

# ColumbiaX: Machine Learning

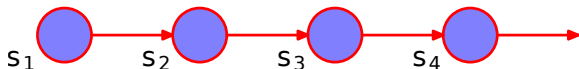
## Lecture 22

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# MARKOV MODELS



The sequence  $(s_1, s_2, s_3, \dots)$  has the *Markov property*, if for all  $t$

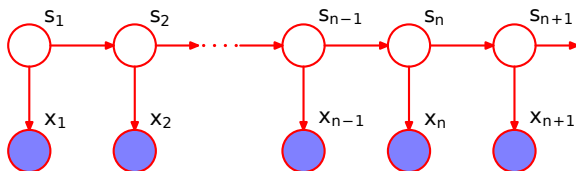
$$p(s_t | s_{t-1}, \dots, s_1) = p(s_t | s_{t-1}).$$

Our first encounter with Markov models assumed a finite state space, meaning we can define an indexing such that  $s \in \{1, \dots, S\}$ .

This allowed us to represent the transition probabilities in a matrix,

$$A_{ij} \quad \Leftrightarrow \quad p(s_t = j | s_{t-1} = i).$$

# HIDDEN MARKOV MODELS



The hidden Markov model modified this by assuming the sequence of states was a *latent process* (i.e., unobserved).

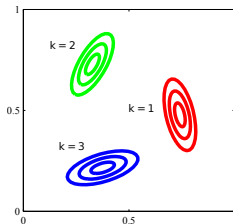
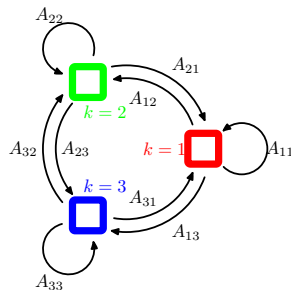
An observation  $x_t$  is associated with each  $s_t$ , where  $x_t \mid s_t \sim p(x \mid \theta_{s_t})$ .

Like a mixture model, this allowed for a few distributions to generate the data. It adds an extra transition rule between distributions.

# DISCRETE STATE SPACES

In both cases, the *state space* was discrete and relatively small in number.

- ▶ For the Markov chain, we gave an example where states correspond to positions in  $\mathbb{R}^d$ .
- ▶ A continuous hidden Markov model might perturb the latent state of the Markov chain.
  - ▶ For example, each  $s_i$  can be modified by continuous-valued noise,  $x_i = s_i + \epsilon_i$ .
  - ▶ But  $s_{1:T}$  is still a *discrete* Markov chain.



# DISCRETE VS CONTINUOUS STATE SPACES

Markov and hidden Markov models both assume a discrete state space.

For Markov models:

- ▶ The state could be a data point  $x_i$  (Markov Chain classifier)
- ▶ The state could be an object (object ranking)
- ▶ The state could be the destination of a link (internet search engines)

For hidden Markov models we can simplify complex data:

- ▶ Sequences of discrete data may come from a few discrete distributions.
- ▶ Sequences of continuous data may come from a few distributions.

What if we model the states as continuous too?

# CONTINUOUS-STATE MARKOV MODEL

Continuous Markov models extend the state space to a continuous domain. Instead of  $s \in \{1, \dots, S\}$ ,  $s$  can take any value in  $\mathbb{R}^d$ .

Again compare:

- ▶ Discrete-state Markov models: The states live in a discrete space.
- ▶ Continuous-state Markov models: The states live in a continuous space.

The simplest example is the process

$$s_t = s_{t-1} + \epsilon_t, \quad \epsilon_t \sim N(0, aI).$$

Each successive state is a perturbed version of the current state.

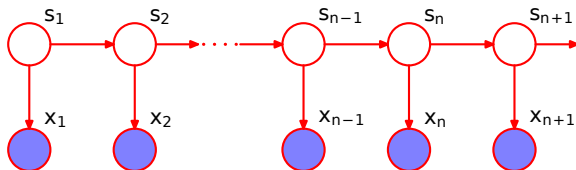
# LINEAR GAUSSIAN MARKOV MODEL

The most basic continuous-state version of the hidden Markov model is called a *linear Gaussian Markov model* (also called the *Kalman filter*).

$$\underbrace{s_t = Cs_{t-1} + \epsilon_{t-1}}_{\text{latent process}}, \quad \underbrace{x_t = Ds_t + \varepsilon_t}_{\text{observed process}}$$

- ▶  $s_t \in \mathbb{R}^p$  is a continuous-state latent (unobserved) Markov process
- ▶  $x_t \in \mathbb{R}^d$  is a continuous-valued observation
- ▶ The process noise  $\epsilon_t \sim N(0, Q)$
- ▶ The measurement noise  $\varepsilon_t \sim N(0, V)$

# EXAMPLE APPLICATIONS



Difference from HMM:  $s_t$  and  $x_t$  are *both* from continuous distributions.

The linear Gaussian Markov model (and its variants) has many applications.

- ▶ Tracking moving objects
- ▶ Automatic control systems
- ▶ Economics and finance (e.g., stock modeling)
- ▶ etc.



## EXAMPLE: TRACKING

We get (very) noisy measurements of an object's position in time,  $x_t \in \mathbb{R}^2$ .

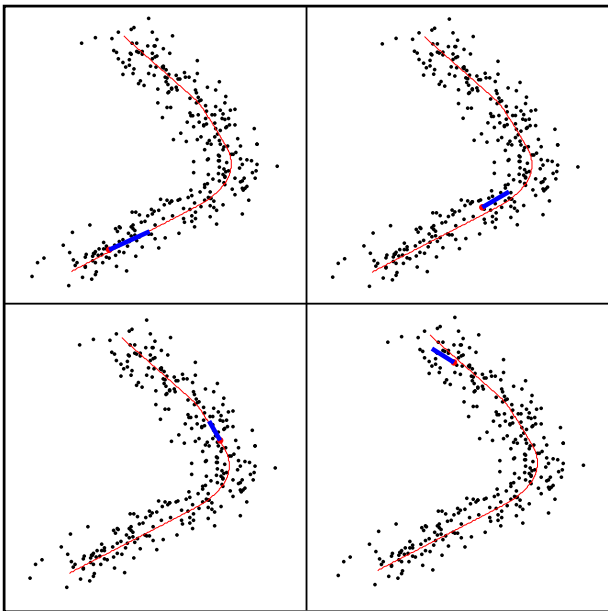
The time-varying state vector is  $s = [\text{pos}_1 \text{ vel}_1 \text{ accel}_1 \text{ pos}_2 \text{ vel}_2 \text{ accel}_2]^T$ .

Motivated by the underlying physics, we model this as:

$$s_{t+1} = \underbrace{\begin{bmatrix} 1 & \Delta t & \frac{1}{2}(\Delta t)^2 & 0 & 0 & 0 \\ 0 & 1 & \Delta t & 0 & 0 & 0 \\ 0 & 0 & e^{-\alpha\Delta t} & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 & \Delta t & \frac{1}{2}(\Delta t)^2 \\ 0 & 0 & 0 & 0 & 1 & \Delta t \\ 0 & 0 & 0 & 0 & 0 & e^{-\alpha\Delta t} \end{bmatrix}}_{\equiv C} s_t + \epsilon_t$$
$$x_{t+1} = \underbrace{\begin{bmatrix} 1 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 & 0 & 0 \end{bmatrix}}_{\equiv D} s_{t+1} + \varepsilon_{t+1}$$

Therefore,  $s_t$  not only approximates where the target is, but where it's going.

## EXAMPLE: TRACKING



# THE LEARNING PROBLEM

As with the hidden Markov model, we're given the sequence  $(x_1, x_2, x_3, \dots)$ , where each  $x \in \mathbb{R}^d$ . The goal is to learn state sequence  $(s_1, s_2, s_3, \dots)$ .

All distributions are Gaussian,

$$p(s_{t+1} = s | s_t) = N(Cs_t, Q), \quad p(x_t = x | s_t) = N(Ds_t, V).$$

Notice that with the discrete HMM we wanted to learn  $\pi$ ,  $A$  and  $B$ , where

- ▶  $\pi$  is the initial state distribution
- ▶  $A$  is the transition matrix among the discrete set of states
- ▶  $B$  contains the state-dependent distributions on discrete-valued data

The situation here is very different.

# THE LEARNING PROBLEM

No “B” to learn: In the linear Gaussian Markov model, each state is unique and so the distribution on  $x_t$  is different for each  $t$ .

No “A” to learn: In addition, each state transition is to a brand new state, so each  $s_t$  has its own unique probability distribution.

What we can learn are the two posterior distributions.

1.  $p(s_t|x_1, \dots, x_t)$  : A distribution on the current state given the past.
  2.  $p(s_t|x_1, \dots, x_T)$  : A distribution on each latent state in the sequence
- ▶ #1: Kalman *filtering* problem. We'll focus on this one today.
  - ▶ #2: Kalman *smoothing* problem. Requires extra step (not discussed).

# THE KALMAN FILTER

**Goal:** Learn the sequence of distributions  $p(s_t|x_1, \dots, x_t)$  given a sequence of data  $(x_1, x_2, x_3, \dots)$  and the model

$$s_{t+1} | s_t \sim N(Cs_t, Q), \quad x_t | s_t \sim N(Ds_t, V).$$

This is the (linear) Kalman filtering problem and is often used for tracking.

**Setup:** We can use Bayes rule to write

$$p(s_t|x_1, \dots, x_t) \propto p(x_t|s_t) p(s_t|x_1, \dots, x_{t-1})$$

and represent the prior as a marginal distribution

$$p(s_t|x_1, \dots, x_{t-1}) = \int p(s_t|s_{t-1}) p(s_{t-1}|x_1, \dots, x_{t-1}) ds_{t-1}$$

# THE KALMAN FILTER

We've decomposed the problem into parts that we do and don't know (yet)

$$p(s_t|x_1, \dots, x_t) \propto \underbrace{p(x_t|s_t)}_{N(Ds_t, V)} \int \underbrace{p(s_t|s_{t-1})}_{N(Cs_{t-1}, Q)} \underbrace{p(s_{t-1}|x_1, \dots, x_{t-1})}_{?} ds_{t-1}$$

Observations and considerations:

1. The left is the posterior on  $s_t$  and the right has the posterior on  $s_{t-1}$ .
2. We want the integral to be in closed form and a known distribution.
3. We want the prior and likelihood terms to lead to a known posterior.
4. We want future calculations, e.g. for  $s_{t+1}$ , to be easy.

We will see how choosing the Gaussian distribution makes this all work.

# THE KALMAN FILTER: STEP 1

## Calculate the marginal for prior distribution

Hypothesize (temporarily) that the unknown distribution is Gaussian,

$$p(s_t|x_1, \dots, x_t) \propto \underbrace{p(x_t|s_t)}_{N(Ds_t, V)} \int \underbrace{p(s_t|s_{t-1})}_{N(Cs_{t-1}, Q)} \underbrace{p(s_{t-1}|x_1, \dots, x_{t-1})}_{N(\mu, \Sigma) \text{ by hypothesis}} ds_{t-1}$$

A property of the Gaussian is that marginals are still Gaussian,

$$\int N(s_t|Cs_{t-1}, Q)N(s_{t-1}|\mu, \Sigma)ds_{t-1} = N(s_t|C\mu, Q + C\Sigma C^T).$$

We know  $C$  and  $Q$  (by design) and  $\mu$  and  $\Sigma$  (by hypothesis).

# THE KALMAN FILTER: STEP 2

## Calculate the posterior

We plug in the marginal distribution for the prior and see that

$$p(s_t|x_1, \dots, x_t) \propto N(x_t|Ds_t, V) N(s_t|C\mu, Q + C\Sigma C^T).$$

Though the parameters look complicated, the posterior is just a Gaussian

$$p(s_t|x_1, \dots, x_t) = N(s_t|\mu', \Sigma')$$

$$\Sigma' = [(Q + C\Sigma C^T)^{-1} + D^T V^{-1} D]^{-1}$$

$$\mu' = \Sigma' (D^T V^{-1} x_t + (Q + C\Sigma C^T)^{-1} C\mu)$$

We can plug the relevant values into these two equations.



# ADDRESSING THE GAUSSIAN ASSUMPTION

By making the assumption of a Gaussian in the prior,

$$p(s_t|x_1, \dots, x_t) \propto \underbrace{p(x_t|s_t)}_{N(x_t|Ds_t, V)} \int \underbrace{p(s_t|s_{t-1})}_{N(s_t|Cs_{t-1}, Q)} \underbrace{p(s_{t-1}|x_1, \dots, x_{t-1})}_{N(\mu, \Sigma) \text{ by hypothesis}} ds_{t-1}$$

we found that the posterior is also Gaussian with a new mean and covariance.

- We therefore only need to define a Gaussian prior on the first state to keep things moving forward. For example,

$$p(s_0) \sim N(0, I).$$

Once this is done, all future calculations are in closed form.

# KALMAN FILTER: ONE FINAL QUANTITY

## Making predictions

We know how to update the sequence of state posterior distributions

$$p(s_t | x_1, \dots, x_t).$$

What about predicting  $x_{t+1}$ ?

$$\begin{aligned} p(x_{t+1} | x_1, \dots, x_t) &= \int p(x_{t+1} | s_{t+1}) p(s_{t+1} | x_1, \dots, x_t) ds_{t+1} \\ &= \int \underbrace{p(x_{t+1} | s_{t+1})}_{N(x_{t+1} | Ds_{t+1}, V)} \int \underbrace{p(s_{t+1} | s_t)}_{N(s_{t+1} | Cs_t, Q)} \underbrace{p(s_t | x_1, \dots, x_t)}_{N(s_t | \mu', \Sigma')} ds_t ds_{t+1} \end{aligned}$$

Again, Gaussians are nice because these operations stay Gaussian.

This is a multivariate Gaussian that looks even more complicated than the previous one (omitted). Simply perform the previous integral twice.

# ALGORITHM: KALMAN FILTERING

The Kalman filtering algorithm can be run in real time.

0. Set the initial state distribution  $p(s_0) = N(0, I)$

1. Prior to observing each new  $x_t \in \mathbb{R}^d$  predict

$$x_t \sim N(\mu_t^x, \Sigma_t^x) \quad (\text{using previously discussed marginalization})$$

2. After observing each new  $x_t \in \mathbb{R}^d$  update

$$p(s_t | x_1, \dots, x_t) = N(\mu_t^s, \Sigma_t^s) \quad (\text{using equations on previous slide})$$

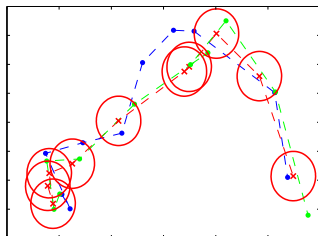
# EXAMPLE

## Learning state trajectory

Green: True trajectory

Blue: Observed trajectory

Red: State distribution



Intuitions about what this is doing:

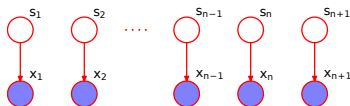
- In the prior distribution notice that we add  $Q$  to the covariance,

$$p(s_t|x_1, \dots, x_{t-1}) = N(s_t|C\mu, Q + C\Sigma C^T).$$

This allows the state  $s_t$  to “drift” away from  $s_{t-1}$ .

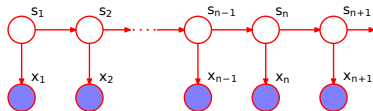
- In the posterior  $p(s_t|x_1, \dots, x_t)$ ,  $x_t$  “pulls” the distribution away.

# SOME FINAL MODEL COMPARISONS



## Gaussian mixture model

- ▶  $s_t \sim \text{Discrete}(\pi)$
- ▶  $x_t | s_t \sim N(\mu_{s_t}, \Sigma_{s_t})$

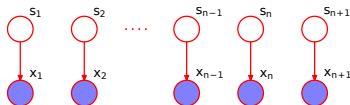


## Continuous hidden Markov model

- ▶  $s_t | s_{t-1} \sim \text{Discrete}(A_{s_{t-1}})$
- ▶  $x_t | s_t \sim N(\mu_{s_t}, \Sigma_{s_t})$

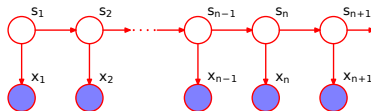
We saw how the transition from GMM  $\rightarrow$  HMM involves using a Markov chain to index the distribution on clusters.

# SOME FINAL MODEL COMPARISONS



## Probabilistic PCA

- ▶  $s_t \sim N(0, Q)$
- ▶  $x_t | s_t \sim N(Ds_t, V)$



## Linear Gaussian Markov model

- ▶  $s_t | s_{t-1} \sim N(Cs_{t-1}, Q)$
- ▶  $x_t | s_t \sim N(Ds_t, V)$

There is a similar relationship between probabilistic PCA and the Kalman filter. (Probabilistic PCA also learns  $D$ , while the Kalman filter doesn't).

# EXTENSIONS

There are a variety of extensions to this framework. The equations in the corresponding algorithms would all look familiar given our discussion.

**Extended Kalman filter:** *Nonlinear Kalman filters* use nonlinear function of the state,  $h(s_t)$ . The EKF approximates  $h(s_t) \approx h(z) + \nabla h(z)(s_t - z)$

$$s_{t+1} \mid s_t \sim N(Ds_t, Q), \quad x_t \mid s_t \sim N(h(s_t), V).$$

**Continuous time:** Sometimes the time between observations varies. Let  $\Delta_t$  be the time between observation  $x_t$  and  $x_{t+1}$ , then model

$$s_{t+1} \mid s_t \sim N(s_t, \Delta_t Q), \quad x_t \mid s_t \sim N(Ds_t, V).$$

**Adding control:** In dynamic models, we can add control to the state using a vector  $u_t$  whose values we choose (e.g., thrusters).

$$s_{t+1} \mid s_t \sim N(Cs_t + Gu_t, Q), \quad x_t \mid s_t \sim N(Ds_t, V).$$