

Prediction and Estimation, Part II

Class 18, 1 Nov 2012

Recap: An automotive example



- Determine automatically, by only *listening* to a running automobile, if it is:

- idling; or

- Travelling at constant velocity; or

- Accelerating; or

- Decelerating

- Assume (for illustration) that we only record energy level

- (SPL) in the sound

- The SPL is measured once per second

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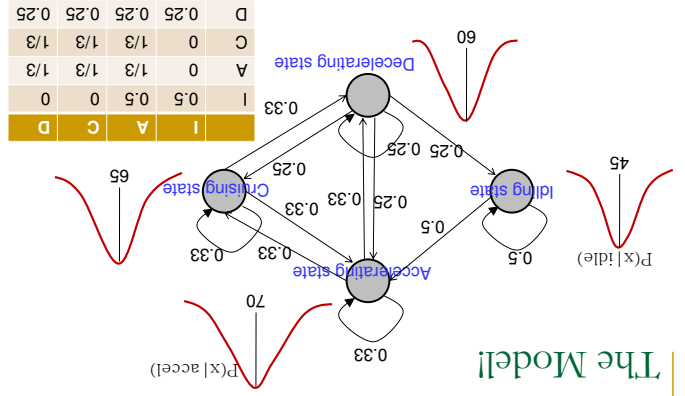
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The Model



- The state-space model

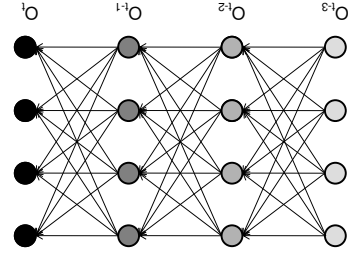
- Assuming all transitions from a state are equally probable

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Prediction with HMMs

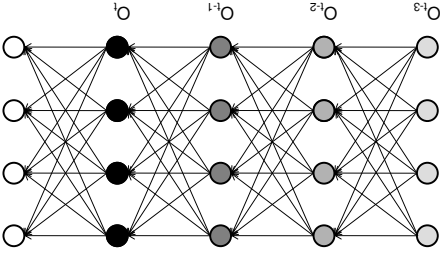


- At t , you have some beliefs for the states

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Prediction with HMMs



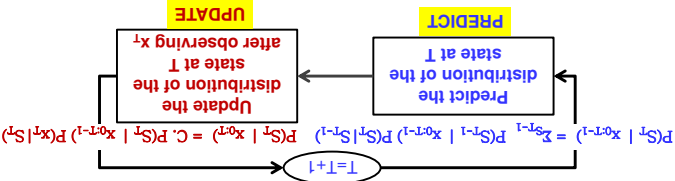
- At t , you have some beliefs for the states
- You predict the beliefs for the state at $t+1$

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Overall procedure



- At $T=0$ the predicted state distribution is the initial state probability

- At each time T , the current estimate of the distribution over states considers *all* observations $x_0 \dots x_T$

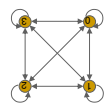
- A natural outcome of the Markov nature of the model

- The prediction+update is identical to the forward computation for HMMs to within a normalizing constant

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Discrete vs. Continuous State Systems

$$P(s'_t | O_{0:t-1}) = \sum_{s_{t-1}} P(s_{t-1} | O_{0:t-1}) P(s'_t | s_{t-1})$$

Prediction at time t

$$P(s'_t | O_{0:t-1}) = \int_{-\infty}^{\infty} P(s_{t-1} | O_{0:t-1}) P(s'_t | s_{t-1}) d s_{t-1}$$

Update after O_t :

$$P(s'_t | O_{0:t}) = C P(s'_t | O_{0:t-1}) P(O_{0:t} | s'_t)$$

$$o'_t = g(s'_t, \gamma'_t)$$

$$s'_t = f(s_{t-1}, \mathcal{E}'_t)$$

- A_t, B_t and Gaussian parameters assumed known
- Probability of observation noise γ is Gaussian
- A linear observation equation
- Sometimes viewed as a driving term μ_ε and additive zero-mean noise
- Probability of state driving term ε is Gaussian
- A linear state dynamics equation

$$o'_t = B'_t s'_t + \gamma'_t$$

$$s'_t = A'_t s_{t-1} + \mathcal{E}'_t$$

$$P(o'_t | \gamma) = \frac{1}{\sqrt{2\pi} \sigma_\gamma} \exp\left(-\frac{1}{2\sigma_\gamma^2} (o'_t - B'_t s'_t)^2\right)$$

$$P(s'_t | \varepsilon) = \frac{1}{\sqrt{2\pi} \sigma_\varepsilon} \exp\left(-\frac{1}{2\sigma_\varepsilon^2} (s'_t - A'_t s_{t-1})^2\right)$$

Special case: Linear Gaussian model

- The state is a continuous valued parameter that is not directly seen
- The state is the position of navlab or the star
- The observations are dependent on the state and are the only way of knowing about the state
- Sensor readings (for navlab) or recorded image (for the telescope)

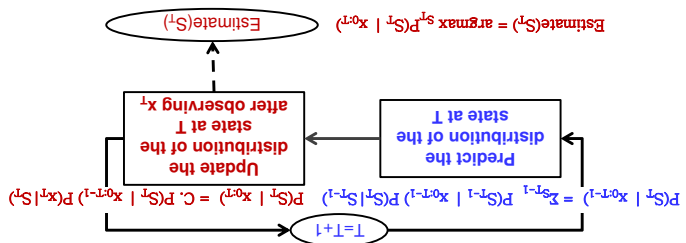


$$o'_t = g(s'_t, \gamma'_t)$$

$$s'_t = f(s_{t-1}, \mathcal{E}'_t)$$

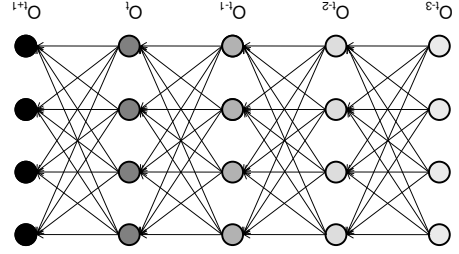
Continuous state system

- The state is estimated from the updated distribution
- The updated distribution is propagated into time, not the state



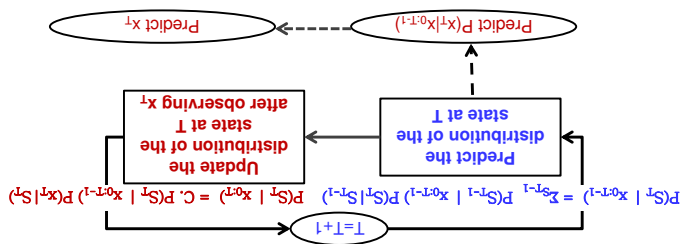
Estimating the state

- At t , you have some beliefs for the states
- You predict the beliefs for the state at $t+1$
- And update these after observing O_{t+1}



Prediction with HMMs

- The probability distribution for the observations at the next time is a mixture:
 - $P(x_{T+1} | x_{0:T-1}) = \sum_{s_T} P(x_{T+1} | s_T) P(s_T | x_{0:T-1})$
- The actual observation can be predicted from $P(x_{T+1} | x_{0:T-1})$



Predicting the next observation

- f) and/or g() may not be nice linear functions
- Conventional Kalman update rules are no longer valid
- z and/or γ may not be Gaussian
- Gaussian based update rules no longer valid

$$s'_t = f(s_{t-1}, \mathcal{E}'_t)$$

$$o'_t = g(s'_t, \gamma'_t)$$

Problems

The Linear Gaussian model (KF)

■ Iterative prediction and update

$$P_0(s) = \text{Gaussian}(s; \bar{s}, R)$$

$$P(s'_t | s_{t-1}) = \text{Gaussian}(s'_t; \mu_\varepsilon + A_t s_{t-1}, \Theta_\varepsilon)$$

$$P(o'_t | s'_t) = \text{Gaussian}(o'_t; B_t s'_t, \Theta_\gamma)$$

$$P(s'_t | o_{0:t}) = \text{Gaussian}(s'_t; \bar{s}_t, R'_t)$$

$$P(s'_t | o_{0:t-1}) = \text{Gaussian}(s'_t; \bar{s}_t, R'_t)$$

$$K_t = R'_t B_t^T (B_t R'_t B_t^T + \Theta_\gamma)^{-1}$$

$$\bar{s}_t = \bar{s}'_t + K_t (o_t - B_t \bar{s}'_t)$$

$$R'_t = (I - K_t B_t) R'_t$$

Linear Gaussian Model

All distributions remain Gaussian

$P(s) = \text{a priori}$
 $P(s_t | s_0) = \text{Transition prob.}$
 $P(o_t | s_t) = \text{State output prob.}$

$$s'_t = A_t s_{t-1} + \mathcal{E}'_t$$

$$o'_t = B_t s'_t + \gamma'_t$$

- Estimation requires knowledge of $P(o|s)$
- Difficult to estimate for nonlinear $g()$
- Even if it can be estimated, may not be tractable with update loop
- Estimation also requires knowledge of $P(s'_t | s_{t-1})$
- Difficult for nonlinear $f()$
- May not be amenable to closed form integration

$$s'_t = f(s_{t-1}, \mathcal{E}'_t)$$

$$o'_t = g(s'_t, \gamma'_t)$$

$$P(s'_t | o_{0:t-1}) = \int_{-\infty}^{\infty} P(s_{t-1} | o_{0:t-1}) P(s'_t | s_{t-1}) P(\mathcal{E}'_t) ds_{t-1}$$

$$P(s'_t | o_{0:t}) = C P(s'_t | o_{0:t-1}) P(o'_t | s'_t)$$

The problem with non-linear functions

The Kalman filter

■ Prediction

$$\bar{s}_t = A_{t-1} \bar{s}_{t-1} + \mu_\varepsilon$$

$$R_t = \Theta_\varepsilon + A_{t-1} R_{t-1} A_{t-1}^T$$

■ Update

$$K_t = R_t B_t^T (B_t R_t B_t^T + \Theta_\gamma)^{-1}$$

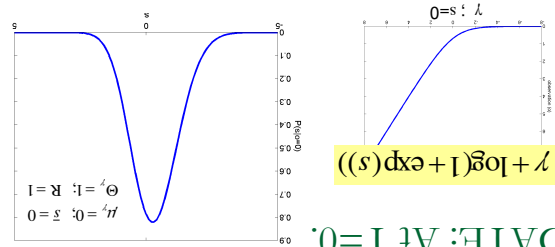
$$\hat{s}_t = \bar{s}_t + K_t (o_t - B_t \bar{s}_t)$$

$$\hat{R}_t = (I - K_t B_t) R_t$$

■ = Not Gaussian

$$P(s_0 | o_0) = C \exp \left(-0.5(s_0 - \sigma)^T R^{-1}(s_0 - \sigma) \right)$$

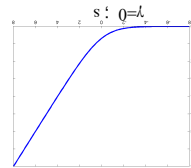
$$P(s_0 | o_0) = CGaussian(o; \mu, \Sigma) = CGaussian(o; \mu, \Sigma) \cdot \Theta^T(s_0)$$



UPDATE: At T=0.

$$P(o | s) = Gaussian(o; \mu, \Sigma) = \log(1 + \exp(s))$$

$$P(\gamma) = Gaussian(\gamma; \mu, \Sigma)$$



■ P(o|s) = ?

$$o = \gamma + \log(1 + \exp(s))$$

Example: a simple nonlinearity

■ Even if a closed form exists initially, it will typically become intractable very quickly

$$P(o_i | s_i) = \sum_{\gamma: g(s_i, \gamma) = o_i} \frac{P(\gamma)}{P(\gamma)}$$

■ The PDF may not have a closed form

$$o_i = g(s_i, \gamma_i)$$

The problem with nonlinearity

■ = intractable

$$P(s_1 | o_0) = \int_{-\infty}^{\infty} C \exp \left(-0.5(s_1 - \mu)^T R^{-1}(s_1 - \mu) - \log(1 + \exp(s_0)) - o_0 \Theta^T(s_1 - s_0) \right) ds_0$$

■ Prediction $P(s_1 | o_0) = \int_{-\infty}^{\infty} P(s_0 | o_0) P(s_1 | s_0) ds_0$

$$P(s_1 | s_{1-1}) = Gaussian(s_1; s_{1-1}, \Theta^T)$$

■ Trivial, linear state transition equation

$$s_1^T = s_{1-1}^T + \mathcal{E}$$

$$P(\mathcal{E}) = Gaussian(\mathcal{E}; 0, \Theta^T)$$

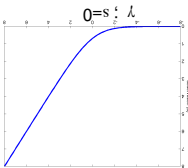
Prediction for T = 1

$$P(s_0 | o_0) = CGaussian(o; \mu, \Sigma) = \log(1 + \exp(s_0))$$

■ Update $P(s_0 | o_0) = CP(o_0 | s_0) P(s_0)$

$$P(s_0) = P_0(s) = Gaussian(s; \bar{s}, R)$$

■ Assume initial probability P(s) is Gaussian



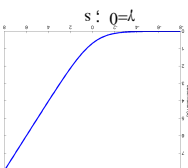
$$o = \gamma + \log(1 + \exp(s))$$

Example: At T=0.

■ Assume γ is Gaussian

$$P(\gamma) = Gaussian(\gamma; \mu, \Sigma)$$

■ P(o|s) = ?



$$o = \gamma + \log(1 + \exp(s))$$

Example: a simple nonlinearity

Update at T=1 and later

- Update at T=1

$$P(s_t | o_{t-1}) = CP(s_t | o_{t-1})P(o_t | s_t)$$

- Intractable

- Prediction for T=2

$$P(s_t | o_{t-1}) = \int_{-\infty}^{\infty} P(s_t | o_{t-1})P(s_t | s_{t-1})P(o_t | s_t) ds_{t-1}$$

- Intractable

- Similar problems arise for the state prediction equation
- $P(s_t | s_{t-1})$ may not have a closed form
- Even if it does, it may become intractable within the prediction and update equations
- Particularly the prediction equation, which includes an integration operation

$$s_t = f(s_{t-1}, \mathcal{E}_t)$$

The State prediction Equation

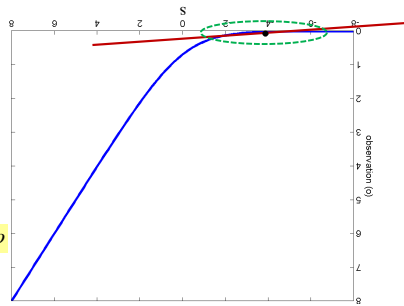
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Simplifying the problem: Linearize

- The tangent at any point is a good local approximation if the function is sufficiently smooth



$$o = \gamma + \log(1 + \exp(s))$$

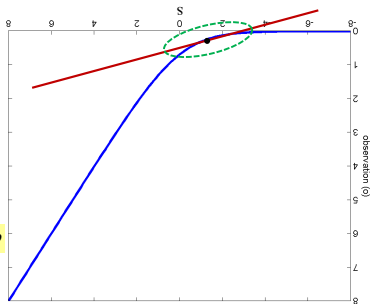
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Simplifying the problem: Linearize

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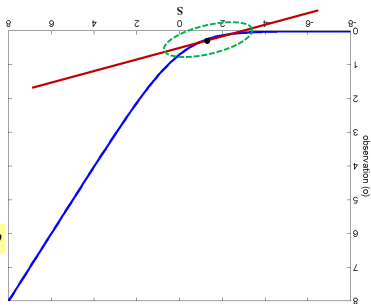
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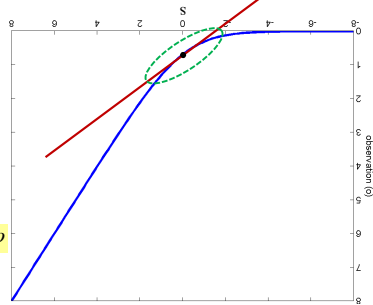
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Simplifying the problem: Linearize

- The tangent at any point is a good local approximation if the function is sufficiently smooth



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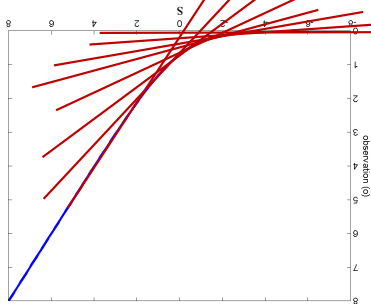
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Simplifying the problem: Linearize

- The tangent at any point is a good local approximation if the function is sufficiently smooth



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Simplifying the problem: Linearize

$$s_t = f(s_{t-1}) + \varepsilon \leftarrow s_t \approx \varepsilon + f(\delta_{t-1}) + J_f(\delta_{t-1})(s_{t-1} - \delta_{t-1})$$

- Linearize around the mean of the updated distribution of s at $t-1$
- Which should be Gaussian

$$P(s_{t-1} | o_{0:t-1}) = \text{Gaussian}(s_{t-1}; \delta_{t-1}, R_{t-1})$$

- Solution: Linearize
- Again, direct use of $f()$ can be disastrous

$$s_t = f(s_{t-1}) + \varepsilon \quad P(\varepsilon) = \text{Gaussian}(\varepsilon; 0, \Theta_\varepsilon)$$

Prediction?

$$P(o | s) = \text{Gaussian}(o; g(s) + J_g(s)(s - \bar{s}), \Theta_\gamma)$$

$$P(\gamma) = \text{Gaussian}(\gamma; 0, \Theta_\gamma)$$

- Observation PDF is Gaussian

$$o \approx \gamma + g(s) \leftarrow o \approx \gamma + g(\bar{s}) + J_g(\bar{s})(s - \bar{s})$$

$$P(s) = \text{Gaussian}(s, R)$$

Linearizing the observation function

$$o \approx \gamma + g(s) \leftarrow o \approx \gamma + g(\bar{s}) + J_g(\bar{s})(s - \bar{s})$$

$$P(s) = \text{Gaussian}(s, R)$$

Linearizing the observation function

- Simple first-order Taylor series expansion
 - $J()$ is the Jacobian matrix
 - Simply a determinant for scalar state
- Expansion around a *prior* (or predicted) mean of the state

$$P(s_t | o_{0:t-1}) = \int_{-\infty}^{\infty} P(s_{t-1} | o_{0:t-1}) P(s_t | s_{t-1}) P(s_{t-1}) ds_{t-1}$$

- The predicted state probability is:

$$P(s_t | s_{t-1}) = \text{Gaussian}(s_t; f(\delta_{t-1}) + J_f(\delta_{t-1})(s_{t-1} - \delta_{t-1}), \Theta_\varepsilon)$$

- The state transition probability is now:

$$P(s_{t-1} | o_{0:t-1}) = \text{Gaussian}(s_{t-1}; \delta_{t-1}, R_{t-1}) \quad P(\varepsilon) = \text{Gaussian}(\varepsilon; 0, \Theta_\varepsilon)$$

$$s_t = f(s_{t-1}) + \varepsilon \leftarrow s_t \approx \varepsilon + f(\delta_{t-1}) + J_f(\delta_{t-1})(s_{t-1} - \delta_{t-1})$$

Prediction

- Gaussian!!
- Note: This is actually only an approximation

$$P(s | o) = \text{Gaussian}(s; \bar{s} + R J_g(\bar{s})^T (J_g(\bar{s}) R J_g(\bar{s})^T + \Theta_\gamma)^{-1} (o - g(\bar{s})) - R J_g(\bar{s})^T (J_g(\bar{s}) R J_g(\bar{s})^T + \Theta_\gamma)^{-1} J_g(\bar{s}) R)$$

$$P(s | o) = \text{Gaussian}(o; g(\bar{s}) + J_g(\bar{s})(s - \bar{s}), \Theta_\gamma) \text{Gaussian}(s; \bar{s}, R)$$

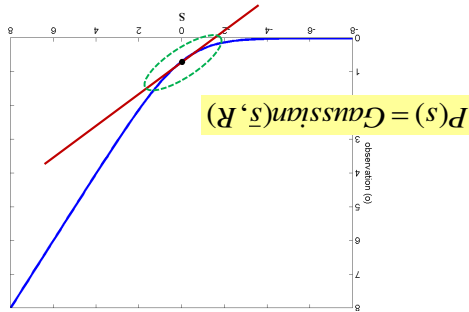
$$P(s) = \text{Gaussian}(s; \bar{s}, R) \quad P(s | o) = CP(o | s)P(s)$$

$$P(o | s) = \text{Gaussian}(o; g(\bar{s}) + J_g(\bar{s})(s - \bar{s}), \Theta_\gamma)$$

$$o \approx \gamma + g(\bar{s}) + J_g(\bar{s})(s - \bar{s})$$

UPDATE.

- $P(s)$ is small approximation error is large
- Most of the probability mass of s is in low-error regions



Most probability is in the low-error region

The Extended Kalman filter

- Prediction
 - $\underline{s}_t = f(\underline{s}_{t-1})$
 - $R_t = \Theta_\varepsilon + A_t R_{t-1} A_t^T$
 - $K_t = R_t B_t^T (B_t R_t B_t^T + \Theta_\gamma)^{-1}$
 - $\hat{s}'_t = \underline{s}_t + K_t (o_t - g(\underline{s}_t))$
 - $\hat{R}'_t = (I - K_t B_t^T) R_t$
- Update

Linearized Prediction and Update

- Prediction for time t
 - $P(s_t | o_{1:t}) = \text{Gaussian}(s_t; \underline{s}_t, R_t)$
 - $R_t = J_f(\underline{s}_{t-1}) R_{t-1} J_f^T(\underline{s}_{t-1}) + \Theta_\varepsilon$
- Update at time t
 - $\hat{s}'_t = \underline{s}_t + R_t J_g^T(\underline{s}_t) (J_g(\underline{s}_t) R_t J_g^T(\underline{s}_t) + \Theta_\gamma)^{-1} J_g(\underline{s}_t) R_t$
 - $\hat{R}'_t = (I - R_t J_g^T(\underline{s}_t) (J_g(\underline{s}_t) R_t J_g^T(\underline{s}_t) + \Theta_\gamma)^{-1} J_g(\underline{s}_t) R_t) R_t$
- Update at time t

Prediction

$P(s_{t-1} | o_{1:t-1}) = \text{Gaussian}(s_{t-1}; \underline{s}_{t-1}, R_{t-1})$

- The predicted state probability is:
 - $P(s_t | o_{1:t-1}) = \int_{-\infty}^{\infty} \text{Gaussian}(s_{t-1}; \underline{s}_{t-1}, R_{t-1}) \text{Gaussian}(s_t; f(\underline{s}_{t-1}) + J_f(\underline{s}_{t-1})(s_{t-1} - \underline{s}_{t-1}), \Theta_\varepsilon) ds_{t-1}$
 - $P(s_t | o_{1:t-1}) = \int_{-\infty}^{\infty} P(s_{t-1} | o_{1:t-1}) P(s_t | s_{t-1}) ds_{t-1}$
 - $P(s_t | s_{t-1}) = \text{Gaussian}(s_t; f(\underline{s}_{t-1}) + J_f(\underline{s}_{t-1})(s_{t-1} - \underline{s}_{t-1}), \Theta_\varepsilon)$
- Gaussian!!
 - $P(s_t | o_{1:t-1}) = \text{Gaussian}(s_t; f(\underline{s}_{t-1}), J_f(\underline{s}_{t-1}) R_{t-1} J_f^T(\underline{s}_{t-1}) + \Theta_\varepsilon)$

□ This is actually only an approximation

The Kalman filter

- Prediction
 - $\underline{s}_t = A_t \underline{s}_{t-1} + \mu_\varepsilon$
 - $R_t = \Theta_\varepsilon + A_t R_{t-1} A_t^T$
- Update
 - $K_t = R_t B_t^T (B_t R_t B_t^T + \Theta_\gamma)^{-1}$
 - $\hat{s}'_t = \underline{s}_t + K_t (o_t - B_t \underline{s}_t)$
 - $\hat{R}'_t = (I - K_t B_t^T) R_t$

Linearized Prediction and Update

- Prediction for time t
 - $P(s_t | o_{1:t}) = \text{Gaussian}(s_t; \underline{s}_t, R_t)$
 - $R_t = A_t R_{t-1} A_t^T + \Theta_\varepsilon$
- Update at time t
 - $\hat{s}'_t = \underline{s}_t + R_t B_t^T (B_t R_t B_t^T + \Theta_\gamma)^{-1} (o_t - g(\underline{s}_t))$
 - $\hat{R}'_t = (I - R_t B_t^T (B_t R_t B_t^T + \Theta_\gamma)^{-1} B_t) R_t$
- Update at time t

The linearized prediction/update

- Given: two non-linear functions for state update and observation generation
 - $o_t = g(s_t) + \gamma$
 - $s_t = f(s_{t-1}) + \varepsilon$
- Note: the equations are *deterministic* non-linear functions of the state variable
 - They are *linear* functions of the noise!
 - Non-linear functions of stochastic noise are slightly more complicated to handle

- EKFs are probably the most commonly used algorithm for tracking and prediction
- Most systems are non-linear
- Specifically, the relationship between state and observation is usually nonlinear
- The approach can be extended to include non-linear functions of noise as well
- The term "Kalman filter" often simply refers to an *extended Kalman filter* in most contexts.
- But..

A different problem: Non-Gaussian PDFs

$$o_t = g(s_t) + \gamma$$

$$s_t = f(s_{t-1}) + \varepsilon$$

- We have assumed so far that:
 - $P_0(s)$ is Gaussian or can be approximated as Gaussian
 - $P(\varepsilon)$ is Gaussian
 - $P(\gamma)$ is Gaussian

This has a happy consequence: All distributions remain Gaussian

- But when any of these are not Gaussian, the results are not so happy

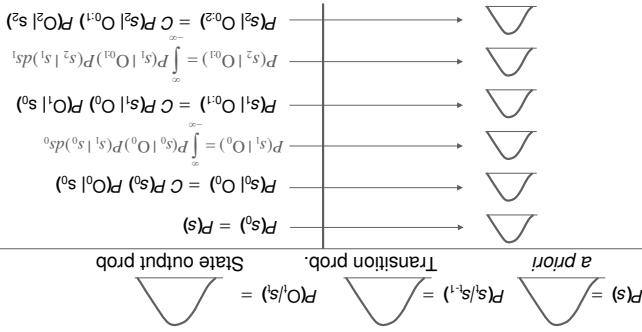
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A different problem: Non-Gaussian PDFs

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$$s_t = f(s_{t-1}) + \varepsilon$$

All distributions remain Gaussian



Linear Gaussian Model

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- Most systems are non-linear
- Specifically, the relationship between state and observation is usually nonlinear
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- But..

EKFs

EKFs have limitations



- If the non-linearity changes too quickly with s , the linear approximation is invalid
- Unstable
- The estimate is often biased
- The true function lies entirely on one side of the approximation
- Various extensions have been proposed:
 - Unscented Kalman filters (UKF)
 - Invariant extended Kalman filters (IEKF)

$$P(o_t | s_t) = P(\gamma = o_t - Bs_t) = \sum_{i=0}^t w_i \text{Gaussian}(o; \mu_i + Bs_t, \Theta_i)$$

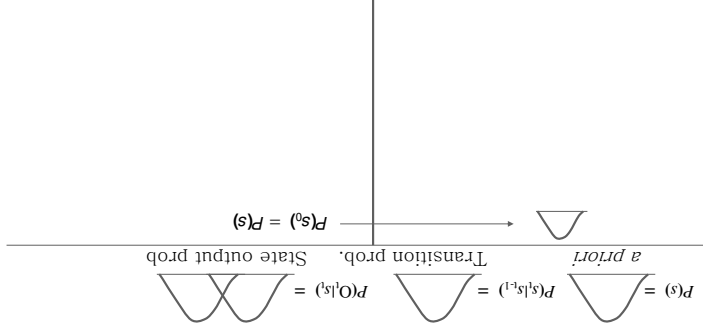
- $P(\gamma)$ is a mixture of only two Gaussians
- o is a linear function of s
- Non-linear functions would be linearized anyway
- $P(o|s)$ is also a Gaussian mixture!

A simple case

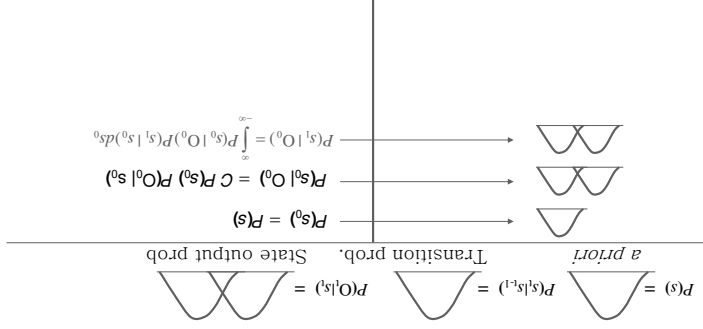
$$P(\gamma) = \sum_{i=0}^t w_i \text{Gaussian}(\gamma; \mu_i, \Theta_i)$$

$$o_t = Bs_t + \gamma$$

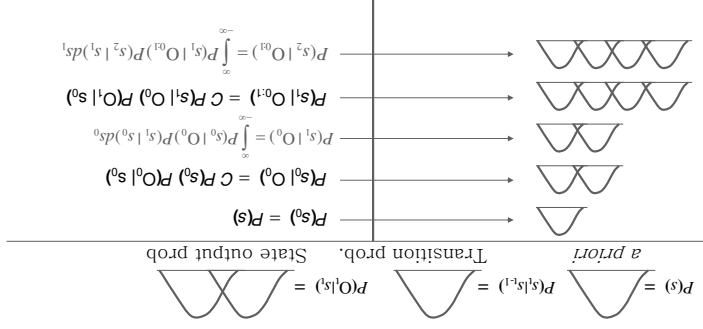
When distributions are not Gaussian



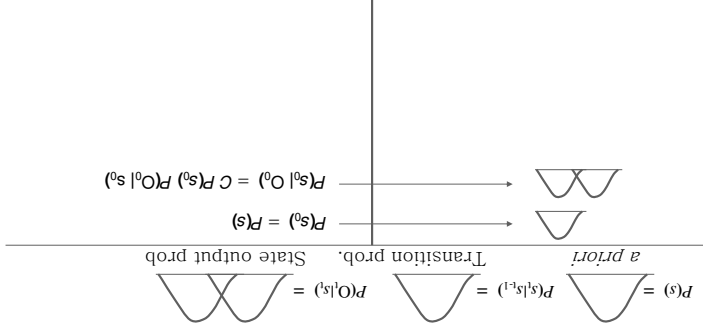
When distributions are not Gaussian



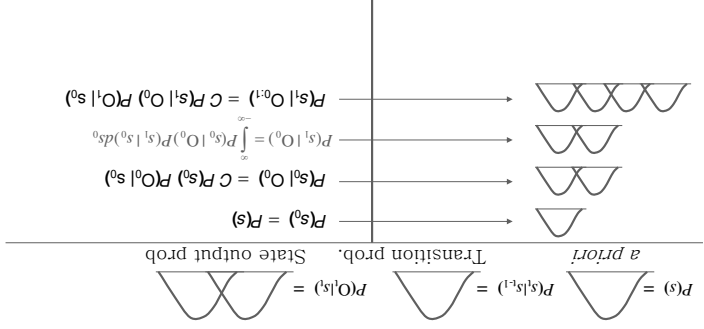
When distributions are not Gaussian



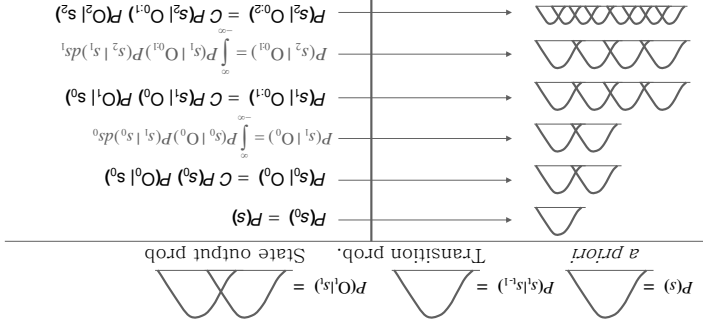
When distributions are not Gaussian



When distributions are not Gaussian



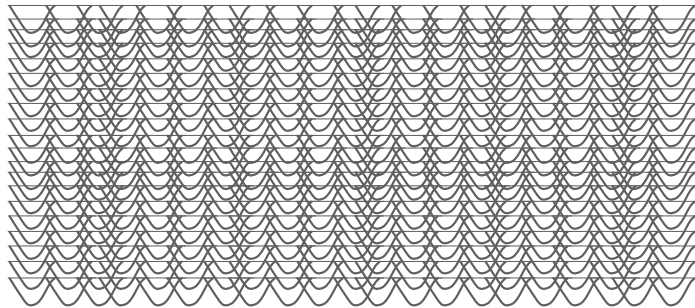
When distributions are not Gaussian



When $P(O_{i+1})$ has more than one Gaussian, after only a few time steps...

When distributions are not Gaussian

$$P(s'_i | O^{(i)}) =$$



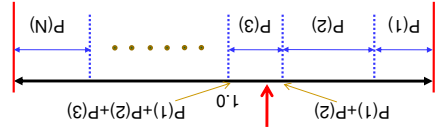
We have too many Gaussians for comfort..

Related Topic: How to sample from a Distribution?

- "Sampling from a Distribution $P(x; \Gamma)$ with parameters Γ "
- Generate random numbers such that
- The distribution of a large number of generated numbers is $P(x; \Gamma)$
- The parameters of the distribution are Γ
- Many algorithms to generate RVs from a variety of distributions
- Generation from a uniform distribution is well studied
- Uniform RVs used to sample from multinomial distributions
- Other distributions: Most commonly, transform a uniform RV to the desired distribution

Sampling from a multinomial

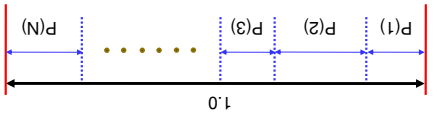
- Given a multinomial over N symbols, with probability of i th symbol = $P(i)$
- Randomly generate symbols from this distribution
- Can be done by sampling from a uniform distribution



Sampling a multinomial

- Segment a range $(0, 1)$ according to the probabilities $P(i)$
- The $P(i)$ terms will sum to 1.0
- Randomly generate a number from a uniform distribution
- Matlab: "rand"
- Generates a number between 0 and 1 with uniform probability
- If the number falls in the i th segment, select the i th symbol

Sampling a multinomial



- Segment a range $(0, 1)$ according to the probabilities $P(i)$
- The $P(i)$ terms will sum to 1.0

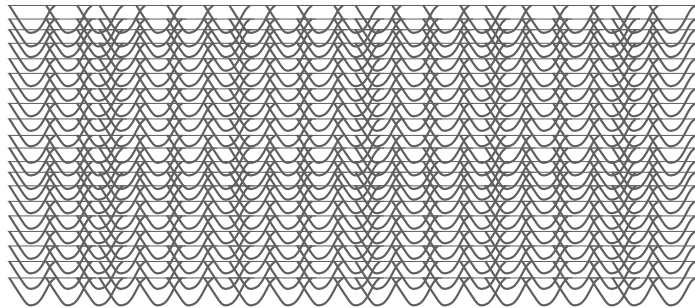
Related Topic: Sampling from a Gaussian

- Many algorithms
- Simplest: add many samples from a uniform RV
- The sum of 12 uniform RVs (uniform in $(0, 1)$) is approximately Gaussian with mean 6 and variance 1
- For scalar Gaussian, mean μ , std dev σ :

$$x = \sum_{i=1}^{12} r_i - 6$$
- Matlab : $x = \mu + \text{randn} * \sigma$
- "randn" draws from a Gaussian of mean=0, variance=1

When distributions are not Gaussian

$$P(s'_i | O^{(i)}) =$$

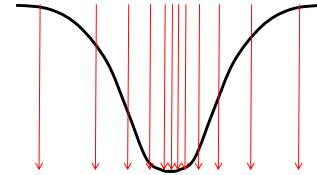


We have too many Gaussians for comfort..

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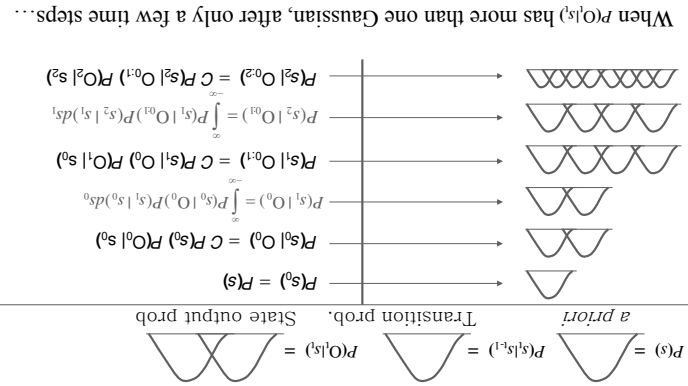
- A large-enough collection of randomly-drawn samples from a distribution will approximately quantize the space of the random variable into equi-probable regions
- We have more random samples from high-probability regions and fewer samples from low-probability regions



Discrete approximation to a distribution

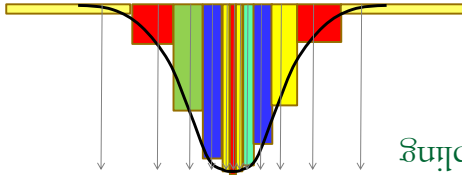
- Multivariate (d-dimensional) Gaussian with mean μ and covariance Θ
 - Compute eigen value matrix Λ and eigenvector matrix E for Θ
 - $\Theta = E \Lambda E^T$
 - Generate d -mean unit-variance numbers x_1, \dots, x_d
 - Arrange them in a vector: $X = [x_1 \dots x_d]^T$
 - Multiply X by the square root of Λ and E , add μ
- $$Y = \mu + E \text{sqrt}(\Lambda) X$$

Related Topic: Sampling from a Gaussian



When $P(O_i | s_i)$ has more than one Gaussian, after only a few time steps...

- A PDF can be approximated as a uniform probability distribution over randomly drawn samples
 - Since each sample represents approximately the same probability mass ($1/M$ if there are M samples)
- $$P(x) \approx \frac{1}{M} \sum_{i=1}^M \delta(x - x_i)$$



Discrete approximation: Random Sampling

- Select a Gaussian by sampling the multinomial distribution of weights: $j \sim \text{multinomial}(w_1, w_2, \dots)$
- Sample from the selected Gaussian

$$\sum_i w_i \text{Gaussian}(X; \mu_i, \Theta_i)$$

Sampling from a Gaussian Mixture

- The complexity of the distribution increases exponentially with time
- This is a consequence of having a *continuous* state space
- Only Gaussian PDFs propagate without increase of complexity
- Discrete-state systems do not have this problem
- The number of states in an HMM stays fixed
- However, discrete state spaces are too coarse
- Solution: Combine the two concepts
- Discretize the state space dynamically

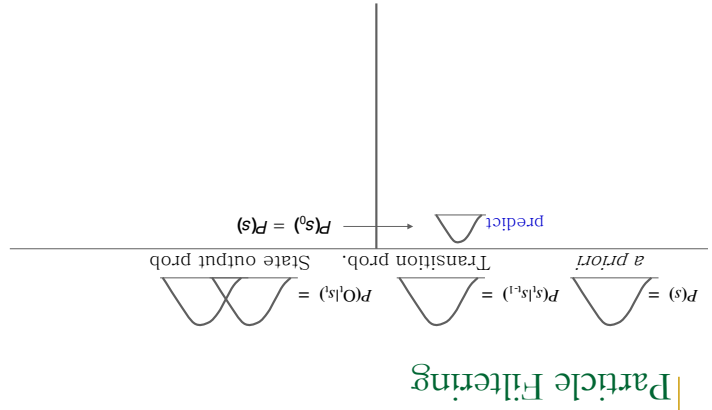
The problem of the exploding distribution

Note: Properties of a discrete distribution

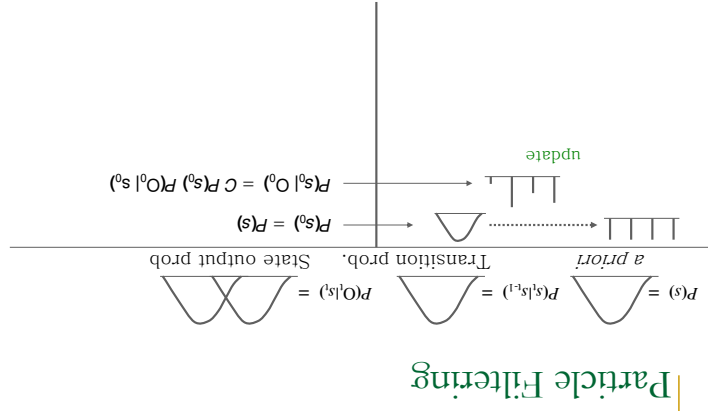
- $$P(x) \approx \frac{1}{M} \sum_{l=0}^{M-1} \delta(x - x_l)$$
- The product of a discrete distribution with another distribution is simply a weighted discrete probability
- $$\int_{-\infty}^{\infty} P(x) P(y|x) dx = \sum_{l=0}^{M-1} w_l P(y|x_l)$$
- The integral of the product is a mixture distribution

Discretizing the state space

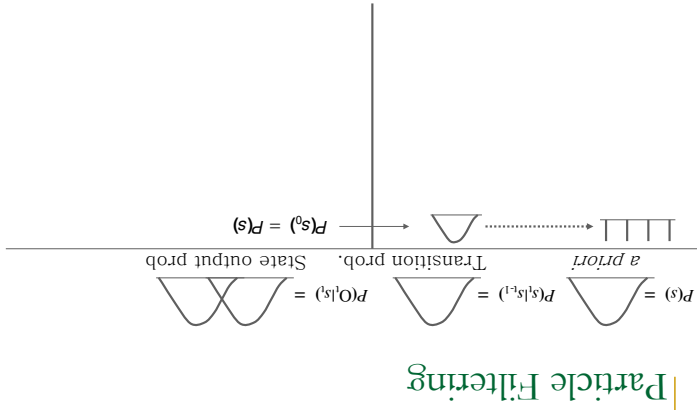
- At each time, discretize the predicted state space
- $$P(s_t | o_{1:t}) \approx \frac{1}{M} \sum_{l=0}^{M-1} \delta(s_t - s_l)$$
- s_l are randomly drawn samples from $P(s_t | o_{1:t})$
 - Propagate the discretized distribution



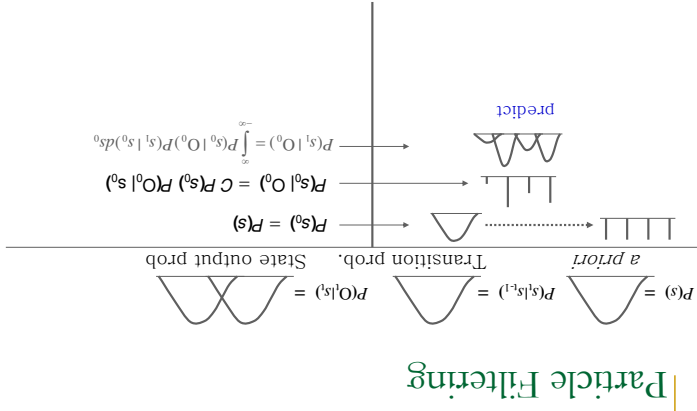
Assuming that we only generate **FOUR** samples from the predicted distributions



Assuming that we only generate **FOUR** samples from the predicted distributions

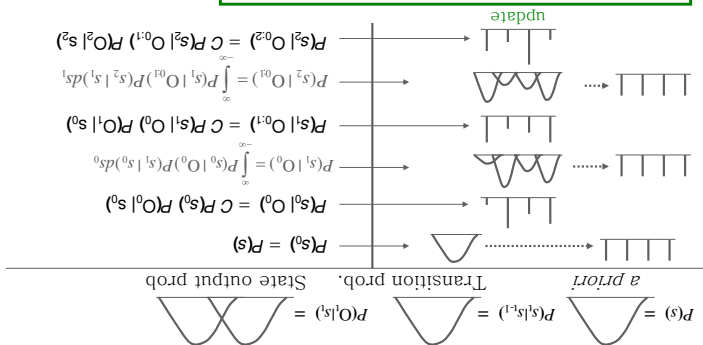


Assuming that we only generate **FOUR** samples from the predicted distributions



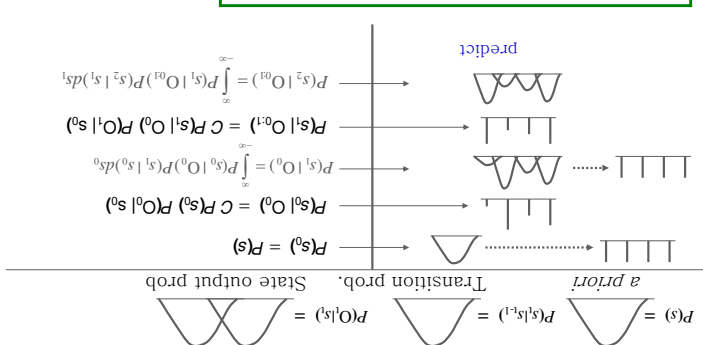
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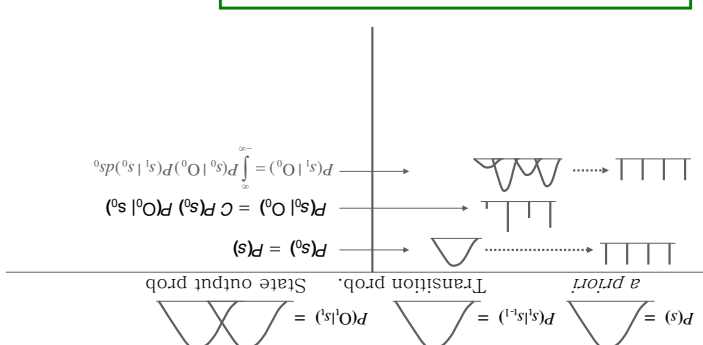
Particle Filtering

Assuming that we only generate **FOUR** samples from the predicted distributions



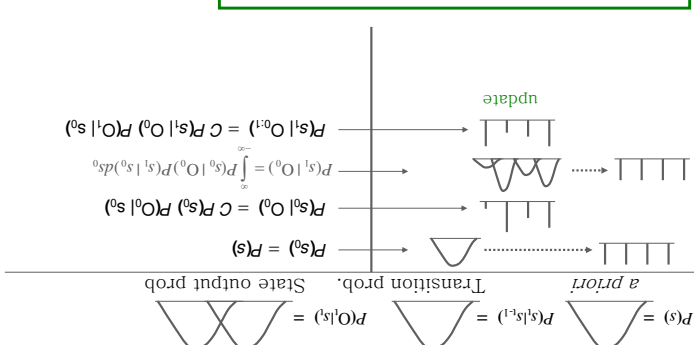
Particle Filtering

Assuming that we only generate **FOUR** samples from the predicted distributions



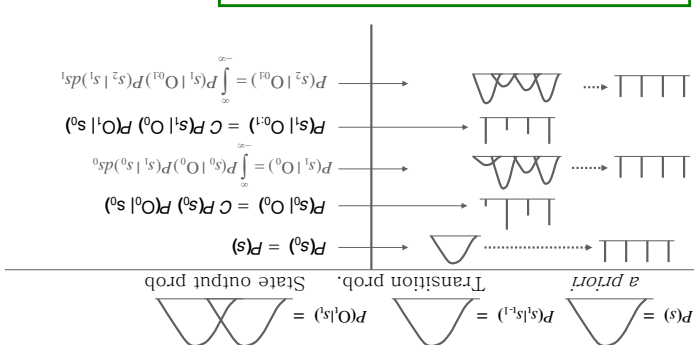
Particle Filtering

Assuming that we only generate **FOUR** samples from the predicted distributions



Particle Filtering

Assuming that we only generate **FOUR** samples from the predicted distributions



Particle Filtering

- Discretize state space at the prediction step
 - By sampling the continuous predicted distribution
 - If appropriately sampled, all generated samples may be considered to be equally probable
 - Sampling results in a **discrete uniform** distribution
- Update step updates the distribution of the quantized state space
 - Results in a **discrete non-uniform** distribution
- Predicted state distribution for the next time instant will again be continuous
 - Must be **discretized** again by sampling
- At any step, the current state distribution will not have more components than the number of samples generated at the previous sampling step
 - The complexity of distributions remains constant

Particle Filtering

Estimating a state

- The algorithm gives us a discrete updated distribution over states:

$$P(s_t | o_{0:t}) = C \sum_{s_{t-1}}^{M-1} P_t^\gamma(o_t - g(s_{t-1})) \delta(s_t - s_{t-1}^\gamma)$$
- The actual state can be estimated as the mean of this distribution

$$\hat{s}_t = C \sum_{s_{t-1}}^{M-1} \hat{s}_t^\gamma P_t^\gamma(o_t - g(s_{t-1}^\gamma))$$
- Alternately, it can be the most likely sample

$$\hat{s}_t = s_{t-1}^\gamma : f = \arg \max_{s_{t-1}} P_t^\gamma(o_t - g(s_{t-1}^\gamma))$$

Particle Filtering

- Predict the state distribution at the next time

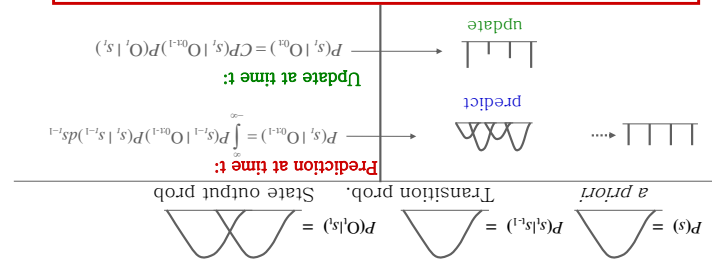
$$P(s_t | o_{0:t-1}) = C \sum_{s_{t-1}}^{M-1} P_t^\gamma(o_{t-1} - g(s_{t-1}^\gamma)) P^\varepsilon(s_t - f(s_{t-1}^\gamma))$$
- Sample the predicted state distribution

$$P(s_t | o_{0:t-1}) \approx \frac{1}{M} \sum_{s_{t-1}}^{M-1} \delta(s_t - s_{t-1}^\gamma) \text{ where } s_{t-1}^\gamma \rightarrow P(s_{t-1} | o_{0:t-1})$$

$$o_t = g(s_t) + \gamma \quad P^\gamma(\gamma)$$

$$s_t = f(s_{t-1}) + \varepsilon \quad P^\varepsilon(\varepsilon)$$

Number of mixture components in predicted distribution governed by number of samples in discrete distribution. By deriving a small (100-1000) number of samples at each time instant, all distributions are kept manageable.



Particle Filtering

Simulations with a Linear Model

- ε_t has a Gaussian distribution with 0 mean, known variance
- x_t has a mixture Gaussian distribution with known parameters
- Simulation:
 - Generate state sequence s_t from model
 - Generate sequence of X_t from model with one X_t term for every s_t term
 - Generate observation sequence O_t from s_t and X_t
 - Attempt to estimate s_t from O_t

$$o_t = s_{t-1} + \varepsilon_t \quad o_t = s_t + x_t$$

Particle Filtering

- Predict the state distribution at t

$$P(s_t | o_{0:t-1}) = C \sum_{s_{t-1}}^{M-1} P_t^\gamma(o_{t-1} - g(s_{t-1}^\gamma)) P^\varepsilon(s_t - f(s_{t-1}^\gamma))$$
- Sample the predicted state distribution at t

$$P(s_t | o_{0:t-1}) \approx \frac{1}{M} \sum_{s_{t-1}}^{M-1} \delta(s_t - s_{t-1}^\gamma) \text{ where } s_{t-1}^\gamma \rightarrow P(s_{t-1} | o_{0:t-1})$$
- Update the state distribution at t

$$P(s_t | o_{0:t}) = C \sum_{s_{t-1}}^{M-1} P_t^\gamma(o_t - g(s_{t-1}^\gamma)) \delta(s_t - s_{t-1}^\gamma)$$

$$o_t = g(s_t) + \gamma \quad P^\gamma(\gamma)$$

$$s_t = f(s_{t-1}) + \varepsilon \quad P^\varepsilon(\varepsilon)$$

Particle Filtering

- At $t = 0$, sample the initial state distribution

$$P(s_0 | o_{-1}) = P(s_0) \approx \frac{1}{M} \sum_{s_0}^{M-1} \delta(s_0 - s_0^\gamma) \text{ where } s_0^\gamma \rightarrow P_0(s)$$
- Update the state distribution with the observation

$$C = \frac{1}{\sum_{s_{t-1}}^{M-1} P_t^\gamma(o_t - g(s_{t-1}^\gamma))}$$

$$o_t = g(s_t) + \gamma \quad P^\gamma(\gamma)$$

$$s_t = f(s_{t-1}) + \varepsilon \quad P^\varepsilon(\varepsilon)$$

Simulation: Synthesizing data

Generate state sequence according to:

$$s_t = s_{t-1} + \varepsilon_t$$

Generate state sequence according to:

$$s_t = s_{t-1} + \varepsilon_t$$

ε_t is Gaussian with mean 0 and variance 10

Generate observation sequence from state sequence according to:

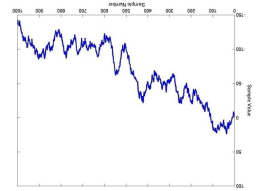
$$o_t = s_t + x_t$$

x_t is mixture Gaussian with parameters:

Means = [-4, 0, 4, 8, 12, 16, 18, 20]

Variances = [10, 10, 10, 10, 10, 10, 10, 10]

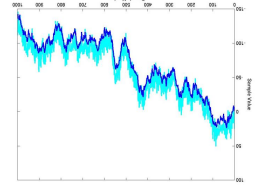
Mixture weights = [0.125, 0.125, 0.125, 0.125, 0.125, 0.125, 0.125, 0.125]



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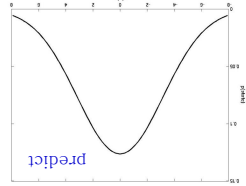
Simulation: Synthesizing data



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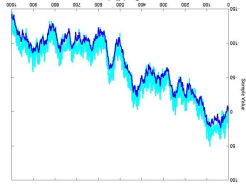
87

Combined figure for more compact representation



PREDICTED STATE DISTRIBUTION
AT TIME = 1

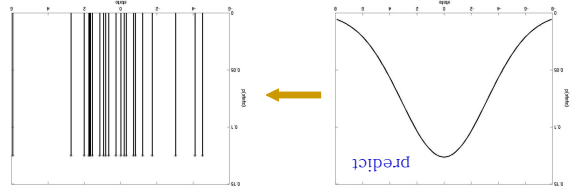
SIMULATION: TIME = 1



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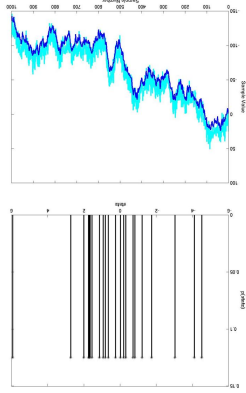
SIMULATION: TIME = 1



PREDICTED STATE DISTRIBUTION
AT TIME = 1

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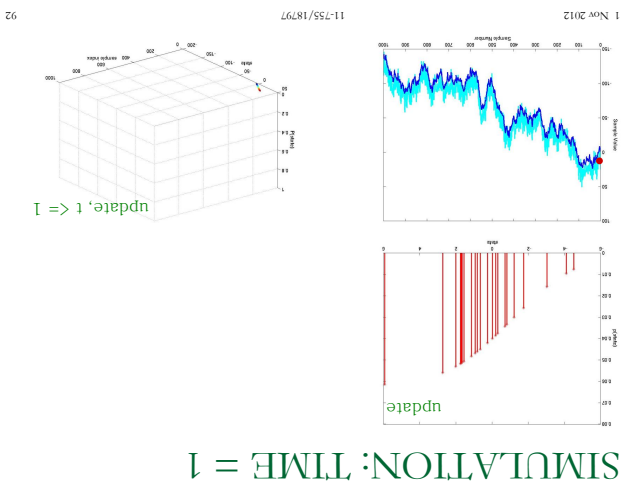
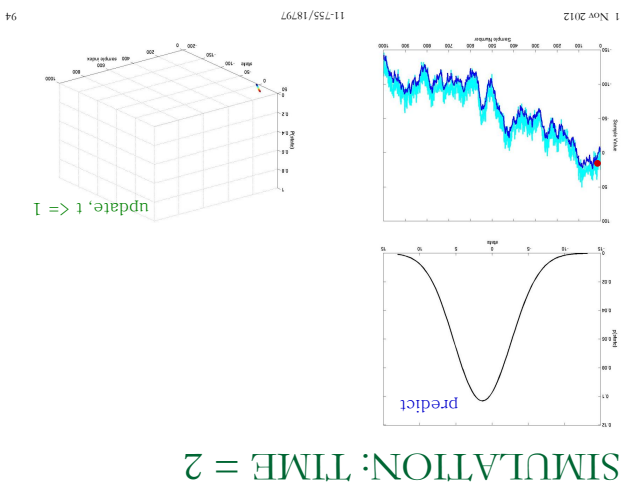
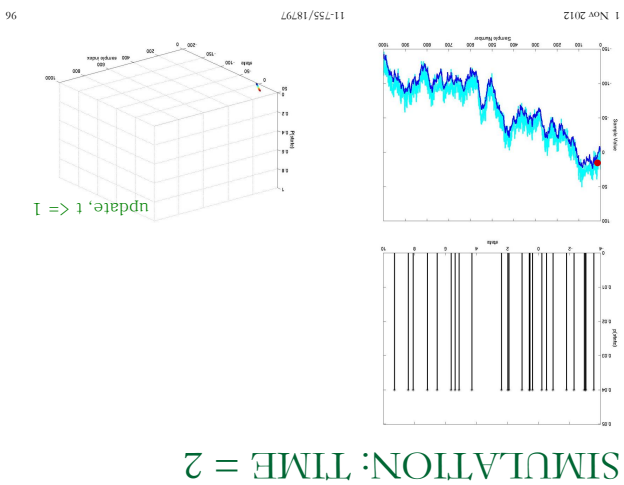
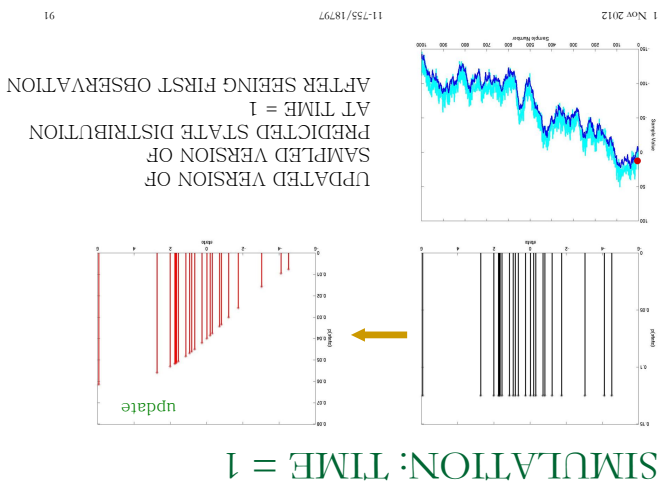
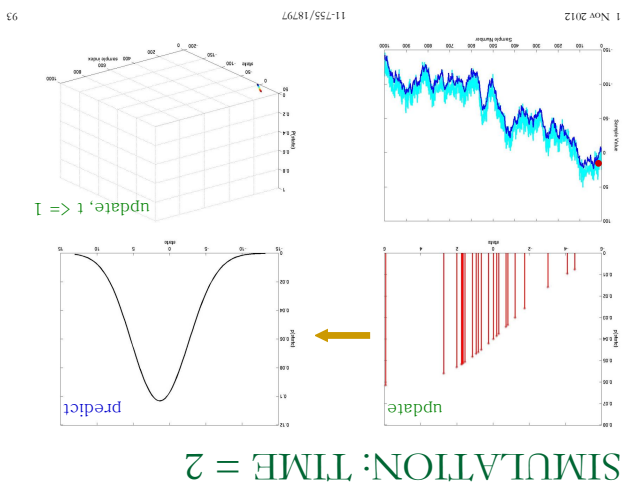
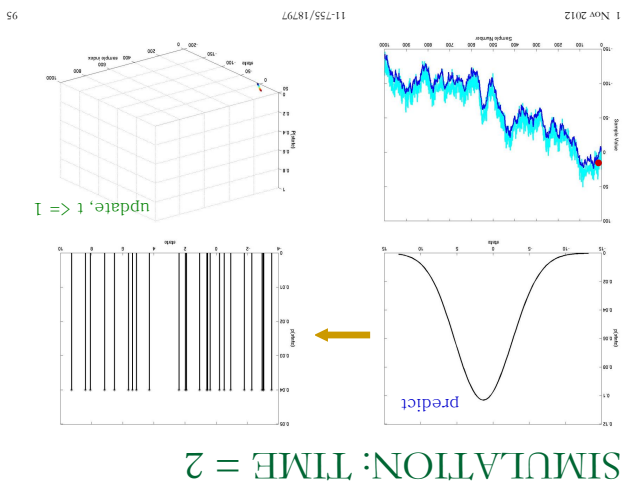


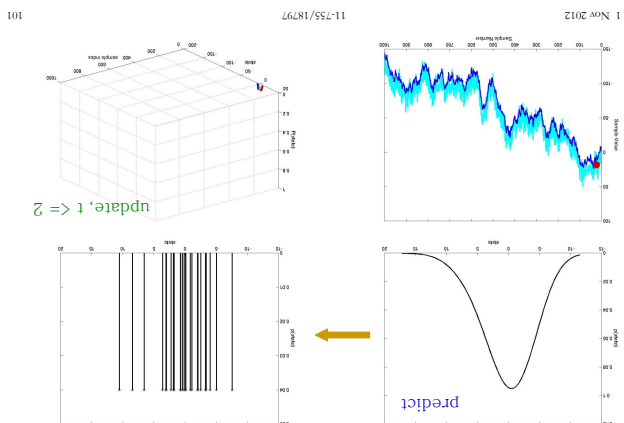
PREDICTED STATE DISTRIBUTION
AT TIME = 1

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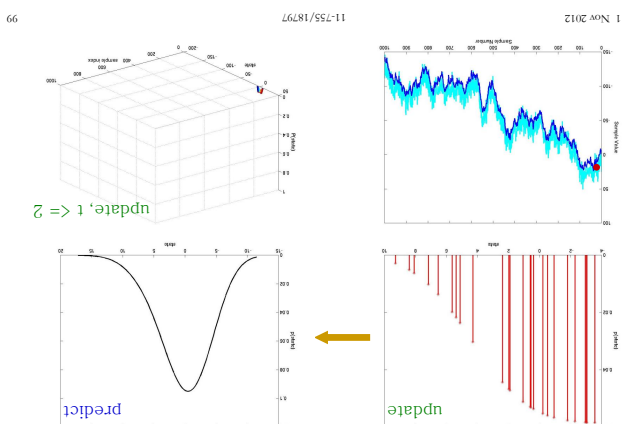
90

Simulation: Synthesizing data

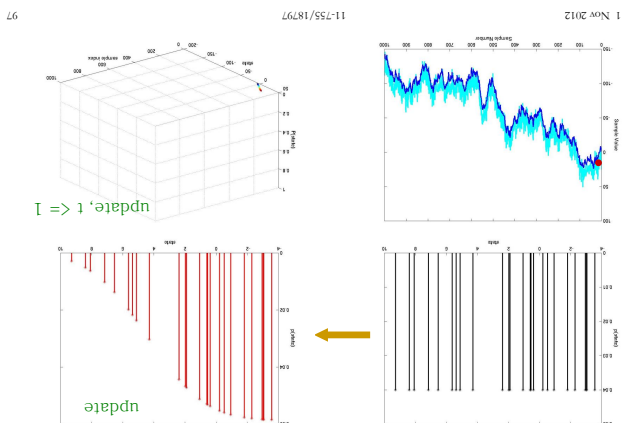




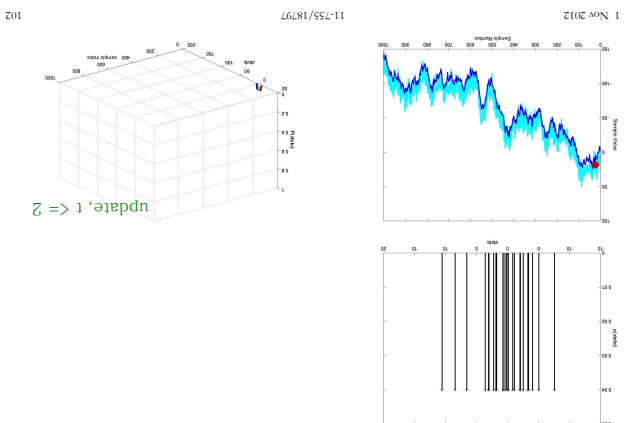
SIMULATION: TIME = 3



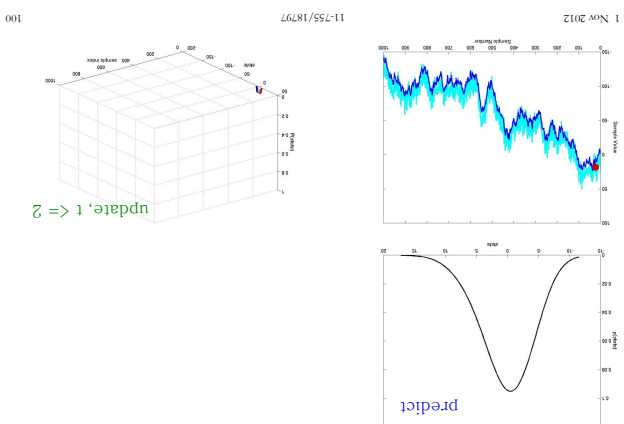
SIMULATION: TIME = 3



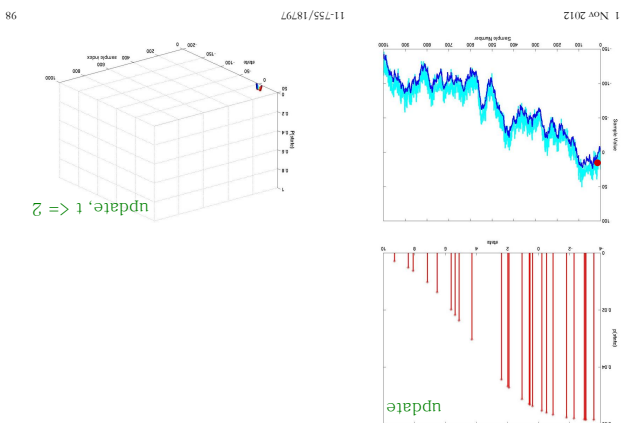
SIMULATION: TIME = 2



SIMULATION: TIME = 3



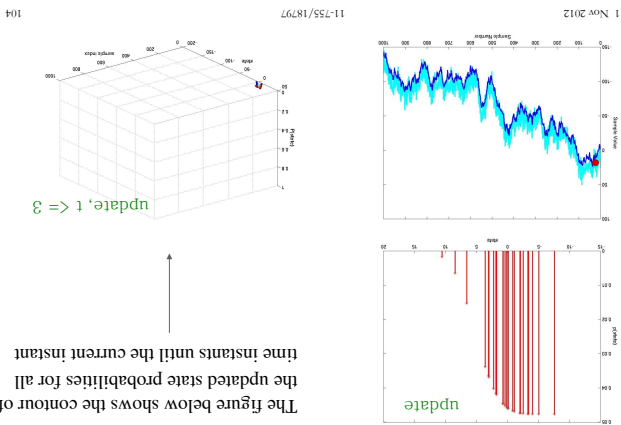
SIMULATION: TIME = 3



SIMULATION: TIME = 2

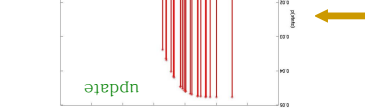
SIMULATION: TIME = 3

The figure below shows the contour of the updated state probabilities for all time instants until the current instant



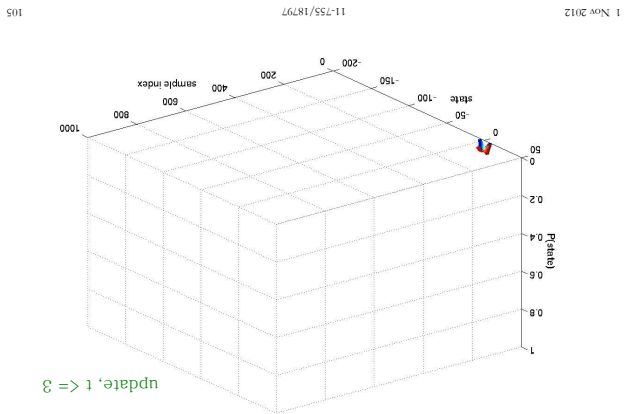
104

SIMULATION: TIME = 3



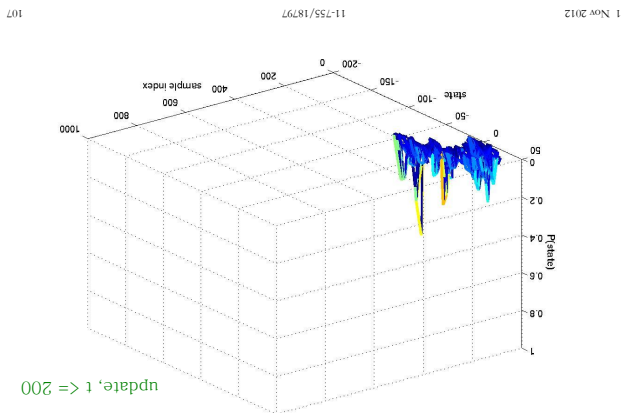
103

Simulation: Updated Probs Until T=3



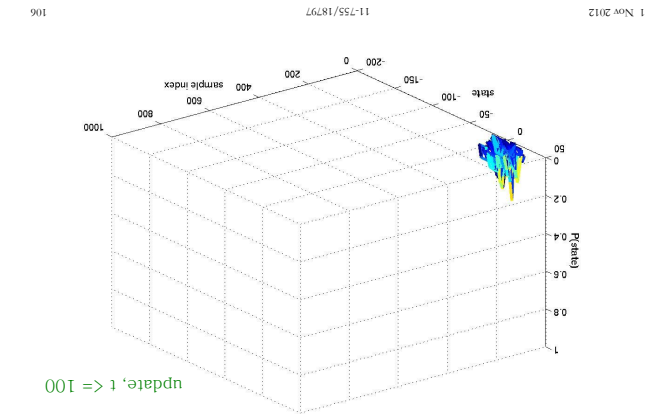
105

Simulation: Updated Probs Until T=200



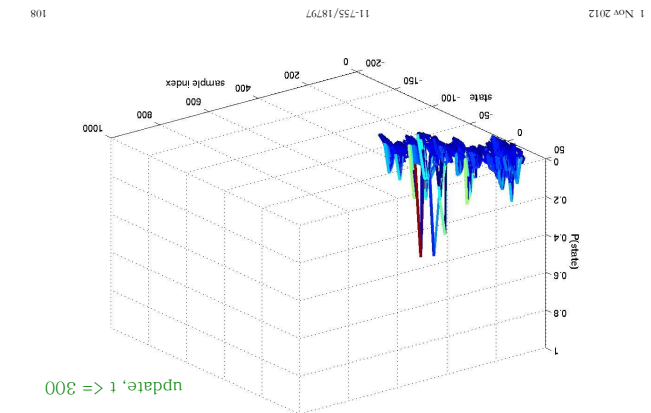
107

Simulation: Updated Probs Until T=100

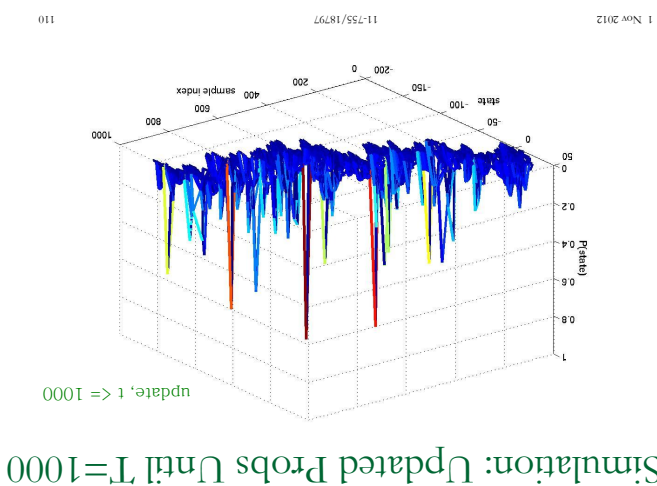
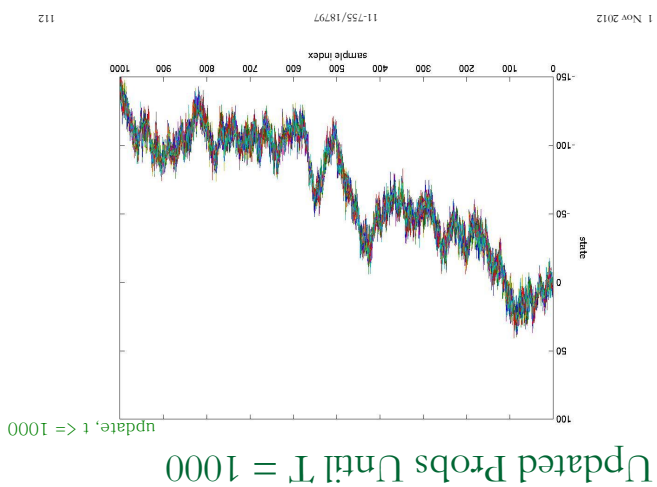
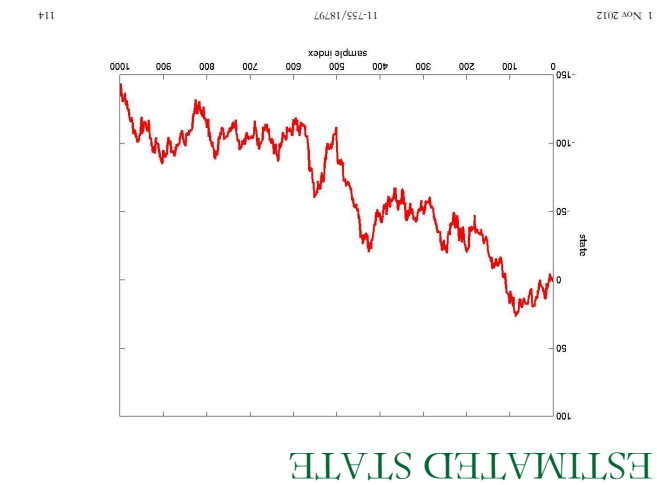
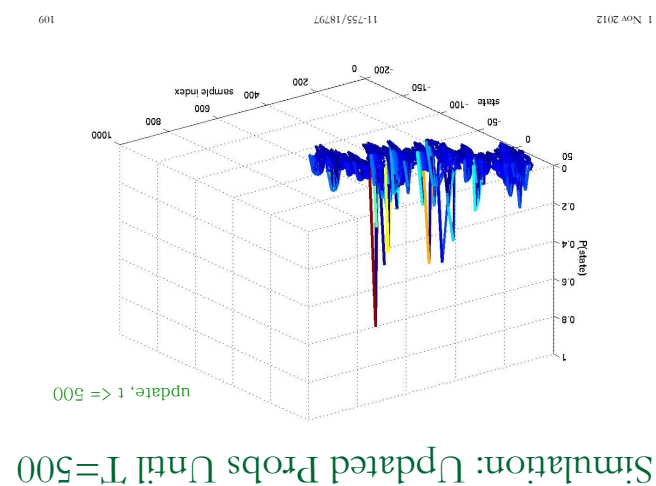
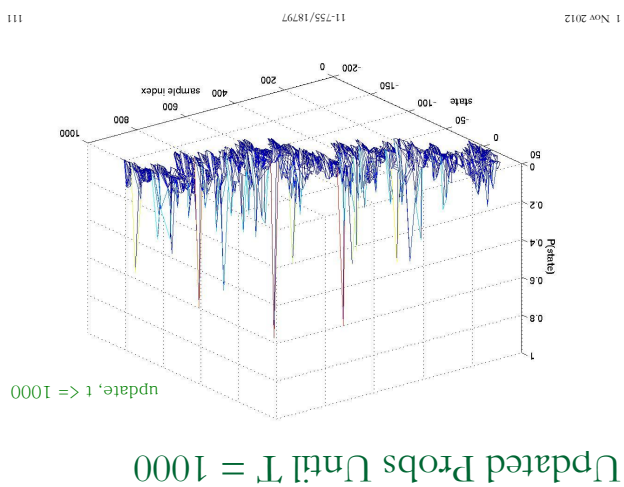
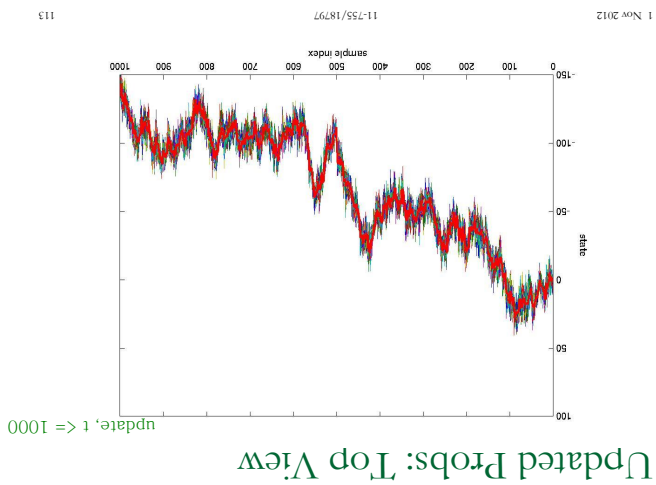


106

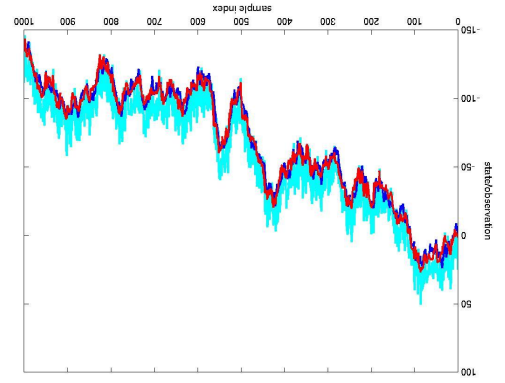
Simulation: Updated Probs Until T=300



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Observation, True States, Estimate



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Particle Filtering

- Generally quite effective in scenarios where EKF/UKF may not be applicable
 - Potential applications include tracking and edge detection in images!
 - Not very commonly used however
- Highly dependent on sampling
 - A large number of samples required for accurate representation
 - Samples may not represent mode of distribution
 - Some distributions are not amenable to sampling
 - Use importance sampling instead: Sample a Gaussian and assign non-uniform weights to samples

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