## Lecture 18:

SEQUENTIAL STATE ESTIMATION IN NONLINEAR, NON GaUSSIAN DYNAMICAL SYSTEMS

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Overview of filtering methods

|  |  |  |  |
| :--- | :--- | :--- | :--- |
| Dynamics | Observation | Belief | Method |
| Trans. mat | Any | Histogram | HMM/ forwards |
| DBN | DBN | Jtree | Thin Junction tree filter (TJTF) |
| DBN | DBN | Prod. histo | Boyen-Koller (BK) filter |
| Lin Gauss | Lin Gauss | Gauss | Kalman filter (KF) |
| NonLin Gauss | NonLin Gauss | Gauss | Extended Kalman filter (EKF) |
| NonLin Gauss | NonLin Gauss | Gauss | Unscented Kalman filter (UKF) |
| Any | Any | Gauss | Assumed Density Filtering (ADF) |
| Mix Lin Gauss | Mix Lin Gauss Mix Gauss | Assumed Density Filtering (ADF) |  |
| Any | Any | Samples | Particle filtering |

- A generic state-space model defines the dynamics $P\left(X_{t} \mid X_{t-1}\right)$ and the observation model $P\left(y_{t} \mid X_{t}\right)$.

- Online inference (filtering) means recursively computing the belief state $P\left(X_{t} \mid y_{1: t}\right)$ using Bayes' rule:

$$
P\left(X_{t} \mid y_{1: t}\right)=\frac{1}{P\left(y_{t} \mid y_{1: t-1}\right)} P\left(y_{t} \mid X_{t}\right) \int P\left(X_{t} \mid x_{t-1}\right) P\left(x_{t-1} \mid y_{1: t-1}\right) d x_{t-1}
$$

- If $X_{t}=\theta$ is a constant, this can be used for recursive parameter estimation.
- There are many different methods, depending on how we represent the dynamical model, the observation model and the belief state.

- Represent belief state as a histogram: $P\left(X_{t}=i \mid y_{1: t}\right)=\alpha_{t}(i)$.
- Predict step:

$$
P\left(X_{t+1}=j \mid y_{1: t}\right)=\sum_{i} P\left(X_{t+1}=j \mid X_{t}=i\right) P\left(X_{t}=i \mid y_{1: t}\right)
$$

- Update step:

$$
P\left(X_{t+1}=j \mid y_{1: t+1}\right)=\frac{p\left(y_{t+1} \mid X_{t+1}=j\right) P\left(X_{t+1}=j \mid y_{1: t}\right)}{p\left(y_{t+1} \mid y_{1: t}\right)}
$$

- A DBN generalizes a state space model by representing the components of $X_{t}$ and $Y_{t}$ and their conditional (in)dependencies using a graph.
- By factorizing the state space in this way, we can substantially reduce the number of free parameters.
- e.g., let $X_{t}$ be a bit vector of length $K$. An HMM transition matrix would have $O\left(2^{K} \times 2^{K}\right)$ parameters. A DBN may have $O(K)$ parameters, depending on the structure.
- For linear Gaussian models, sparse graphs $\equiv$ sparse matrices, so DBNs are not needed as much (but are still helpful).


## $\underline{\text { Sparse graphs } \nRightarrow \text { SPARSE DISCRETE TRANSITION MATRICES }}$

- Any discrete DBN can be flattened into an HMM.
- But the resulting transition matrix will not necessarily have 0 s in it.

$$
\begin{aligned}
& P\left(Q_{t}^{(1)}=j_{1}, Q_{t}^{(2)}=j_{2} \mid Q_{t-1}^{(1)}=i_{1}, Q_{t-1}^{(2)}=i_{2}\right) \\
& =P\left(Q_{t}^{(1)}=j_{1} \mid Q_{t-1}^{(1)}=i_{1}\right) \times P\left(Q_{t}^{(2)}=j_{2} \mid Q_{t-1}^{(2)}=i_{2}\right)
\end{aligned}
$$

- For discrete state spaces, the graph structure provides a compact representation of the model in a way that cannot be easily captured in the form of the parameter matrices.


Belief state $P\left(X_{t} \mid y_{1: t}\right)$ has size $O\left(2^{N}\right)$ for $N$ binary chains, because the common observed child $Y_{t}$ couples all the parents (explaining away).

## Linear Gaussian DBNs

- For linear Gaussian models, sparse graphs $\equiv$ sparse matrices, so DBNs are not needed as much (but are still helpful).
- Consider a Vector Auto Regressive process of order 2:
$X_{t}=A_{1} X_{t-1}+A_{2} X_{t-2}+\epsilon_{t}$ where $\epsilon_{t} \sim \mathcal{N}(0, \Sigma)$.
- If $A_{k}(i, j)=0$, then the directed arc $X_{t-k}(i) \rightarrow X_{t}(j)$ is missing (for $k=1,2$ ).
- If $\Sigma^{-1}(i, j)=0$, then the undirected arc $X_{t}(i)-X_{t}(j)$ is missing.



Belief state for coupled HMMs
Even though CHMMs are sparse, all nodes eventually become correlated, so $P\left(X_{t} \mid y_{1: t}\right)$ has size $O\left(2^{N}\right)$.


## Assumed density filtering (ADF)

- ADF forces the belief state to live in some restricted family $\mathcal{F}$, e.g., product of histograms, Gaussian.
- Given a prior $\tilde{\alpha}_{t-1} \in \mathcal{F}$, do one step of exact Bayesian updating to get $\hat{\alpha}_{t} \notin \mathcal{F}$. Then do a projection step to find the closest approximation in the family:

$$
\tilde{\alpha}_{t}=\arg \min _{q \in \mathcal{F}} D\left(\hat{\alpha}_{t} \| q\right)
$$

- If $\mathcal{F}$ is the exponential family, we can solve the KL minimization by moment matching.
exact
approx
- The BK algorithm is ADF applied to a DBN.
- The simplest approximation is to let $\mathcal{F}$ be a product of marginals:

$$
\alpha_{t} \approx \tilde{\alpha}_{t}=\prod_{i=1}^{N} P\left(X_{t}^{i} \mid y_{1: t}\right)
$$

- We start with a prior that is fully factored, $\prod_{i} P\left(X_{t-1}^{i} \mid y_{1: t-1}\right)$.
- We do one step of exact updating using a 2-slice junction tree. This will couple some nodes, but not as many as in the $T$-slice (unrolled) jtree.
- Then we compute the marginals $P\left(X_{t}^{i} \mid y_{1: t}\right)$ (projection step).


## SLAM (SImUlTANEOUS LOCALIZATION AND MAPPING)

- State is location of robot and landmarks $X_{t}=\left(R_{t}, L_{t}^{1: N}\right)$
- Measure location of subset of landmarks at each time step.
- Assume everything is linear Gaussian.
- Use Kalman filter to solve optimally.


Thin Junction tree filter (TJTF)

- The BK algorithm is ADF applied to a DBN.
- The approximating family $\mathcal{F}$ is chosen a priori.
- Above we considered products of marginals, but BK also considered products of overlapping cliques (i.e., junction tree).
- For some problems, the structure of the approximating family should be chosen dynamically, after seeing the evidence.
- TJTF dynamically thins the junction tree, to keep $\mathcal{F}$ tractable.
- This is useful e.g., for SLAM (simultaneous localization and mapping) problem in mobile robotics.


## Review: Kalman filter

- LDS model: $x_{t}=A x_{t-1}+v_{t}, \quad y_{t}=C x_{t}+w_{t}$
- Time update (prediction step):

$$
x_{t \mid t-1}=A x_{t-1 \mid t-1}, \quad P_{t \mid t-1}=A P_{t-1 \mid t-1} A^{T}+Q, \quad y_{t \mid t-1}=C x_{t \mid t-1}
$$

- Measurement update (correction step):

$$
\begin{aligned}
\tilde{y}_{t} & =y_{t}-\hat{y}_{t \mid t-1}(\text { error } / \text { innovation }) \\
P_{\tilde{y}_{t}} & =C P_{t \mid t-1} C^{T}+R \text { (covariance of error) } \\
P_{x_{t} y_{t}} & =P_{t \mid t-1} C^{T} \text { (cross covariance) } \\
K_{t} & =P_{x_{t} y_{t}} P_{\tilde{y}_{t}}^{-1} \text { (Kalman gain matrix) } \\
x_{t \mid t} & =x_{t \mid t-1}+K_{t}\left(y_{t}-y_{t \mid t-1}\right) \\
P_{t \mid t} & =P_{t \mid t-1}-K_{t} P_{x_{t} y_{t}}^{T}
\end{aligned}
$$

- Let $X_{t} \in \mathbb{R}^{N_{x}}$ and $Y_{t} \in \mathbb{R}^{N_{y}}$.
- Computing $P_{t \mid t-1}=A P_{t-1 \mid t-1} A^{T}+Q$ takes $O\left(N_{x}^{2}\right)$ time, assuming dense $P$ and dense $A$.
- Computing $K_{t}=P_{x_{t} y_{t}} P_{\tilde{y}_{t}}^{-1}$ takes $O\left(N_{y}^{3}\right)$ time.
- So overall time is, in general, $\max \left\{N_{x}^{2}, N_{y}^{3}\right\}$.
- KF uses moment form, ie. mean $\mu$ and covariance $\Sigma$.
- Initial uncertainty means covariance can be infinite $\Sigma_{i i}=\infty$.
- It is therefore common to use the information form, which works in terms of canonical (natural) parameters $S=\Sigma^{-1}$ and $\xi=\Sigma^{-1} \mu$.
- Substituting into KF equations and using the matrix inversion lemma yields:

$$
\begin{aligned}
U_{t} & =\left(S_{t \mid t}+A^{T} Q A\right)^{-1} \\
\xi_{t+1 \mid t+1} & =Q^{-1} A U \xi_{t \mid t}+C^{T} R^{-1} y_{t+1} \\
S_{t+1 \mid t+1} & =Q^{-1}-Q^{-1} A U A^{T} Q^{-1}+C^{T} R^{-1} C
\end{aligned}
$$

- Now initial uncertainty means the precision is zero $\Sigma_{i i}^{-1}=0$.
- But now complexity is $O\left(N_{x}^{3}\right)$.

SLAM: MEasurement (update) step


$$
\begin{aligned}
& P\left(X_{t} \mid y_{0: t}\right)=\mathcal{N}\left(\mu_{t}, \Sigma_{t}\right):
\end{aligned}
$$

Observe, moralize, absorb evidence. Correlates $R$ with active $L$.

(a)

(b)

(c)

(d)

Marginalizing out old robot position correlates all active landmarks with each other.

## Thin junction tree filters for SLAM

- The precision matrix $\Sigma^{-1}$ will eventually become non-zero everywhere (although edge strengths may be small).
- Hence the covariance matrix $\Sigma$ becomes dense.
- Hence Kalman filter will take $O\left(N^{2}\right)$ time.
- Junction tree filtering is exact, and hence has the same complexity as KF .
- However, we can adaptively reduce the size of "fat" cliques; this is similar to pruning weak edges in the graphical model (i.e., set elements of $\Sigma^{-1}$ to 0 ).
- This yields an $O(N)$ algorithm.
- Further approximations yield an $O(1)$ algorithm.
- In general, we have the following time complexities:
- Computing $P_{t \mid t-1}=A P_{t-1 \mid t-1} A^{T}+Q$ takes $O\left(N_{x}^{2}\right)$ time.
- Computing $K_{t}=P_{x_{t} y_{t}} P_{\tilde{y}_{t}}^{-1}$ takes $O\left(N_{y}^{3}\right)$ time.
- Computing $P_{t \mid t}=P_{t \mid t-1}-K_{t} P_{x y}^{T}$ takes $O\left(N_{x}^{2}\right)$ time.
- For SLAM, the landmarks are stationary, so only the $R R$ and $R L$ components of $P_{t \mid t-1}$ need be updated in $O\left(N_{x}\right)$ time.
- Unfortunately, computing $P_{t \mid t}=P_{t \mid t-1}-K_{t} P_{x y}^{T}$ takes $O\left(N_{x}^{2}\right)$ time, since $P_{t \mid t}$ becomes dense.
- Doesn't scale to many landmarks.


## NONLINEAR SYSTEMS

- In robotics and other problems, the motion model and the observation model are often nonlinear:

$$
x_{t}=f\left(x_{t-1}\right)+v_{t}, \quad y_{t}=g\left(x_{t}\right)+w_{t}
$$

- An optimal closed form solution to the filtering problem is no longer possible.
- The nonlinear functions $f$ and $g$ are sometimes represented by neural networks (multi-layer perceptrons or radial basis function networks).
- The parameters of $f, g$ may be learned offline using EM, where we do gradient descent (back propagation) in the M step, c.f. learning a MRF/CRF with hidden nodes.
- Or we may learn the parameters online by adding them to the state space: $\tilde{x}_{t}=\left(x_{t}, \theta\right)$. This makes the problem even more nonlinear.
- The basic idea of the EKF is to linearize $f$ and $g$ using a second order Taylor expansion, and then apply the standard KF.
- i.e., we approximate a stationary nonlinear system with a non-stationary linear system.

$$
\begin{aligned}
x_{t} & =f\left(\hat{x}_{t-1 \mid t-1}\right)+A_{\hat{x}_{t \mid t-1}}\left(x_{t-1}-\hat{x}_{t-1 \mid t-1}\right)+v_{t} \\
y_{t} & =g\left(\hat{x}_{t \mid t-1}\right)+C_{\hat{x}_{t \mid t-1}}\left(x_{t}-\hat{x}_{t \mid t-1}\right)+w_{t}
\end{aligned}
$$

where $\hat{x}_{t \mid t-1}=f\left(\hat{x}_{t-1 \mid t-1}\right)$ and $\left.A_{\hat{x}} \stackrel{\text { def }}{=} \frac{\partial f}{\partial x}\right|_{\hat{x}}$ and $\left.C_{\hat{x}} \stackrel{\text { def }}{=} \frac{\partial g}{\partial x}\right|_{\hat{x}}$.

- The noise covariance ( $Q$ and $R$ ) is not changed, i.e., the additional error due to linearization is not modeled.
- UKF does not require computing derivatives of $f$ or $g$, and is accurate to second order.
- The UKF applies the unscented transform twice, once to compute $P\left(X_{t} \mid y_{1: t-1}\right)$ and once to compute $P\left(X_{t} \mid y_{1: t}\right)$.
- The unscented transform passes the mean $\pm \sigma$ in each dimension, and fits an ellipse to the resulting points.

THE NEED FOR MULTIMODAL BELIEF STATES



Switching LDS (Jump Linear system)


Combination of HMM and LDS.

$$
\begin{aligned}
P\left(X_{t}=x_{t} \mid X_{t-1}=x_{t-1}, S_{t}=i\right) & =\mathcal{N}\left(x_{t} ; A_{i} x_{t-1}, Q_{i}\right) \\
P\left(Y_{t}=y \mid X_{t}=x\right) & =\mathcal{N}(y ; C x, R) \\
P\left(S_{t}=j \mid S_{t-1}=i\right) & =M(i, j)
\end{aligned}
$$

Belief state has $O\left(2^{t}\right)$ Gaussian modes:

$$
P\left(X_{t}, S_{t} \mid y_{1: t}\right)=\sum_{s_{1: t-1}} \int d x_{1: t-1} P\left(X_{1: t}, S_{1: t} \mid y_{1: t}\right)
$$

Generalized pseudo Bayesian $=$ ADF per mode of $S_{t}$.


- Optimal belief state has $O\left(2^{t}\right)$ modes.
- Common to use nearest neighbor approximation.
- For each time slice, can enforce that at most one source causes each observation (maximal matching problem, solvable in $O\left(N^{3}\right)$ time using Hungarian algorithm).
- Correspondence problem also arises in shape matching and stereo vision.


## Importance sampling

- Assume we want to compute $\Phi=E \phi(x)=\int d^{N} x P(x) \phi(x)$.
- Drawing samples from $P(x)$ may be too hard, but we can evaluate $P^{*}(x)$ where $P(x)=P^{*}(x) / Z_{P}$.
- Assume we can sample from a proposal density $Q(x)$ and can evaluate $Q^{*}(x)$, where $Q(x)=Q^{*}(x) / Z_{Q}$.
- We assign each sample an importance weight $w_{r}=P^{*}\left(x^{r}\right) / Q^{*}\left(x^{r}\right)$ and then approximate $\Phi=\frac{\sum_{r} w_{r} \phi\left(x^{r}\right)}{\sum_{r} w_{r}}$.
- Does not work well if $Q(x)$ is small where $\left|\phi(x) P^{*}(x)\right|$ is large; $Q(x)$ should have heavy tails.

- Suppose the target density is $P\left(x_{1: t} \mid y_{1: t}\right)$ and the proposal is $q\left(x_{1: t} \mid y_{1: t}\right)$, so $w_{t}^{i} \propto P\left(x_{1: t}^{i} \mid y_{1: t}\right) / Q\left(x_{1: t}^{i} \mid y_{1: t}\right)$.
- The probability of a sample path can be computed recursively using Bayes' rule:

$$
\begin{aligned}
w_{t}^{i} & \propto \frac{P\left(y_{t} \mid x_{t}^{i}\right) P\left(x_{t}^{i} \mid x_{t-1}^{i}\right) P\left(x_{1: t-1}^{i} \mid y_{1: t-1}\right)}{Q\left(x_{t}^{i} \mid x_{1: t-1}^{i}, y_{1: t}\right) Q\left(x_{1: t-1}^{i} \mid y_{1: t-1}\right)} \\
& =\frac{P\left(y_{t} \mid x_{t}^{i}\right) P\left(x_{t}^{i} \mid x_{t-1}^{i}\right)}{Q\left(x_{t}^{i} \mid x_{1: t-1}^{i}, y_{1: t}^{i}\right)} w_{t-1}^{i} \\
& =\hat{w}_{t}^{i} w_{t-1}^{i}
\end{aligned}
$$

- For online problems, we typically use $Q\left(x_{t} \mid x_{1: t-1}^{i}, y_{1: t}\right)=Q\left(x_{t} \mid x_{t-1}^{i}, y_{1: t}\right)$ so we don't have to store the entire history. Hence

$$
\hat{w}_{t}^{i}=\frac{P\left(y_{t} \mid x_{t}^{i}\right) P\left(x_{t}^{i} \mid x_{t-1}^{i}\right)}{Q\left(x_{t}^{i} \mid x_{t-1}^{i}, y_{1: t}\right)}
$$

## Proposal distribution for PF

- The simplest proposal is to sample from the prior $Q\left(x_{t} \mid x_{t-1}^{i}, y_{1: t}\right)=$ $P\left(X_{t} \mid x_{t-1}^{i}\right)$. In vision, this is called the condensation algorithm.
- Recall that the incremental weight is

$$
\hat{w}_{t}^{i}=\frac{P\left(y_{t} \mid x_{t}^{i}\right) P\left(x_{t}^{i} \mid x_{t-1}^{i}\right)}{Q\left(x_{t}^{i} \mid x_{t-1}^{i}, y_{1: t}\right)}
$$

- So for condensation, $\hat{w}_{t}^{i}=P\left(y_{t} \mid x_{t}^{i}\right)$.
- It is better to look at the evidence before proposing:

$$
q\left(x_{t} \mid x_{t-1}^{i}, y_{t}\right)=P\left(x_{t} \mid x_{t-1}^{i}, y_{t}\right)=\frac{P\left(y_{t} \mid x_{t}\right) P\left(x_{t} \mid x_{t-1}^{i}\right)}{\int d x_{t} P\left(y_{t} \mid x_{t}\right) P\left(x_{t} \mid x_{t-1}^{i}\right)}
$$

- In this case, the incremental weight is the denominator $\hat{w}_{t}^{i}=P\left(y_{t} \mid x_{t-1}^{i}\right)$.
(SISR)
- As time increases, one sample path will turn out to be exponentially more likely than any other, so all the weights except one go to 0 .
- This is called sample impoverishment.
- Whenever the effective number of samples $N_{\text {eff }}=1 / \sum_{i}\left(w_{t}^{i}\right)^{2}$ drops below a threshold, we resample with replacement.
- The resampled weights are set to $1 / N$, since the past weights are reflected in the empirical frequency.
- There are various ways to do the resampling in $O(N)$ time.
- PF is the same as SISR.


## Unscented particle filtering

- Often it is too hard to compute the optimal proposal $P\left(X_{t} \mid x_{t-1}^{i}, y_{1: t}\right)$ exactly.
- But sometimes we can approximate this.
- Consider a nonlinear system with Gaussian process noise and linearGaussian observations:

$$
\begin{aligned}
P\left(X_{t} \mid x_{t-1}^{i}\right) & =\mathcal{N}\left(X_{t} ; f_{t}\left(x_{t-1}^{i}\right), Q_{t}\right) \\
P\left(Y_{t} \mid X_{t}\right) & =\mathcal{N}\left(y_{t} ; C_{t} X_{t}, R_{t}\right)
\end{aligned}
$$

- Then we can compute $Q\left(X_{t} \mid x_{t-1}^{i}, y_{1: t}\right)$ using an EKF/UKF (with a delta function prior on $x_{t-1}^{i}$ ), and sample from this.
- Sampling in high dimensional spaces causes high variance in the estimate.
- RBPF idea: sample some variables $R$, and conditional on that, compute expected value of rest $X$ analytically.
- So-called because of RB theorem, which is based on this identity:

$$
\operatorname{Var}[\tau(X, R)]=\operatorname{Var}[E(\tau(X, R) \mid R)]+E[\operatorname{Var}(\tau(X, R) \mid R)]
$$

- Hence $\operatorname{Var}[E(\tau(X, R) \mid R)] \leq \operatorname{Var}[\tau(X, R)]$, so $\tau^{\prime}(X, R)=E(\tau(X, R) \mid R)$ is a lower variance estimator.
- Key idea: if you always know the robot's location, the posterior over landmarks factorizes, so KF takes $O(N)$ time.
- So sample $R_{1: t}$, and for each particle/ trajectory, run a Kalman filter.


