | \mathbf{z}_{1:k-1}) d\mathbf{x}_{k-1} \\ &= \int \mathcal{N}(\mathbf{x}_k; \mathbf{F}\mathbf{x}_{k-1}, Q) \mathcal{N}(\mathbf{x}_{k-1}; \hat{\mathbf{x}}_{k-1}, \mathbf{P}_{k-1}) d\mathbf{x}_{k-1} \\ &= \mathcal{N}(\mathbf{x}_k; \mathbf{F}\hat{\mathbf{x}}_{k-1}, \mathbf{F}\mathbf{P}_{k-1}\mathbf{F}^T + Q) \\ &\quad \cdot \int \mathcal{N}(\mathbf{x}_{k-1}; \text{some vector, some covariance matrix}) d\mathbf{x}_{k-1} \\ &= \mathcal{N}(\mathbf{x}_k; \hat{\mathbf{x}}_{k|k-1}, \mathbf{P}_{k|k-1}). \end{aligned}$$

- $\hat{\mathbf{x}}_{k-1}$  is the previous state estimate.
- $\mathbf{P}_{k-1}$  is the previous covariance.
- $\hat{\mathbf{x}}_{k|k-1} = \mathbf{F}\hat{\mathbf{x}}_{k-1}$  is the predicted state estimate.
- $\mathbf{P}_{k|k-1} = \mathbf{F}\mathbf{P}_{k-1}\mathbf{F}^T + Q$  is the predicted covariance.

# The update step of the Kalman filter

The posterior density is given by

$$\begin{aligned} p(\mathbf{x}_k | \mathbf{z}_{1:k}) &\propto p(\mathbf{z}_k | \mathbf{x}_k) p(\mathbf{x}_k | \mathbf{z}_{1:k-1}) \\ &= \mathcal{N}(\mathbf{z}_k; \mathbf{H}\mathbf{x}_k, \mathbf{R}) \mathcal{N}(\mathbf{x}_k; \hat{\mathbf{x}}_{k|k-1}, \mathbf{P}_{k|k-1}) \\ &= \mathcal{N}(\mathbf{z}_k; \mathbf{H}\hat{\mathbf{x}}_{k|k-1}, \mathbf{H}\mathbf{P}_{k|k-1}\mathbf{H}^\top + \mathbf{R}) \mathcal{N}(\mathbf{x}_k; \hat{\mathbf{x}}_k, \mathbf{P}_k) \\ &\propto \mathcal{N}(\mathbf{x}_k; \hat{\mathbf{x}}_k, \mathbf{P}_k). \end{aligned}$$

- $\hat{\mathbf{x}}_k = \hat{\mathbf{x}}_{k|k-1} + \mathbf{W}_k(\mathbf{z}_k - \mathbf{H}\hat{\mathbf{x}}_{k|k-1})$  is the posterior state estimate.
- $\mathbf{P}_k = (\mathbf{I} - \mathbf{W}_k\mathbf{H})\mathbf{P}_{k|k-1}$  is the posterior covariance.
- $\mathbf{W}_k = \mathbf{P}_{k|k-1}\mathbf{H}^\top(\mathbf{H}\mathbf{P}_{k|k-1}\mathbf{H}^\top + \mathbf{R})^{-1}$  is the Kalman gain.

## More about the covariance

### Joseph form

$$\mathbf{P}_k = (\mathbf{I} - \mathbf{W}_k \mathbf{H}) \mathbf{P}_{k|k-1} (\mathbf{I} - \mathbf{W}_k \mathbf{H})^\top + \mathbf{W}_k \mathbf{R} \mathbf{W}_k^\top$$

### Information form

$$\mathbf{P}_k^{-1} = \mathbf{H}^\top \mathbf{R}^{-1} \mathbf{H} + \mathbf{P}_{k|k-1}^{-1}$$

### Orthogonality properties

- The estimation errors  $\tilde{\mathbf{x}}_k = \hat{\mathbf{x}}_k - \mathbf{x}_k$  do not constitute a white sequence:

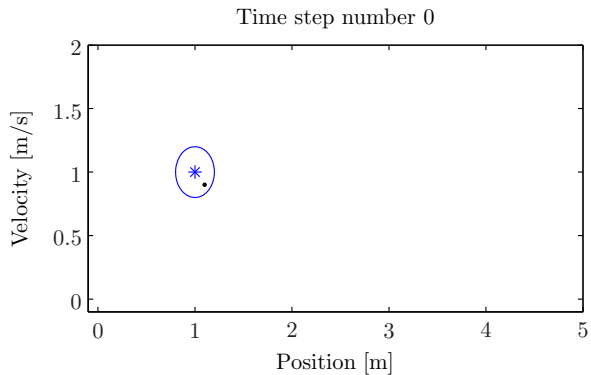
$$E[\tilde{\mathbf{x}}_k \tilde{\mathbf{x}}_{k-1}^\top] = (\mathbf{I} - \mathbf{W}_k \mathbf{H}) \mathbf{F} \mathbf{P}_k.$$

- The innovations  $\boldsymbol{\nu}_k = \mathbf{z}_k - \mathbf{H} \hat{\mathbf{x}}_{k|k-1}$  on the other hand are a white sequence:

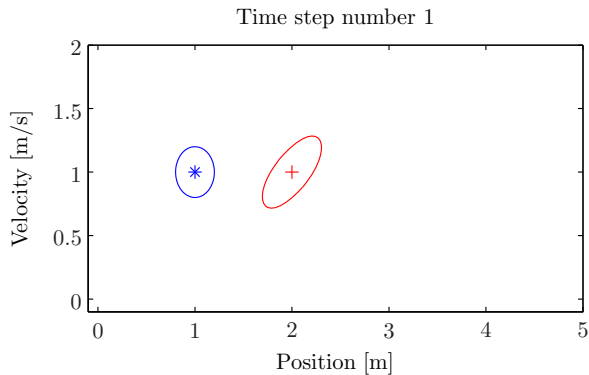
$$E[\boldsymbol{\nu}_k \boldsymbol{\nu}_j^\top] = \mathbf{0} \text{ if } k \neq j \Leftrightarrow p(\mathbf{z}_{1:k}) = \prod_{j=1}^k p(\boldsymbol{\nu}_j).$$



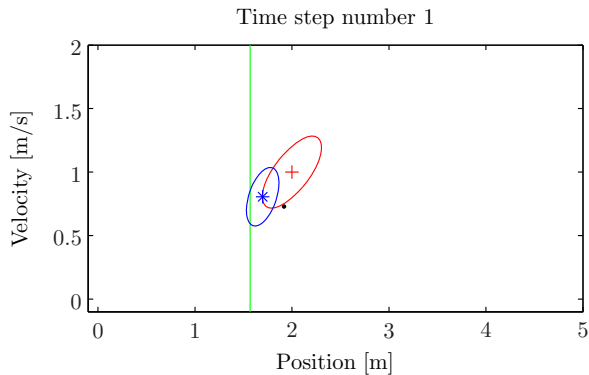
## Example run of the Kalman filter



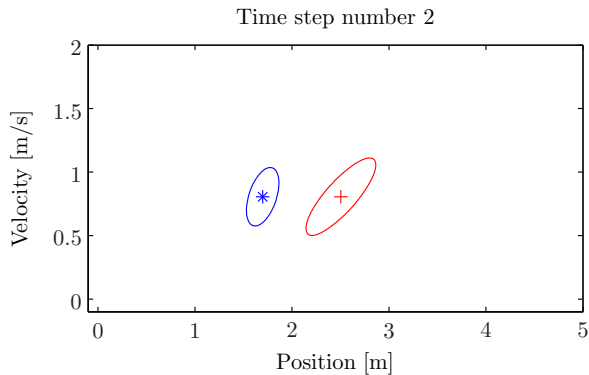
## Example run of the Kalman filter



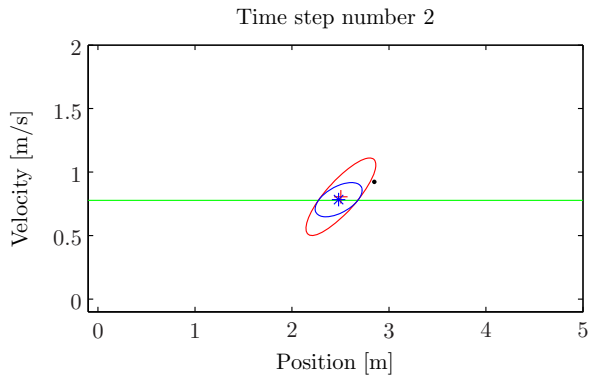
## Example run of the Kalman filter



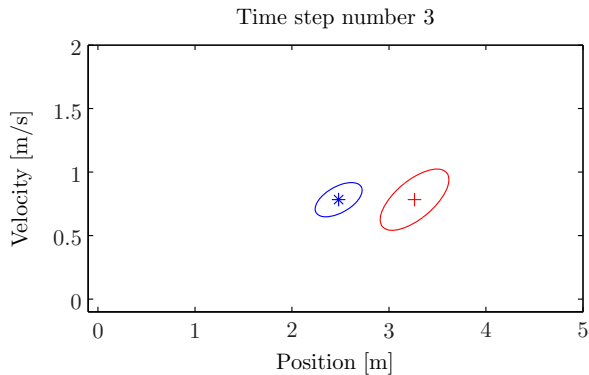
## Example run of the Kalman filter



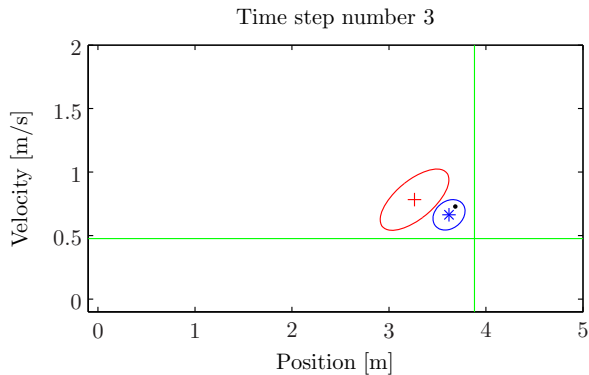
## Example run of the Kalman filter



## Example run of the Kalman filter



## Example run of the Kalman filter



## The tuning problem

Given the discrete time model

$$\mathbf{x}_k = \mathbf{F}\mathbf{x}_{k-1} + \mathbf{u}_k + \mathbf{v}_k, \quad \mathbf{z}_k = \mathbf{H}\mathbf{x}_k + \mathbf{w}_k, \quad \mathbf{v}_k \sim \mathcal{N}(\mathbf{0}, \mathbf{Q}), \quad \mathbf{w}_k \sim \mathcal{N}(\mathbf{0}, \mathbf{R})$$

we must determine the values of the noise matrices  $\mathbf{Q}$  and  $\mathbf{R}$  that ...

- faithfully represent the uncertainties of the process and measurement models.
- give the Kalman filter optimal **accuracy** and **robustness**.

### The process noise covariance

- The matrix  $\mathbf{Q}$  says something about how the system is expected to evolve between two time steps.
- But the system dynamics are generally modeled in continuous time.
- Therefore we need to relate  $\mathbf{Q}$  to a continuous-time model of the form

$$\dot{\mathbf{x}} = \mathbf{A}\mathbf{x} + \mathbf{B}\mathbf{u} + \mathbf{G}\mathbf{n}$$

### The measurement noise covariance

- The matrix  $\mathbf{R}$  says something about how accurate our measurement devices (sensors) are.
- This is fully encapsulated by the discrete-time model.



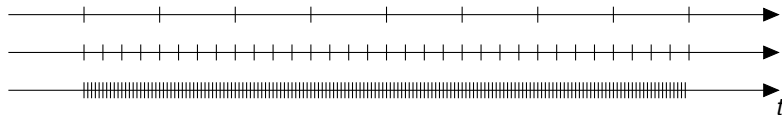
# Stochastic processes

- Consider a stochastic vector

$$\mathbf{x} = [x(t_1) \quad x(t_2) \quad \dots \quad x(t_n)]^T$$

where  $x(t_k)$  is the value of the stochastic variable  $x$  at time  $t_k$ .

- Let the discretization length  $T = t_k - t_{k-1}$  go towards zero.
- Every realization of  $\mathbf{x}$  will then be equivalent to a **function**  $x(t)$ . Such a random function is known as a **stochastic process**.

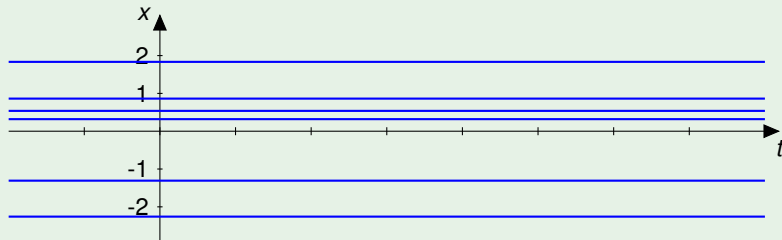


- To fully specify  $x(t)$  in the general case we would need the joint distribution of all tuples  $x(t_1), \dots, x(t_k)$  for any number  $k$ .
- We restrict our attention to stochastic process which can be defined in terms of their construction or in terms of first- and second-order moments.

# A very simple stochastic process

## A random constant

Let the function  $x(t)$  be given by  $x(t) = a$  where  $a \sim \mathcal{N}(0, 1)$ . Different realizations of this stochastic process can be depicted as follows:



Any number that depends on  $x(t)$ , such as a time integral of  $x(t)$ , will be a random variable. Let

$$y = \int_0^t x(\tau) d\tau.$$

Then it can be shown that  $y \sim \mathcal{N}(0, t^2)$ .

# A not so simple stochastic process

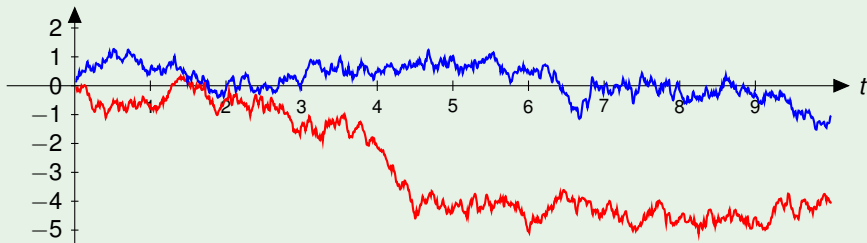
## The Wiener process

- Define the stochastic process  $x(t)$  by

$$x(nT) = \sum_{i=1}^n x_i \quad \text{where } x_i \sim \mathcal{N}(0, T) \text{ i.i.d.}$$

- Then we define the Wiener process  $b(t)$  as the limit

$$b(t) = \lim_{T \rightarrow 0} x(t).$$



# More about the Wiener process

## Alternative definition

Mathematicians like to define the Wiener process in terms of 4 fundamental properties:

- 1  $b(0) = 0$ .
- 2  $b(t)$  has independent increments. That is: If  $t_1 < t_2$ , then  $b(t_2) - b(t_1)$  is independent of the past values  $b(s)$  for  $s < t_1$ .
- 3  $b(t)$  has Gaussian increments: If  $t_1 < t_2$  then  $b(t_2) - b(t_1) \sim \mathcal{N}(0, t_2 - t_1)$
- 4  $b(t)$  is continuous in  $t$ .

## Statistics of the Wiener process

- The expectation of the Wiener process is always 0.
- The variance of the Wiener process at any particular time is

$$E[b(t)^2] = \text{Var} \left[ \sum_{i=1}^n x_i \right] = nT = \frac{t}{T} T = t.$$

## White Gaussian noise

We define **continuous-time white Gaussian noise** as the derivative of the Wiener process

$$n(t) = \lim_{\Delta \rightarrow 0} \frac{b(t + \Delta) - b(t)}{\Delta}.$$

- We always use white noise as a driving mechanism in stochastic continuous time models.
- For this to make sense, the contributions from a white noise process over a limited time interval must be finite and non-zero:

$$\Rightarrow 0 < \text{Var} \left[ \int_0^s n(t) dt \right] < \infty.$$

- Making matters complicated, this requirement in turn implies that

$$\text{Var} [n(t)] = \infty.$$

White noise is a mathematical abstraction because it has infinite energy.

## The autocorrelation function

Motivation: We want to have a useful description of important stochastic processes such as white Gaussian noise and its relatives.

### Definition: Autocorrelation function (ACF)

The ACF of a stochastic process  $\mathbf{x}(t)$  is  $R(t_1, t_2) = E[\mathbf{x}(t_1)\mathbf{x}(t_2)^T]$ .

### Definition: Wide-sense stationarity

A stochastic process  $\mathbf{x}(t)$  is said to be wide-sense stationary if its expectation is constant and its ACF can be written as a function of  $\tau = t_2 - t_1$ :

$$R(\tau) = E[\mathbf{x}(t)\mathbf{x}(t + \tau)^T].$$

### Example: ACF of white Gaussian noise

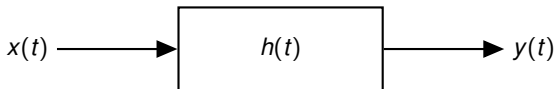
The ACF of the white noise process defined on the previous slide is

$$R(\tau) = \delta(\tau).$$

See the proof of Theorem 4.3.2 in the book for a derivation of this result.

## Stochastic linear systems

What happens to white noise (or any other stochastic process) when it is used as input to a system with a given impulse response?



### Convolution formulas for the ACF

Let  $x(t)$  be a scalar real-valued stochastic process with ACF  $R_{xx}(t_1, t_2)$  and let

$$y(t) = \int_{-\infty}^{\infty} h(t - \alpha)x(\alpha)d\alpha$$

where  $h(t)$  also is scalar real-valued. The the ACF of  $y(t)$  is given by

$$R_{xy}(t_1, t_2) = \int_{-\infty}^{\infty} R_{xx}(t_1, t_2 - \alpha)h(\alpha)d\alpha$$
$$R_{yy}(t_1, t_2) = \int_{-\infty}^{\infty} R_{xy}(t_1 - \alpha, t_2)h(\alpha)d\alpha.$$

# The Gauss-Markov process

- Consider a system with impulse response  $h(t) = e^{-ct}u(t)$ .
- We send white noise  $n(t)$  into the system, starting at  $t = 0$ .
- What is then the ACF of the output?

$$R_{xy}(t_1, t_2) = qe^{-c(t_2-t_1)}u(t_1)u(t_2 - t_1)$$

$$R_{yy}(t_1, t_2) = \frac{q}{2c}(1 - e^{-2ct_1})e^{-c(t_2-t_1)}.$$

- The formulas are valid if  $0 < t_1 < t_2$ .
- In the limit as  $t_1 \rightarrow \infty$  the Gauss-Markov process becomes a stationary process with ACF

$$\frac{q}{2c}e^{-c|t_2-t_1|}.$$



## Continuous time modeling: Accelerometer with bias

The Gauss-Markov process can be used to model a slowly varying accelerometer bias.

- Let the state vector be

$$\mathbf{x} = \begin{bmatrix} \text{Position of the vehicle} \\ \text{Velocity of the vehicle} \\ \text{Bias of the accelerometer} \end{bmatrix}$$

- The system is described by a state-space model of the form  $\dot{\mathbf{x}} = \mathbf{Ax} + \mathbf{Bu} + \mathbf{Gn}$  where

$$\mathbf{A} = \begin{bmatrix} 0 & 1 & 0 \\ 0 & 0 & 1 \\ 0 & 0 & -c \end{bmatrix}, \quad \mathbf{B} = \begin{bmatrix} 0 \\ 1 \\ 0 \end{bmatrix}, \quad \mathbf{G} = \begin{bmatrix} 0 & 0 \\ 1 & 0 \\ 0 & 1 \end{bmatrix}$$

and where

$$\mathbf{n} \sim \mathcal{N}(\mathbf{0}, \mathbf{D}\delta(t - \tau)) \text{ where } \mathbf{D} = \begin{bmatrix} \sigma_a^2 & 0 \\ 0 & \sigma_b^2 \end{bmatrix}.$$

- Notice that the accelerometer readings are treated as a control input and not as measurements.

## Continuous time modeling: The CV model in 2 dimensions

- This is perhaps the most common model used in sensor fusion.
- The model is of the form  $\dot{\mathbf{x}} = \mathbf{A}\mathbf{x} + \mathbf{G}\mathbf{n}$  where the matrices are given by

$$\mathbf{A} = \begin{bmatrix} 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \end{bmatrix} \quad \text{and} \quad \mathbf{G} = \begin{bmatrix} 0 & 0 \\ 0 & 0 \\ 1 & 0 \\ 0 & 1 \end{bmatrix}$$

and the process noise is given by

$$\mathbf{n} \sim \mathcal{N}(\mathbf{0}, \mathbf{D}\delta(t - \tau)) \quad \text{where} \quad \mathbf{D} = \begin{bmatrix} \sigma_a^2 & 0 \\ 0 & \sigma_a^2 \end{bmatrix}.$$

- We see that the process noise strength is solely given by the number  $\sigma_a$ , which is a measure of root-mean-square acceleration.
- Since the model essentially integrates white noise the two positional states become independent Wiener processes.

## Discretization

Consider the linear continuous-time state space model

$$\dot{\mathbf{x}} = \mathbf{A}\mathbf{x} + \mathbf{B}\mathbf{u} + \mathbf{G}\mathbf{n}, \quad \mathbf{n} \sim \mathcal{N}(\mathbf{0}, \mathbf{D}\delta(t - \tau)).$$

A discrete time solution can be written

$$\mathbf{x}_k = \mathbf{F}\mathbf{x}_{k-1} + \mathbf{u}_k + \mathbf{v}_k$$

where

$$\mathbf{F} = e^{\mathbf{A}(t_k - t_{k-1})}, \quad \mathbf{u}_k = \int_{t_{k-1}}^{t_k} e^{\mathbf{A}(t_k - \tau)} \mathbf{B}\mathbf{u}(\tau) d\tau \quad \text{and} \quad \mathbf{v}_k = \int_{t_{k-1}}^{t_k} e^{\mathbf{A}(t_k - \tau)} \mathbf{G}\mathbf{n}(\tau) d\tau.$$

### From continuous to discrete time process noise covariance

Let the discretization time be fixed at  $T = t_k - t_{k-1}$ . The covariance matrix of  $\mathbf{v}_k$  in the discrete-time model is then given by

$$\mathbf{Q} = E[\mathbf{v}_k \mathbf{v}_k^T] = \int_0^T e^{(T-\tau)\mathbf{A}} \mathbf{G} \mathbf{D} \mathbf{G}^T e^{(T-\tau)\mathbf{A}^T} d\tau$$

# Evaluating the discrete-time process noise matrix

## First/nth order approximations

- The simplest possible approximation is  $\mathbf{Q} \approx \mathbf{G}\mathbf{D}\mathbf{G}^T T$ .
- Not recommended because it may be singular even if the true  $\mathbf{Q}$  is SPD.
- Higher-order approximations are often used.
- Notice the linear dependence on the sampling interval  $T$ . This reflects the fact that the variance of the Wiener process increases linearly with time.

## Van Loan's formula

Define the matrices  $\mathbf{V}_1$  and  $\mathbf{V}_2$  according to

$$\exp\left(\begin{bmatrix} -\mathbf{A} & \mathbf{G}\mathbf{D}\mathbf{G}^T \\ \mathbf{0} & \mathbf{A}^T \end{bmatrix} T\right) = \begin{bmatrix} \times & \mathbf{V}_2 \\ \mathbf{0} & \mathbf{V}_1 \end{bmatrix}.$$

Then we can find  $\mathbf{Q}$  according to  $\mathbf{Q} = \mathbf{V}_1^T \mathbf{V}_2$ .

## How to tune the process noise covariance

There are at least 3 approaches that can be used to determine suitable values for the process noise covariance.

### Purely physical considerations

- In a CV model we may use the largest accelerations observed as a guideline for how large  $\sigma_a$  needs to be.
- In the accelerometer model we can calculate  $\sigma_a$  as a continuous-time equivalent of the specified accuracy of the accelerometer.

### Consistency analysis

Set the process noise as high as required to make the data or state estimates **plausible**.

### Maximum likelihood estimation

Find the **most likely** value of the process noise strength given the data.

**The process noise strength should not depend on the measurement model.**

# The concept of filter consistency

A filter is said to be consistent if

- 1 The state errors are acceptable as zero mean.
- 2 The state errors have magnitude commensurate with the state covariance yielded by the filter.
- 3 The innovations are acceptable as zero mean.
- 4 The innovations have magnitude commensurate with the innovation covariance yielded by the filter.
- 5 The innovations are acceptable as white.

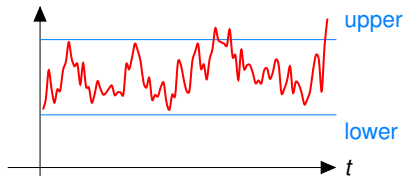
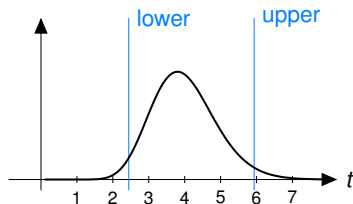
Criteria 2 and 4 are most important, and are tested by means of the normalized estimation error squared (NEES) and the normalized innovations squared (NIS):

$$\epsilon_k = (\hat{\mathbf{x}}_k - \mathbf{x}_k)^T \mathbf{P}_k^{-1} (\hat{\mathbf{x}}_k - \mathbf{x}_k)$$
$$\epsilon_k^\nu = \boldsymbol{\nu}_k^T \mathbf{S}_k^{-1} \boldsymbol{\nu}_k = (\mathbf{z}_k - \hat{\mathbf{z}}_{k|k-1})^T \mathbf{S}_k^{-1} (\mathbf{z}_k - \hat{\mathbf{z}}_{k|k-1}).$$

## $\chi^2$ -test for filter consistency (NEES)

- Suppose that  $\mathbf{x} \in \mathbb{R}^d$ .
- We perform  $N$  Monte-Carlo simulations of our Kalman filter.
- Then a 95% confidence interval for the average (ANEES) value of  $\epsilon_k$  is given by

$$\text{lower} = \text{chi2inv}(0.025, Nd)/N \quad \text{and} \quad \text{upper} = \text{chi2inv}(0.975, Nd)/N$$



- If ANEES lies below the  $\chi^2$  limits we can expect the filter to be overly conservative.
- If ANEES lies above the  $\chi^2$  limits we can expect the filter to be overconfident and put too little emphasis on the measurements.

Use NIS in a similar manner if working on real data without ground truth.

## Example of consistency-based tuning

Finding a suitable value for  $\sigma_a$  in the CV model originally used in the radar-based Autosea tracker.<sup>1</sup>

**Table 2** Process noise evaluation via AIS filter consistency. The  $(r_1, r_2)$  interval is the two-sided 95% probability concentration region for the  $\chi^2$  distribution related to the corresponding NIS. This varies with according to the AIS data record length  $N$ . The NIS values that are closest to being covariance-consistent, i.e. closest to the 95% probability region, are emphasised in bold

Name	$\sigma_a = 0.05$		$\sigma_a = 0.5$		$(r_1, r_2)$	$N$
	NIS	AI	NIS	AI		
GLUTRA	<b>4.67</b>	-0.02	0.90	0.01	(3.47, 4.55)	109
SULA	<b>3.61</b>	-0.22	0.51	-0.10	(3.49, 4.56)	106
KORSFJORD	71.8	-1.33	<b>4.31</b>	-0.44	(3.52, 4.51)	127
TR.FJORD II	11.3	-0.62	<b>3.24</b>	-0.16	(3.76, 4.24)	533
TELEMETRON	371	-0.04	<b>4.45</b>	-0.01	(3.77, 4.23)	579



<sup>1</sup>Wilthil et al. (2017): "A target tracking system for ASV collision avoidance based on the PDAF", Springer.



# Testing the whiteness of the innovations

## Whiteness test in Monte-Carlo simulations

Let  $\nu_k$  be one of the innovation states in the vector  $\nu_k$ . Let  $N$  be the number of Monte-Carlo simulations. Then the distribution of the **sample autocorrelation**

$$\rho_{kj} = \frac{\sum_{i=1}^N \nu_k^{(i)} \nu_j^{(i)}}{\sqrt{\sum_{i=1}^N (\nu_k^{(i)})^2 \sum_{i=1}^N (\nu_j^{(i)})^2}}$$

should tend to  $\mathcal{N}(0, 1/N)$  for all  $k \neq j$  when  $N$  is large.

## Single-run whiteness test

The variance of the **time-average autocorrelation** should tend towards  $1/K$ :

$$\bar{\rho}_j = \frac{\sum_{k=1}^K \nu_k \nu_{k+j}}{\sqrt{\sum_{k=1}^K \nu_k^2 \sum_{k=1}^K \nu_{k+j}^2}}$$

## Tuning the measurement noise covariance

For exteroceptive sensors, the appropriate values in  $\mathbf{R}$  depend on

- The sensor resolution.
- The extent of targets or landmarks.

### Example: Point targets with pixellated sensor

- 2-dimensional sensor with square cells of fixed resolution  $\Delta x$ .
- $\mathbf{x} = [x, y, v_x, v_y]^T$ .
- Measurement matrix  $\mathbf{H} = [\mathbf{I}_2, \mathbf{0}]$

$$p(\mathbf{z} - \mathbf{H}\mathbf{x}|\mathbf{x}) = \begin{cases} 1/\Delta x^2 & \text{if } \|\mathbf{z} - \mathbf{H}\mathbf{x}\|_\infty < \Delta x/2 \\ 0 & \text{otherwise.} \end{cases}$$

This distribution can be approximated by a Gaussian

$$p(\mathbf{z}|\mathbf{x}) = \mathcal{N}(\mathbf{z}; \mathbf{H}\mathbf{x}, \mathbf{R})$$

with the same covariance:

$$\mathbf{R} = \begin{bmatrix} \frac{\Delta x^2}{12} & 0 \\ 0 & \frac{\Delta x^2}{12} \end{bmatrix}.$$

## More about the measurement model

Mild nonlinearities in the measurement model can sometimes be removed by converting the measurements. We must then also convert  $\mathbf{R}$  accordingly.

### Nonlinear model.

The model is of the form  $\mathbf{z}_k = \mathbf{h}(\mathbf{x}_k) + \mathbf{w}_k$ ,  $\mathbf{w}_k \sim \mathcal{N}(\mathbf{0}, \mathbf{R})$  where

$$\mathbf{h}(\mathbf{x}_k) = \begin{bmatrix} \sqrt{x^2 + y^2} \\ \text{atan2}(y, x) \end{bmatrix}$$
$$\mathbf{R} = \begin{bmatrix} \sigma_r^2 & 0 \\ 0 & \sigma_\theta^2 \end{bmatrix}$$

### Converted measurements.

The model is of the form  $\mathbf{z}_k = \mathbf{H}\mathbf{x}_k + \mathbf{w}_k$ ,  $\mathbf{w}_k \sim \mathcal{N}(\mathbf{0}, \mathbf{R}_c)$  where

$$\mathbf{H} = \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \end{bmatrix}$$
$$\mathbf{R}_c = \mathbf{J}\mathbf{R}\mathbf{J}^T, \quad \mathbf{J} = \frac{\partial}{\partial \mathbf{z}_k} \mathbf{h}^{-1}(\mathbf{z}_k)$$

For more sophisticated conversion techniques see Lerro & Bar-Shalom (1993): "Tracking with debiased consistent converted measurements versus EKF", IEEE-TAES.

Measurement conversion is generally preferable to EKF techniques because the linearization in an EKF can lead to filter instability.

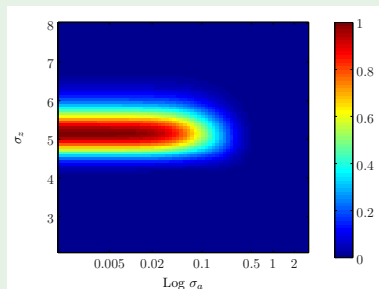
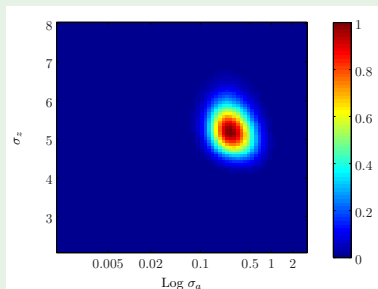
# Maximum likelihood estimation of system parameters

## Theorem for maximum likelihood estimation

Assume that both  $\mathbf{Q}$  and  $\mathbf{R}$  depend on an unknown parameter vector  $\mathbf{q}$ . The maximum likelihood estimate of  $\mathbf{q}$ , if it exists, can then be found as

$$\mathbf{q}_{\text{ML}} = \arg \max_{\mathbf{q}} \sum_k \log \mathcal{N}(\mathbf{z}_k; \mathbf{H}\hat{\mathbf{x}}_{k|k-1}, \mathbf{S}_k)$$

## Example: Estimating $\sigma_a$ and $\sigma_z$ for CV model



The likelihood function, normalized relative to its maximum, for two realizations of the CV model.

## LTV and LTI systems

We point out the main distinction in the continuous case. The discrete case is similar.

### Linear time-variant systems

The system can be of the form

$$\begin{aligned}\dot{\mathbf{x}} &= \mathbf{A}(t)\mathbf{x} + \mathbf{B}(t)\mathbf{u} + \mathbf{G}(t)\mathbf{n} \\ \mathbf{z} &= \mathbf{H}(t)\mathbf{x} + \mathbf{w}\end{aligned}$$

The system matrices are allowed to depend on  $t$ .

### Linear time-invariant systems

The system can be of the form

$$\begin{aligned}\dot{\mathbf{x}} &= \mathbf{A}\mathbf{x} + \mathbf{B}\mathbf{u} + \mathbf{G}\mathbf{n} \\ \mathbf{z} &= \mathbf{H}\mathbf{x} + \mathbf{w}\end{aligned}$$

The system matrices are not allowed to depend on  $t$ .

- In both cases, the matrices are assumed known.
- Uncertainty in the matrices can be modeled as a non-linear system.

# Observability for LTI systems

## Continuous-time observability

Consider a continuous-time LTI system of the form

$$\dot{\mathbf{x}} = \mathbf{A}\mathbf{x} + \mathbf{B}u$$

$$\mathbf{z} = \mathbf{H}\mathbf{x}.$$

The observability matrix of the system is

$$\mathbf{Q}_O = \begin{bmatrix} \mathbf{H} \\ \mathbf{H}\mathbf{A} \\ \vdots \\ \mathbf{H}\mathbf{A}^d \end{bmatrix}.$$

We say that the pair  $[\mathbf{A}, \mathbf{H}]$  is observable if  $\mathbf{Q}_O$  is of full rank.

## Discrete-time observability

Simply replace  $\mathbf{A}$  and  $\mathbf{B}$  with their discrete-time equivalents.

## Example: Observability for CV model

Consider the system model

$$\mathbf{A} = \begin{bmatrix} 0 & 1 \\ 0 & 0 \end{bmatrix} \qquad \mathbf{H} = \begin{bmatrix} h_1 & h_2 \end{bmatrix}.$$

The observability matrix is then

$$\mathbf{Q}_O = \begin{bmatrix} \mathbf{H} \\ \mathbf{H}\mathbf{A} \end{bmatrix} = \begin{bmatrix} h_1 & h_2 \\ 0 & h_1 \end{bmatrix}.$$

### Case 1: Only position measurements

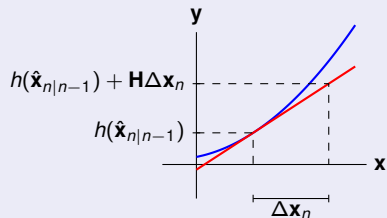
$h_2 = 0 \Rightarrow \mathbf{Q}_O = \mathbf{I} \Rightarrow \mathbf{Q}_O$  is of full rank  $\Rightarrow$  The system is observable.

### Case 2: Only velocity measurements

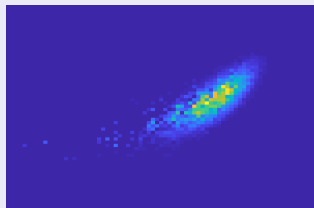
$h_1 = 0 \Rightarrow$  Second row of  $\mathbf{Q}_O$  is zero  $\Rightarrow$  The system is not observable.

# The road ahead

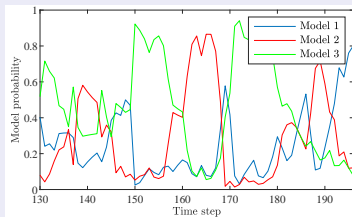
## The EKF



## Particle filters



## Interacting Multiple Models



## Target tracking

