

PROBABILISTIC MACHINE LEARNING

LECTURE 21

HIDDEN MARKOV MODELS

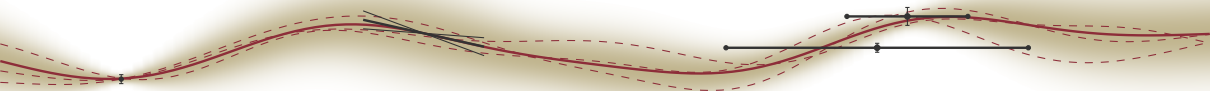
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Reminder: Goal for this week – *Time Series* as a problem class.

- ▶ Application Layer: Data arriving as a stream
- ▶ Model Structure Layer: Markov Chains / Hidden Markov Models
- ▶ Concrete Model Layer: Gauss–Markov Models
- ▶ Algorithm Layer: Kalman Filter & RTS Smoother

Today:

Theory: What is the connection between Gaussian processes and Gauss–Markov Processes?

Parameters: Can we learn the parameters of a Gauss–Markov model?

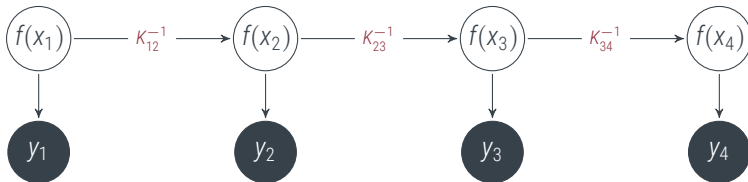
Generalization: What if the world isn't Gaussian?



Recap: Markov Chains

Processes with the *state* as a local memory

$$p(f) = \mathcal{GP}(f; 0, k) \quad p\left(\begin{bmatrix} f_1 \\ f_2 \\ f_3 \\ f_4 \end{bmatrix}\right) = \mathcal{N}\left(0, \begin{bmatrix} K_{11}^{-1} & K_{12}^{-1} & 0 & 0 \\ K_{12}^{-1} & K_{22}^{-1} & K_{23}^{-1} & 0 \\ 0 & K_{23}^{-1} & K_{33}^{-1} & K_{34}^{-1} \\ 0 & 0 & K_{34}^{-1} & K_{44}^{-1} \end{bmatrix}^{-1}\right) \quad p(y | f) = \prod_i \mathcal{N}(y_i; f_i, \sigma^2)$$



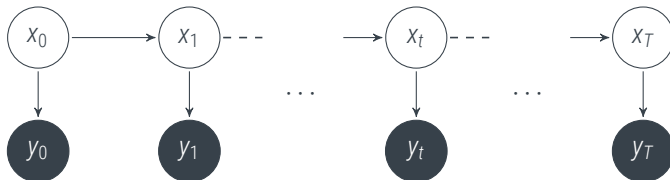
Recap: Markov Chains

Endow the algorithmic structure of filtering and smoothing

Assume:

$$p(x_t | X_{0:t-1}) = p(x_t | x_{t-1})$$

and $p(y_t | X) = p(y_t | x_t)$



Filtering: $\mathcal{O}(T)$

predict: $p(x_t | Y_{0:t-1}) = \int p(x_t | x_{t-1}) p(x_{t-1} | Y_{0:t-1}) dx_{t-1}$ (Chapman-Kolmogorov Eq.)

update: $p(x_t | Y_{0:t}) = \frac{p(y_t | x_t) p(x_t | Y_{0:t-1})}{p(y_t)}$ (Bayes' Theorem)

Smoothing: $\mathcal{O}(T)$

smooth: $p(x_t | Y) = p(x_t | Y_{0:t}) \int p(x_{t+1} | x_t) \frac{p(x_{t+1} | Y)}{p(x_{t+1} | Y_{1:t})} dx_{t+1}$ (backward pass)

Time Series:

- ▶ **Markov Chains** formalize the notion of a stochastic process with a *local finite memory*
- ▶ Inference over Markov Chains separates into three operations, that can be performed in *linear* time:

Filtering: $\mathcal{O}(T)$

$$\text{predict:} \quad p(x_t \mid Y_{0:t-1}) = \int p(x_t \mid x_{t-1}) p(x_{t-1} \mid Y_{0:t-1}) dx_{t-1} \quad (\text{Chapman-Kolmogorov Eq.})$$

$$\text{update:} \quad p(x_t \mid Y_{0:t}) = \frac{p(y_t \mid x_t) p(x_t \mid Y_{0:t-1})}{p(y_t)} \quad (\text{Bayes' Theorem})$$

Smoothing: $\mathcal{O}(T)$

$$\text{smooth:} \quad p(x_t \mid Y) = p(x_t \mid Y_{0:t}) \int p(x_{t+1} \mid x_t) \frac{p(x_{t+1} \mid Y)}{p(x_{t+1} \mid Y_{0:t})} dx_{t+1} \quad (\text{backward pass})$$

```

1 procedure INFERENCE( $Y, p(x_0), p(x_t | x_{t-1}) \forall t, p(y_t | x_t) \forall t$ )
2   for  $i=1, \dots, n$  do                                     // Filtering
3      $p(x_t | y_{1:t-1}) = \int p(x_t | x_{t-1}) p(x_{t-1} | Y_{0:t-1}) dx_{t-1}$  // Chapman-Kolmogorov eq.
4      $p(x_t | y_{1:t}) = p(y_t | x_t) p(x_t | Y_{0:t-1}) / p(y_t)$  // Update
5   end for
6   for  $i=n-1, \dots, 0$  do                                   // Smoothing
7      $p(x_t | Y) = p(x_t | Y_{0:t}) \int p(x_{t+1} | x_t) p(x_{t+1} | Y) p(x_{t+1} | Y_{1:t}) dx_{t+1}$ 
8   end for
9   return  $p(x_t | Y) \forall t = 0, \dots, n$  // return all marginals
10 end procedure

```

Gauss–Markov Models

Local structure for univariate Gaussian models

$$p(x(t_{i+1}) \mid X_{1:i}) = \mathcal{N}(x_{i+1}; Ax_i, Q) \quad \text{and} \quad p(x_0) = \mathcal{N}(x_0; m_0, P_0) \quad \text{and} \quad p(y_i \mid X) = \mathcal{N}(y_i; Hx_i, R)$$

```

1 procedure FILTER( $m_0, P_0, A, Q, H, R, y$ )
2   for  $t = 1, 2, \dots, T$  do                                     //  $\mathcal{O}(N)$ 
3      $m_t^- = Am_{t-1}$                                            // predictive mean
4      $P_t^- = AP_{t-1}A^\top + Q$                                    // predictive covariance.  $\mathcal{O}(|X|^3)$ 
5      $z = y - Hm_t^-$                                            // residual
6      $S = HP_t^-H^\top + R$                                        // innovation covariance
7      $K = P_t^-H^\top S^{-1}$                                        // gain.  $\mathcal{O}(|y|^3)$  Note you probably don't want to compute  $S^{-1}$  explicitly...
8      $m_t = m_t^- + Kz$                                            // updated mean
9      $P_t = (I - KH)P_t^-$                                        // updated covariance
10  end for
11  return  $(m_t, P_t), (m_t^-, P_t^-)$ 
12 end procedure

```

The entire filtering pass through N time steps has complexity $\mathcal{O}(N \cdot (|X|^3 + |y|^3))$.

Gauss–Markov Models

Local structure for univariate Gaussian models

$$p(x(t_{i+1}) \mid X_{1:i}) = \mathcal{N}(x_{i+1}; Ax_i, Q) \quad \text{and} \quad p(x_0) = \mathcal{N}(x_0; m_0, P_0) \quad \text{and} \quad p(y_i \mid X) = \mathcal{N}(y_i; Hx_i, R)$$

```

1 procedure SMOOTHER( $m_0, P_0, A, Q, H, R, y$ )
2    $(m_t, P_t), (m_t^-, P_t^-) \leftarrow \text{FILTER}(m_0, P_0, A, Q, H, R, y)$ 
3   for  $t = T - 1, T - 2, \dots, 1$  do
4      $G_t = P_t A^\top (P_{t+1}^-)^{-1}$ 
5      $m_t^s = m_t + G_t (m_{t+1}^s - m_{t+1}^-)$ 
6      $P_t^s = P_t + G_t (P_{t+1}^s - P_{t+1}^-) G_t^\top$ 
7   end for
8   return  $(m_t^s, P_t^s)$ 
9 end procedure

```

// RTS gain. Complexity $\mathcal{O}(|X|^3)$

// smoothed mean

// smoothed covariance

Time Series:

- ▶ **Markov Chains** formalize the notion of a stochastic process with a *local finite memory*
- ▶ Inference over Markov Chains separates into three operations, that can be performed in *linear* time.
- ▶ If all relationships are *linear* and *Gaussian*,

$$p(x(t_i) \mid x(t_{i-1})) = \mathcal{N}(x_i; Ax_{i-1}, Q) \quad p(y_t \mid x_t) = \mathcal{N}(y_t; Hx_t, R)$$

then inference is analytic and given by the **Kalman Filter** and the **Rauch–Tung–Striebel Smoother**.



Question 1:

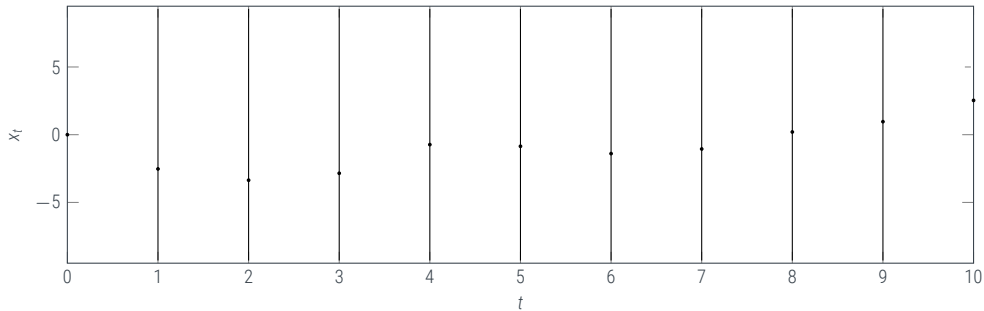
Is there a continuous time limit?



$$p(x_{i+1} \mid x_i) \sim \mathcal{N}(x_{i+1}; x_i, Q)$$

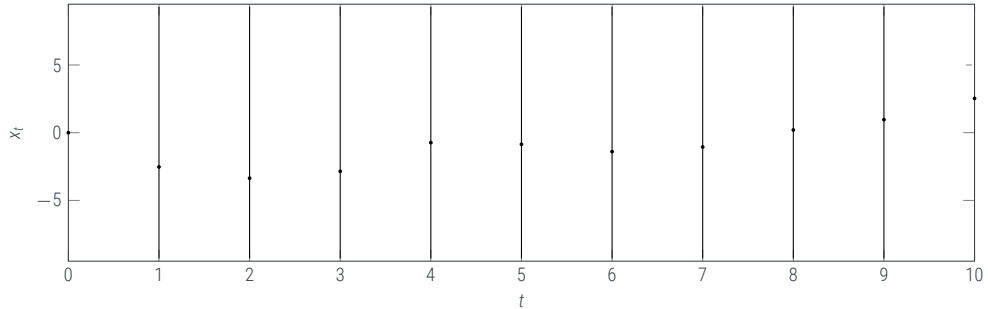


$$p(x_{i+1} | x_i) \sim \mathcal{N}(x_{i+1}; x_i, Q)$$





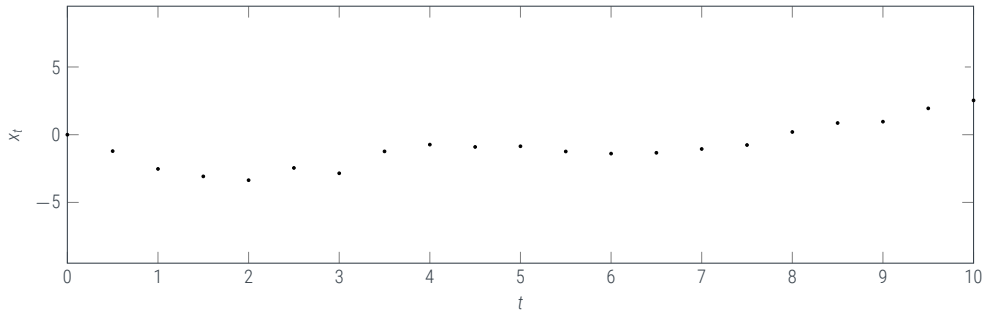
$$p(x(t + \delta t) \mid x(t)) \sim \mathcal{N}(x(t + \delta t); x(t), Q)$$



$$\delta t = 1 \quad Q = 1$$

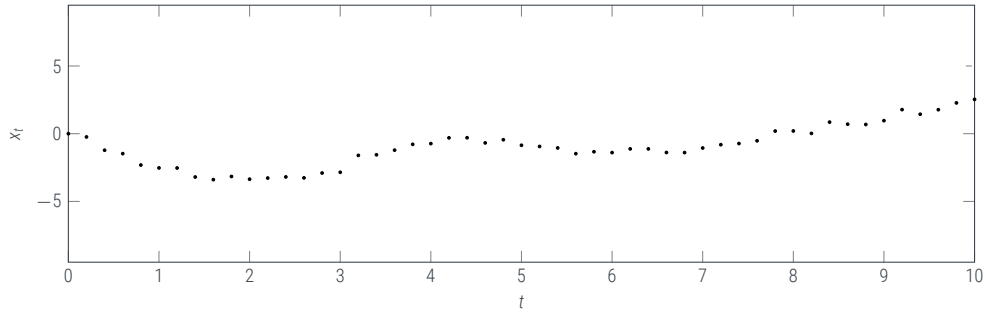


$$p(x(t + \delta t) \mid x(t)) \sim \mathcal{N}(x(t + \delta t); x(t), Q)$$



$$\delta t = 1/2 \quad Q = 1/2$$

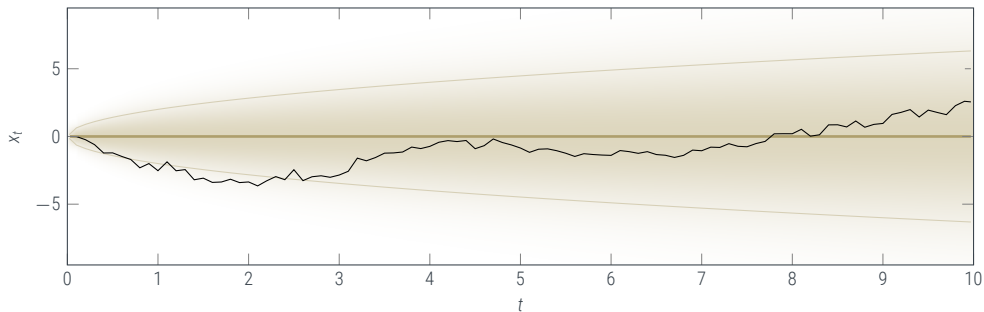
$$p(x(t + \delta t) \mid x(t)) \sim \mathcal{N}(x(t + \delta t); x(t), Q)$$



$$\delta t = 1/4 \quad Q = \delta t$$



$$p(x(t + \delta t) | x(t)) \sim \mathcal{N}(x(t + \delta t); x(t), Q)$$



$$\delta t \rightarrow 0 \quad Q_{\delta t} = ???$$

The Wiener process: One specific limit of one specific discrete model

A different way to write things

$$x(t + \delta t) = x(t) + \Delta\omega(t), \quad \text{with not-really-defined } \Delta\omega,$$

The Wiener process: One specific limit of one specific discrete model

A different way to write things

$$\begin{aligned} x(t + \delta t) &= x(t) + \Delta\omega(t), & \text{with not-really-defined } \Delta\omega, \\ \Leftrightarrow \frac{x(t + \delta t) - x(t)}{\delta t} &= \frac{\Delta\omega(t)}{\delta t}. \end{aligned}$$

The Wiener process: One specific limit of one specific discrete model

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What about the limits?

The Wiener process: One specific limit of one specific discrete model

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What about the limits?

Continuous Time

The Wiener process: One specific limit of one specific discrete model

A different way to write things

$$x(t + \delta t) = x(t) + \Delta\omega(t), \quad \text{with not-really-defined } \Delta\omega,$$

$$\Leftrightarrow \frac{x(t + \delta t) - x(t)}{\delta t} = \frac{\Delta\omega(t)}{\delta t}.$$

What about the limits?

$$\lim_{\delta t \rightarrow 0} \frac{x(t + \delta t) - x(t)}{\delta t} = \frac{dx(t)}{dt},$$

The Wiener process: One specific limit of one specific discrete model

A different way to write things

$$\begin{aligned} x(t + \delta t) &= x(t) + \Delta\omega(t), & \text{with not-really-defined } \Delta\omega, \\ \Leftrightarrow \frac{x(t + \delta t) - x(t)}{\delta t} &= \frac{\Delta\omega(t)}{\delta t}. \end{aligned}$$

What about the limits?

$$\begin{aligned} \lim_{\delta t \rightarrow 0} \frac{x(t + \delta t) - x(t)}{\delta t} &= \frac{dx(t)}{dt}, \\ \lim_{\delta t \rightarrow 0} \frac{\Delta\omega(t)}{\delta t} &= ??? \end{aligned}$$

Continuous Time

The Wiener process: One specific limit of one specific discrete model

A different way to write things

$$\begin{aligned}
 x(t + \delta t) &= x(t) + \Delta\omega(t), & \text{with not-really-defined } \Delta\omega, \\
 \Leftrightarrow \quad \frac{x(t + \delta t) - x(t)}{\delta t} &= \frac{\Delta\omega(t)}{\delta t}.
 \end{aligned}$$

What about the limits?

$$\begin{aligned}
 \lim_{\delta t \rightarrow 0} \frac{x(t + \delta t) - x(t)}{\delta t} &= \frac{dx(t)}{dt}, \\
 \lim_{\delta t \rightarrow 0} \frac{\Delta\omega(t)}{\delta t} &= w(t), \quad \text{where } w(t) \sim \mathcal{N}(0, 1). \quad (\text{"weak" derivative})
 \end{aligned}$$

The Wiener process: One specific limit of one specific discrete model

A different way to write things

$$\begin{aligned} x(t + \delta t) &= x(t) + \Delta\omega(t), & \text{with not-really-defined } \Delta\omega, \\ \Leftrightarrow \frac{x(t + \delta t) - x(t)}{\delta t} &= \frac{\Delta\omega(t)}{\delta t}. \end{aligned}$$

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This is one of the key properties of the *Wiener process* (aka. *Brownian motion*).

Note that: *This is not a proper definition of the Wiener process!* This would go beyond the scope of this course. See [Särkkä & Solin, *Applied Stochastic Differential Equations*, 2019] for a thorough introduction.

For our purposes the (linear, time-invariant) **Stochastic Differential Equation (SDE)**

$$dx(t) = Fx(t) dt + L d\omega(t),$$

together with $x(0) = x_0$, describes the local behaviour of the (unique) Gaussian process with

$$\mathbb{E}(x(t)) =: m(t) = e^{Ft}x_0 \quad \text{cov}(x(t), x(t')) = \int_{\min(t, t')}^{\max(t, t')} e^{F(\max(t, t') - \tau)} LL^T e^{F^T(\max(t, t') - \tau)} d\tau$$

This GP is known as the **solution** of the SDE. It gives rise to the discrete-time stochastic recurrence relation $p(x(t_{i+1}) \mid x(t_i)) = \mathcal{N}(x(t_{i+1}); A_i x(t_i), Q_i)$ with $(\Delta t_i := t_{i+1} - t_i)$

$$A_i = e^{F\Delta t_i} \quad \text{and} \quad Q_i = \int_0^{\Delta t_i} e^{F(\Delta t_i - \tau)} LL^T e^{F^T(\Delta t_i - \tau)} d\tau.$$

Matrix exponential: $e^X := \sum_{i=0}^{\infty} \frac{X^i}{i!}$. Thus : $e^0 = I$, $(e^X)^{-1} = e^{-X}$, $X = VDV^{-1} \Rightarrow Ve^DV^{-1}$, $e^{\text{diag}_i d_i} = \text{diag}_i e^{d_i}$, $\det e^X = e^{\text{tr} X}$.

What this means:

- ▶ LTI-SDEs have a correspondence to GPs (so-called *Gauss–Markov processes*)
- ▶ LTI-SDEs can be discretized *exactly* to get discrete, linear Gaussian transition models

⇒ Gauss–Markov process inference can be done *in linear time* via filtering and smoothing!

$$dx(t) = Fx(t) dt + L d\omega_t$$

$$\mathbb{E}(x(t)) =: m(t) = e^{Ft}x_0 \quad \text{cov}(x(t), x(t')) =: k(t, t') = \int_{\min(t, t')}^{\max(t, t')} e^{F(\max(t, t') - \tau)} LL^T e^{F^T(\max(t, t') - \tau)} d\tau$$

The Connection to GPs

Some well-studied examples

$$dx(t) = Fx(t) dt + L d\omega_t$$

$$\mathbb{E}(x(t)) =: m(t) = e^{Ft} x_0 \quad \text{cov}(x(t), x(t')) =: k(t, t') = \int_{\min(t, t')}^{\max(t, t')} e^{F(\max(t, t') - \tau)} L L^T e^{F^T(\max(t, t') - \tau)} d\tau$$

The (scaled) Wiener process

$$\begin{aligned}
 F = 0, L = \theta & \Rightarrow & m(t) &= x_0 & k(t, t') &= \theta^2 \min(t, t') \\
 A_i &= I & Q_i &= \theta^2 (t_{i+1} - t_i)
 \end{aligned}$$

The Connection to GPs

Some well-studied examples

$$dx(t) = Fx(t) dt + L d\omega_t$$

$$\mathbb{E}(x(t)) =: m(t) = e^{Ft} x_0 \quad \text{cov}(x(t), x(t')) =: k(t, t') = \int_{\min(t, t')}^{\max(t, t')} e^{F(\max(t, t') - \tau)} L L^\top e^{F^\top (\max(t, t') - \tau)} d\tau$$

The Ornstein-Uhlenbeck process

$$F = -\frac{1}{\lambda}, L = \frac{2\theta}{\sqrt{\lambda}} \quad \Rightarrow \quad \begin{aligned} m(t) &= x_0 e^{-\frac{t}{\lambda}} & k(t, t') &= \theta^2 \left(e^{-\frac{|t-t'|}{\lambda}} - e^{-\frac{t+t'}{\lambda}} \right) \\ A_i &= e^{-\Delta t_i / \lambda} & Q_i &= \theta^2 \left(1 - e^{-2\Delta t_i / \lambda} \right) \end{aligned}$$

Non-Scalar State-Space Models

Integrators and Polynomial Splines

$$dx(t) = Fx(t) dt + L d\omega_t$$

- So far, we have seen examples with $x(t) \in \mathbb{R}$.
- But F and L can also be matrices. Consider the example

$$x = \begin{bmatrix} x_{(1)} \\ x_{(2)} \end{bmatrix} \quad F = \begin{bmatrix} 0 & 1 \\ 0 & 0 \end{bmatrix} \quad L = \begin{bmatrix} 0 \\ 1 \end{bmatrix}$$

That is:

$$\begin{bmatrix} dx_{(1)}(t) \\ dx_{(2)}(t) \end{bmatrix} = \begin{bmatrix} x_{(2)}(t) dt + 0 d\omega \\ 0 dt + d\omega \end{bmatrix} \Rightarrow x_{(1)}(t) = \int_0^t x_{(2)}(t) dt + [x_0]_1$$

Non-Scalar State-Space Models

Integrators and Polynomial Splines

$$dx(t) = Fx(t) dt + L d\omega_t$$

The Wiener velocity (aka. "once-integrated Wiener process")

$$F = \begin{bmatrix} 0 & 1 \\ 0 & 0 \end{bmatrix}, L = \begin{bmatrix} 0 \\ 1 \end{bmatrix} \quad \Rightarrow \quad m(t) = e^{Ft} x_0, \quad k(t, t') = \frac{\min^3(t, t')}{3} + |t - t'| \frac{\min^2(t, t')}{2}$$

$$A_i = \begin{bmatrix} 1 & \Delta t_i \\ 0 & 1 \end{bmatrix}, \quad Q_i = \begin{bmatrix} \frac{\Delta t_i^3}{3} & \frac{\Delta t_i^2}{2} \\ \frac{\Delta t_i^2}{2} & \Delta t_i \end{bmatrix}$$

Q1 summary:

- ▶ Certain Gaussian processes can be written as LTI-SDEs
 - ▶ (integrated) Wiener process
 - ▶ (integrated) Ornstein–Uhlenbeck process
 - ▶ Matern processes
 - ▶ Even the square-exponential kernel can be approximated by an LTI-SDE
- ▶ LTI-SDEs can be discretized *exactly* to get discrete, linear Gaussian transition models
- ▶ Inference in linear Gauss–Markov models (and thus in Gauss–Markov processes) can be done *in linear time* via filtering and smoothing



Question 2:
Can we learn the model?

Recap (Lecture 10) Learning the Kernel for GPs

Bayesian Hierarchical Inference

$$p(f | \theta) = \mathcal{GP}(f; m_\theta, k_\theta) \quad \text{e.g.} \quad m_\theta(\bullet) = \phi(\bullet)^\top \theta, \text{ or } k_\theta(\bullet, \circ) = \theta_1 \exp\left(-\frac{(\bullet - \circ)^2}{2\theta_2^2}\right).$$

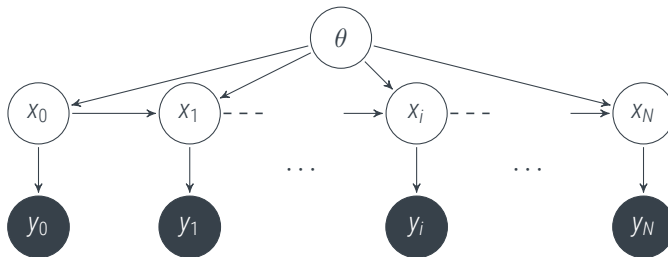
- The **evidence** in Bayes' theorem is the **marginal likelihood for the model**

$$p(f | y, x, \theta) = \frac{p(y | f, x, \theta)p(f |, \theta)}{\int p(y | f, x, \theta)p(f |, \theta) df} = \frac{p(y | f, x, \theta)p(f |, \theta)}{p(y | x, \theta)}$$

- For Gaussians *and* Gaussian processes, the evidence has **analytic form**:

$$\underbrace{\mathcal{N}(y; \phi_X^{\theta^\top} w + b, \Lambda)}_{p(y|x, \theta)} \cdot \underbrace{\mathcal{N}(w, \mu, \Sigma)}_{p(f)} = \underbrace{\mathcal{N}(w; m_{\text{post}}^\theta, V_{\text{post}}^\theta)}_{p(f|y, x, \theta)} \cdot \underbrace{\mathcal{N}(y; \phi_X^{\theta^\top} \mu + b, \phi_X^{\theta^\top} \Sigma \phi_X^\theta + \Lambda)}_{p(y|\theta, x)}$$

$$\mathcal{N}(y; f^\theta(X), \Lambda^\theta) \cdot \mathcal{GP}(f, \mu^\theta, k^\theta) = \mathcal{GP}(f; m_{\text{post}}^\theta, V_{\text{post}}^\theta) \cdot \mathcal{N}(y; \mu^\theta(X), \Lambda^\theta + k^\theta(X, X))$$



For Gauss–Markov Models, is there a way to compute the model evidence in $\mathcal{O}(N)$?

- Assume unknown **model hyper-parameters** θ (define $x_{-1} = \emptyset$):

$$p(\mathbf{y}, \mathbf{x}, \theta) = p(\theta) \cdot p(\mathbf{y}, \mathbf{x} \mid \theta) = p(\theta) \prod_{i=0}^N p(x_i \mid x_{i-1}, \theta) p(y_i \mid x_i, \theta)$$

- to learn θ , we need the **evidence** (aka. **marginal/type-II likelihood**) (define $y_{-1} = \emptyset$)

$$p(\mathbf{y} \mid \theta) = \prod_{i=0}^N p(y_i \mid y_{0:i-1}, \theta)$$

- the terms in the product decompose into local predictions:

$$\begin{aligned} p(y_i \mid y_{0:i-1}, \theta) &= \int p(y_i, x_i \mid y_{0:i-1}, \theta) dx_i &&= \int p(y_i \mid x_i, y_{0:i-1}, \theta) p(x_i \mid y_{0:i-1}, \theta) dx_i \\ &= \int p(y_i \mid x_i, \theta) p(x_i \mid y_{0:i-1}, \theta) dx_i &&\text{which, for linear Gaussian systems, is} \\ &= \int \mathcal{N}(y_i; Hx_i, R) \mathcal{N}(x_i; m_i^-, P_i^-) dx_i &&= \mathcal{N}(y_i; Hm_i^-, HP_i^- H^\top + R) = \mathcal{N}(z_i; 0, S_i) \end{aligned}$$

Learning parameters of a Gauss–Markov Model

For the LTI system

$$p(x_i \mid x_{i-1}, \theta) = \mathcal{N}(x_i; Ax_{i-1}, Q), \quad \text{and} \quad p(y_i \mid x_i, \theta) = \mathcal{N}(y_i; Hx_i, R),$$

the (log) evidence is given by

$$\begin{aligned}
 p(\mathbf{y} \mid \theta) &= \prod_{i=1}^N p(y_i; y_{0:i-1}, \theta) \\
 &= \prod_{i=0}^N \mathcal{N}(y_i; Hm_i^-, HP_i^- H^\top + R) \\
 \log p(\mathbf{y} \mid \theta) &= -\frac{1}{2} \sum_{i=1}^N \left(z_i^\top S_i^{-1} z_i + \log |S_i| + \log 2\pi \right)
 \end{aligned}$$

In principle, this could also be used to learn A, Q, R, H directly, but there's a more elegant way to do this for linear Gaussian systems. For more, cf. Ghahramani & Hinton, 1996, and Särkkä 2013.



Question 3:

What if the world is not linear Gaussian?

Hidden Markov Models

Generalization to non-Gaussian relationships

Name	Distribution	Algorithm
Markovian System:	$p(\mathbf{y}, \mathbf{x}) = \prod_{i=0}^N p(x_i x_{i-1})p(y_i x_i)$	General Bayesian filtering and smoothing
Linear Gaussian System:	$p(\mathbf{y}, \mathbf{x}) = \prod_{i=0}^N \mathcal{N}(x_i; A_i x_{i-1}, Q_i) \mathcal{N}(y_i; H x_i, R)$	Kalman filter, Rauch-Tung-Striebel smoother
Nonlinear Gaussian System:	$p(\mathbf{y}, \mathbf{x}) = \prod_{i=0}^N \mathcal{N}(x_i; a(x_{i-1}), Q_i) \mathcal{N}(y_i; h(x_i), R)$	e.g. Extended/Unscented/Particle filter etc.
Non-Gaussian observations:	$p(\mathbf{y}, \mathbf{x}) = \prod_{i=0}^N \mathcal{N}(x_i; A_i x_{i-1}, Q_i) p(y_i h(x_i))$	
Hidden Markov Model (e.g.):	$p(\mathbf{y}, \mathbf{x}) = \prod_{i=0}^N p(x_i \Pi x_{i-1}) \mathcal{N}(y_i; h(x_i), R)$	

- For continuous systems with nonlinear dynamics and/or non-linear observations, a number of *approximately Gaussian filters* have been developed. For more see, e.g.

- **Simo Särkkä**. *Bayesian Filtering and Smoothing*. Cambridge University Press, 2013

https://users.aalto.fi/~ssarkka/pub/cup_book_online_20131111.pdf

Because streaming data is a common data type, time series are an entire sub-field of their own, studied in a diverse range of domains. There is no time to cover them all in this course.

Summary:

Markov Chains capture **finite memory** of a time series through conditional independence

Gauss–Markov models map this state to linear algebra

Kalman filter is the name for the corresponding algorithm

SDEs (Stochastic Differential Equations) are the continuous-time limit of discrete-time stochastic recurrence relations (in particular, linear SDEs are the continuous-time generalization discrete-time linear Gaussian systems)

Parameters of the model can be learnt by optimizing the (log) evidence, which is also $\mathcal{O}(N)$.

Non-Gaussian models can be learnt by approximate inference, analogous to GP models.

Please cite this course, as

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Gauss–Markov models form the algorithmic scaffold for time-series models .