

Lecture 2: From Linear Regression to Kalman Filter and Beyond

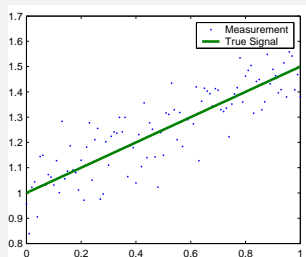
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Batch Linear Regression [1/2]



- Consider the **linear regression model**

$$y_k = a_1 + a_2 t_k + \epsilon_k,$$

with $\epsilon_k \sim \mathcal{N}(0, \sigma^2)$ and $\mathbf{a} = (a_1, a_2) \sim \mathcal{N}(\mathbf{m}_0, \mathbf{P}_0)$.

- In **probabilistic notation** this is:

$$p(y_k | \mathbf{a}) = \mathcal{N}(y_k | \mathbf{H}_k \mathbf{a}, \sigma^2)$$

$$p(\mathbf{a}) = \mathcal{N}(\mathbf{a} | \mathbf{m}_0, \mathbf{P}_0),$$

where $\mathbf{H}_k = (1 \ t_k)$.

- The **Bayesian batch solution** by the Bayes' rule:

$$\begin{aligned} p(\mathbf{a} | y_{1:N}) &\propto p(\mathbf{a}) \prod_{k=1}^N p(y_k | \mathbf{a}) \\ &= N(\mathbf{a} | \mathbf{m}_0, \mathbf{P}_0) \prod_{k=1}^N N(y_k | \mathbf{H}_k \mathbf{a}, \sigma^2). \end{aligned}$$

- The **posterior** is Gaussian

$$p(\mathbf{a} | y_{1:N}) = N(\mathbf{a} | \mathbf{m}_N, \mathbf{P}_N).$$

- The **mean and covariance** are given as

$$\begin{aligned} \mathbf{m}_N &= \left[\mathbf{P}_0^{-1} + \frac{1}{\sigma^2} \mathbf{H}^T \mathbf{H} \right]^{-1} \left[\frac{1}{\sigma^2} \mathbf{H}^T \mathbf{y} + \mathbf{P}_0^{-1} \mathbf{m}_0 \right] \\ \mathbf{P}_N &= \left[\mathbf{P}_0^{-1} + \frac{1}{\sigma^2} \mathbf{H}^T \mathbf{H} \right]^{-1}, \end{aligned}$$

where $\mathbf{H}_k = (1 \ t_k)$ and $\mathbf{H} = (\mathbf{H}_1; \mathbf{H}_2; \dots; \mathbf{H}_N)$, and

- Assume that we have already computed the posterior distribution, which is **conditioned on the measurement up to $k - 1$** :

$$p(\mathbf{a} | y_{1:k-1}) = \mathbf{N}(\mathbf{a} | \mathbf{m}_{k-1}, \mathbf{P}_{k-1}).$$

- Assume that we get the **k th measurement y_k** . Using the equations from the previous slide we get

$$\begin{aligned} p(\mathbf{a} | y_{1:k}) &\propto p(y_k | \mathbf{a}) p(\mathbf{a} | y_{1:k-1}) \\ &\propto \mathbf{N}(\mathbf{a} | \mathbf{m}_k, \mathbf{P}_k). \end{aligned}$$

- The **mean and covariance** are given as

$$\begin{aligned} \mathbf{m}_k &= \left[\mathbf{P}_{k-1}^{-1} + \frac{1}{\sigma^2} \mathbf{H}_k^T \mathbf{H}_k \right]^{-1} \left[\frac{1}{\sigma^2} \mathbf{H}_k^T y_k + \mathbf{P}_{k-1}^{-1} \mathbf{m}_{k-1} \right] \\ \mathbf{P}_k &= \left[\mathbf{P}_{k-1}^{-1} + \frac{1}{\sigma^2} \mathbf{H}_k^T \mathbf{H}_k \right]^{-1}. \end{aligned}$$

- By the **matrix inversion lemma** (or Woodbury identity):

$$\mathbf{P}_k = \mathbf{P}_{k-1} - \mathbf{P}_{k-1} \mathbf{H}_k^T \left[\mathbf{H}_k \mathbf{P}_{k-1} \mathbf{H}_k^T + \sigma^2 \right]^{-1} \mathbf{H}_k \mathbf{P}_{k-1}.$$

- Now the equations for the **mean and covariance** reduce to

$$\mathbf{S}_k = \mathbf{H}_k \mathbf{P}_{k-1} \mathbf{H}_k^T + \sigma^2$$

$$\mathbf{K}_k = \mathbf{P}_{k-1} \mathbf{H}_k^T \mathbf{S}_k^{-1}$$

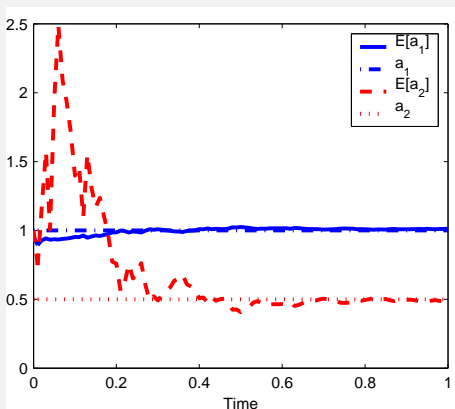
$$\mathbf{m}_k = \mathbf{m}_{k-1} + \mathbf{K}_k [y_k - \mathbf{H}_k \mathbf{m}_{k-1}]$$

$$\mathbf{P}_k = \mathbf{P}_{k-1} - \mathbf{K}_k \mathbf{S}_k \mathbf{K}_k^T.$$

- Computing these for $k = 0, \dots, N$ gives **exactly the linear regression solution** – but without a matrix inversion!
- A special case of **Kalman filter**.

Recursive Linear Regression [3/3]

Convergence of the recursive solution to the batch solution – on the last step the solutions are exactly equal:



General batch solution:

- Specify the **measurement model**:

$$p(\mathbf{y}_{1:N} | \boldsymbol{\theta}) = \prod_k p(\mathbf{y}_k | \boldsymbol{\theta}).$$

- Specify the **prior distribution** $p(\boldsymbol{\theta})$.
- Compute **posterior distribution** by the Bayes' rule:

$$p(\boldsymbol{\theta} | \mathbf{y}_{1:N}) = \frac{1}{Z} p(\boldsymbol{\theta}) \prod_k p(\mathbf{y}_k | \boldsymbol{\theta}).$$

- Compute point estimates, moments, predictive quantities etc. from the posterior distribution.

General recursive solution:

- Specify the **measurement likelihood** $p(\mathbf{y}_k | \theta)$.
- Specify the **prior distribution** $p(\theta)$.
- Process measurements $\mathbf{y}_1, \dots, \mathbf{y}_N$ **one at a time**, starting from the prior:

$$\begin{aligned} p(\theta | \mathbf{y}_1) &= \frac{1}{Z_1} p(\mathbf{y}_1 | \theta) p(\theta) \\ p(\theta | \mathbf{y}_{1:2}) &= \frac{1}{Z_2} p(\mathbf{y}_2 | \theta) p(\theta | \mathbf{y}_1) \\ &\vdots \\ p(\theta | \mathbf{y}_{1:N}) &= \frac{1}{Z_N} p(\mathbf{y}_N | \theta) p(\theta | \mathbf{y}_{1:N-1}). \end{aligned}$$

- The posterior at the last step is the **same as the batch solution**.

Advantages of Recursive Solution

- The recursive solution can be considered as the **online learning** solution to the Bayesian learning problem.
- **Batch** Bayesian inference is a **special case of recursive** Bayesian inference.
- The **parameter** can be modeled to **change** between the measurement steps \Rightarrow **basis of filtering theory**.

- Let assume **Gaussian random walk** between the measurements in the linear regression model:

$$\begin{aligned}p(y_k | \mathbf{a}_k) &= \text{N}(y_k | \mathbf{H}_k \mathbf{a}_k, \sigma^2) \\p(\mathbf{a}_k | \mathbf{a}_{k-1}) &= \text{N}(\mathbf{a}_k | \mathbf{a}_{k-1}, \mathbf{Q}) \\p(\mathbf{a}_0) &= \text{N}(\mathbf{a}_0 | \mathbf{m}_0, \mathbf{P}_0).\end{aligned}$$

- Again, assume that we already know

$$p(\mathbf{a}_{k-1} | y_{1:k-1}) = \text{N}(\mathbf{a}_{k-1} | \mathbf{m}_{k-1}, \mathbf{P}_{k-1}).$$

- The **joint distribution** of \mathbf{a}_k and \mathbf{a}_{k-1} is (due to Markovianity of dynamics!):

$$p(\mathbf{a}_k, \mathbf{a}_{k-1} | y_{1:k-1}) = p(\mathbf{a}_k | \mathbf{a}_{k-1}) p(\mathbf{a}_{k-1} | y_{1:k-1}).$$

- Integrating over \mathbf{a}_{k-1} gives:

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

- This equation for **Markov processes** is called the **Chapman-Kolmogorov equation**.
- Because the distributions are Gaussian, the **result is Gaussian**

$$p(\mathbf{a}_k | y_{1:k-1}) = \mathbf{N}(\mathbf{a}_k | \mathbf{m}_k^-, \mathbf{P}_k^-),$$

where

$$\mathbf{m}_k^- = \mathbf{m}_{k-1}$$

$$\mathbf{P}_k^- = \mathbf{P}_{k-1} + \mathbf{Q}.$$

- As in the pure recursive estimation, we get

$$\begin{aligned} p(\mathbf{a} | y_{1:k}) &\propto p(y_k | \mathbf{a}) p(\mathbf{a} | y_{1:k-1}) \\ &\propto \mathbf{N}(\mathbf{a} | \mathbf{m}_k, \mathbf{P}_k). \end{aligned}$$

- After applying the matrix inversion lemma, **mean and covariance** can be written as

$$\begin{aligned} \mathbf{S}_k &= \mathbf{H}_k \mathbf{P}_k^- \mathbf{H}_k^T + \sigma^2 \\ \mathbf{K}_k &= \mathbf{P}_k^- \mathbf{H}_k^T \mathbf{S}_k^{-1} \\ \mathbf{m}_k &= \mathbf{m}_k^- + \mathbf{K}_k [y_k - \mathbf{H}_k \mathbf{m}_k^-] \\ \mathbf{P}_k &= \mathbf{P}_k^- - \mathbf{K}_k \mathbf{S}_k \mathbf{K}_k^T. \end{aligned}$$

- Again, we have derived a special case of the **Kalman filter**.
- The **batch version** of this solution would be **much more complicated**.

State Space Notation

- In the previous section we formulated the model as

$$p(\mathbf{a}_k | \mathbf{a}_{k-1}) = N(\mathbf{a}_k | \mathbf{a}_{k-1}, \mathbf{Q})$$

$$p(y_k | \mathbf{a}_k) = N(y_k | \mathbf{H}_k \mathbf{a}_k, \sigma^2)$$

- But in **Kalman filtering and control theory** the vector of parameters \mathbf{a}_k is usually called “state” and denoted as \mathbf{x}_k .
- More standard **state space notation**:

$$p(\mathbf{x}_k | \mathbf{x}_{k-1}) = N(\mathbf{x}_k | \mathbf{x}_{k-1}, \mathbf{Q})$$

$$p(y_k | \mathbf{x}_k) = N(y_k | \mathbf{H}_k \mathbf{x}_k, \sigma^2)$$

- Or equivalently

$$\mathbf{x}_k = \mathbf{x}_{k-1} + \mathbf{q}$$

$$y_k = \mathbf{H}_k \mathbf{x}_k + r,$$

where $\mathbf{q} \sim N(\mathbf{0}, \mathbf{Q})$, $r \sim N(0, \sigma^2)$.

- The **canonical Kalman filtering model** is

$$p(\mathbf{x}_k | \mathbf{x}_{k-1}) = N(\mathbf{x}_k | \mathbf{A}_{k-1} \mathbf{x}_{k-1}, \mathbf{Q}_{k-1})$$

$$p(\mathbf{y}_k | \mathbf{x}_k) = N(\mathbf{y}_k | \mathbf{H}_k \mathbf{x}_k, \mathbf{R}_k).$$

- More often, this model can be seen **in the form**

$$\mathbf{x}_k = \mathbf{A}_{k-1} \mathbf{x}_{k-1} + \mathbf{q}_{k-1}$$

$$\mathbf{y}_k = \mathbf{H}_k \mathbf{x}_k + \mathbf{r}_k.$$

- The Kalman filter actually calculates the following distributions:

$$p(\mathbf{x}_k | \mathbf{y}_{1:k-1}) = N(\mathbf{x}_k | \mathbf{m}_k^-, \mathbf{P}_k^-)$$

$$p(\mathbf{x}_k | \mathbf{y}_{1:k}) = N(\mathbf{x}_k | \mathbf{m}_k, \mathbf{P}_k).$$

- **Prediction step** of the Kalman filter:

$$\mathbf{m}_k^- = \mathbf{A}_{k-1} \mathbf{m}_{k-1}$$

$$\mathbf{P}_k^- = \mathbf{A}_{k-1} \mathbf{P}_{k-1} \mathbf{A}_{k-1}^T + \mathbf{Q}_{k-1}.$$

- **Update step** of the Kalman filter:

$$\mathbf{S}_k = \mathbf{H}_k \mathbf{P}_k^- \mathbf{H}_k^T + \mathbf{R}_k$$

$$\mathbf{K}_k = \mathbf{P}_k^- \mathbf{H}_k^T \mathbf{S}_k^{-1}$$

$$\mathbf{m}_k = \mathbf{m}_k^- + \mathbf{K}_k [\mathbf{y}_k - \mathbf{H}_k \mathbf{m}_k^-]$$

$$\mathbf{P}_k = \mathbf{P}_k^- - \mathbf{K}_k \mathbf{S}_k \mathbf{K}_k^T.$$

- These equations will be derived from the general Bayesian filtering equations in the next lecture.

- Generic discrete-time state space models

$$\mathbf{x}_k = \mathbf{f}(\mathbf{x}_{k-1}, \mathbf{q}_k)$$

$$\mathbf{y}_k = \mathbf{h}(\mathbf{x}_k, \mathbf{r}_k).$$

- Generic Markov models

$$\mathbf{y}_k \sim p(\mathbf{y}_k | \mathbf{x}_k)$$

$$\mathbf{x}_k \sim p(\mathbf{x}_k | \mathbf{x}_{k-1}).$$

- Approximation methods: Extended Kalman filters (EKF), Unscented Kalman filters (UKF), sequential Monte Carlo (SMC) filters aka particle filters.

- In **continuous-discrete filtering models**, dynamics are modeled in continuous time, measurements at discrete time steps.
- The continuous time versions of Markov models are called as **stochastic differential equations**:

$$\frac{d\mathbf{x}}{dt} = \mathbf{f}(\mathbf{x}, t) + \mathbf{w}(t)$$

where $\mathbf{w}(t)$ is a continuous time Gaussian white noise process.

- Approximation methods: Extended Kalman filters, Unscented Kalman filters, sequential Monte Carlo, particle filters.

Summary

- **Linear regression problem** can be solved as **batch problem** or **recursively** – the latter solution is a special case of **Kalman filter**.
- A generic **Bayesian estimation problem** can also be solved as **batch problem** or **recursively**.
- If we let the linear regression **parameter change** between the measurements, we get a simple **linear state space model** – again solvable with **Kalman filtering model**.
- By **generalizing this idea** and the solution we get the **Kalman filter** algorithm.
- By further generalizing to **non-Gaussian models** results in a generic **probabilistic state space model**.

Batch and recursive linear regression.