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# Prediction and Estimation, Part II

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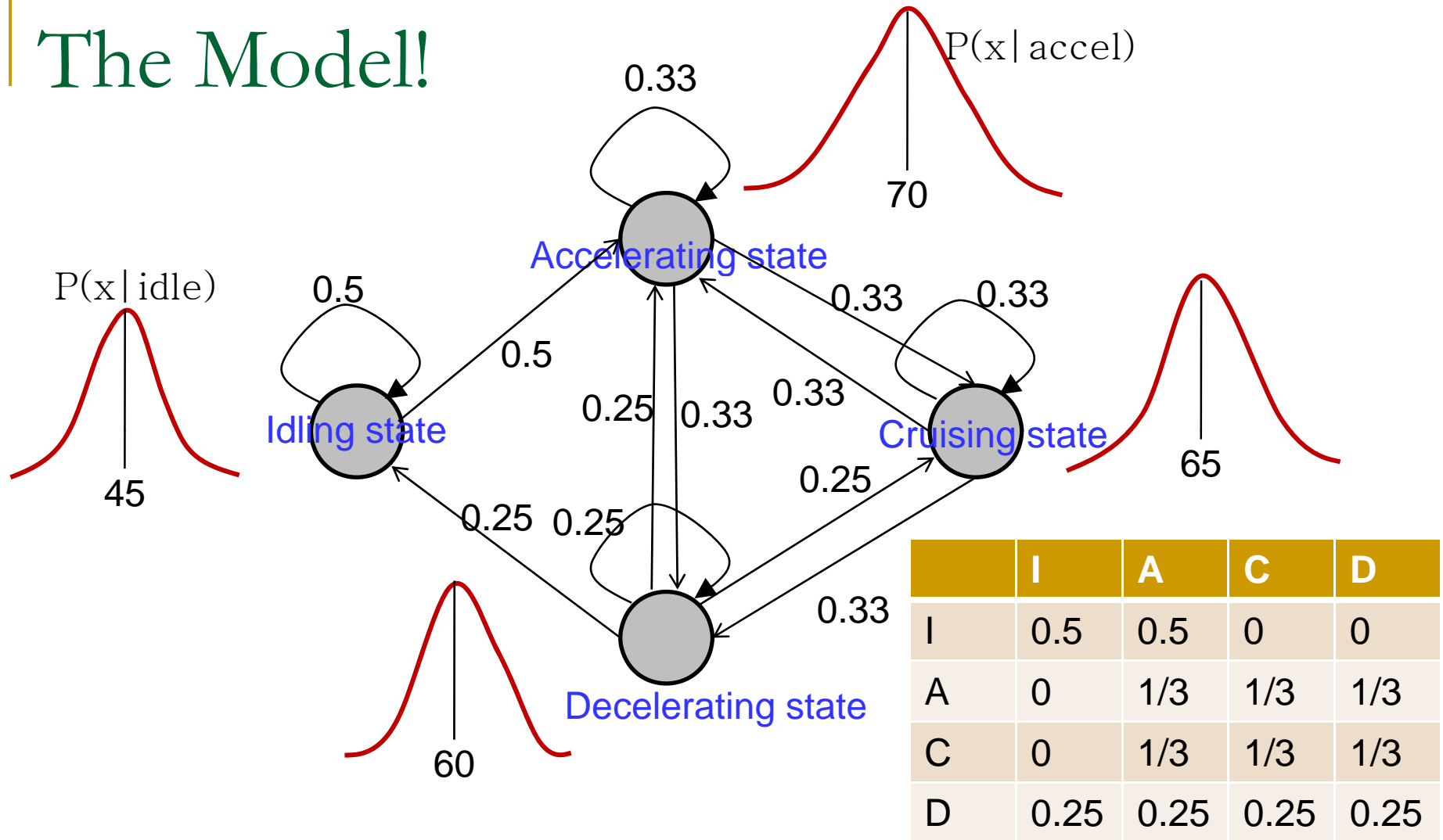
Class 27. 23 Nov 2010

# Recap: An automotive example



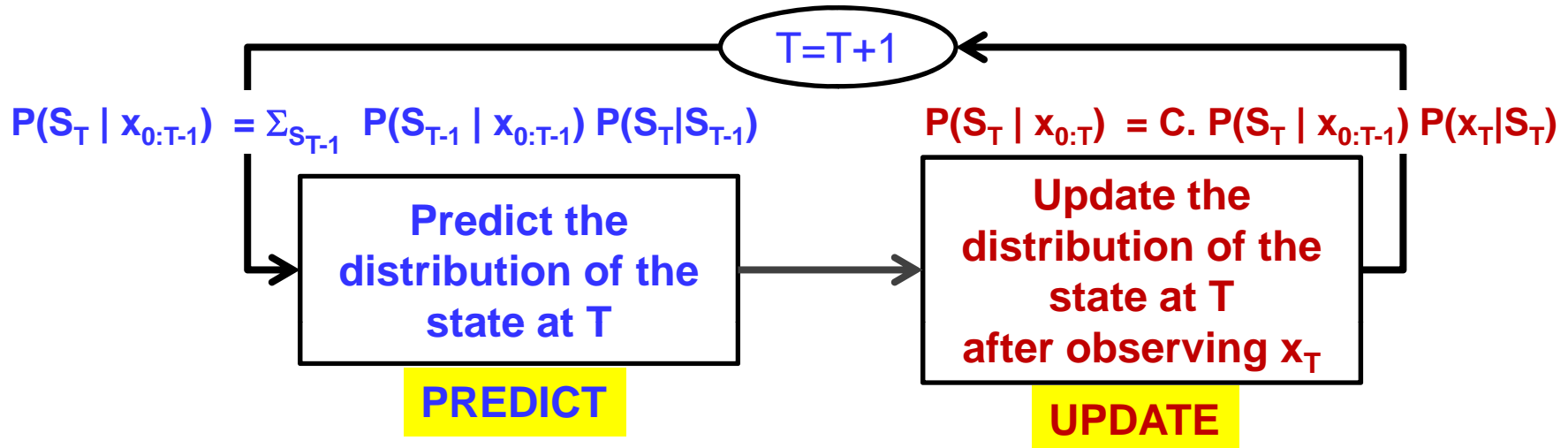
- Determine automatically, by only *listening* to a running automobile, if it is:
  - Idling; or
  - Travelling at constant velocity; or
  - Accelerating; or
  - Decelerating
- Assume (for illustration) that we only record energy level (SPL) in the sound
  - The SPL is measured once per second

# The Model!



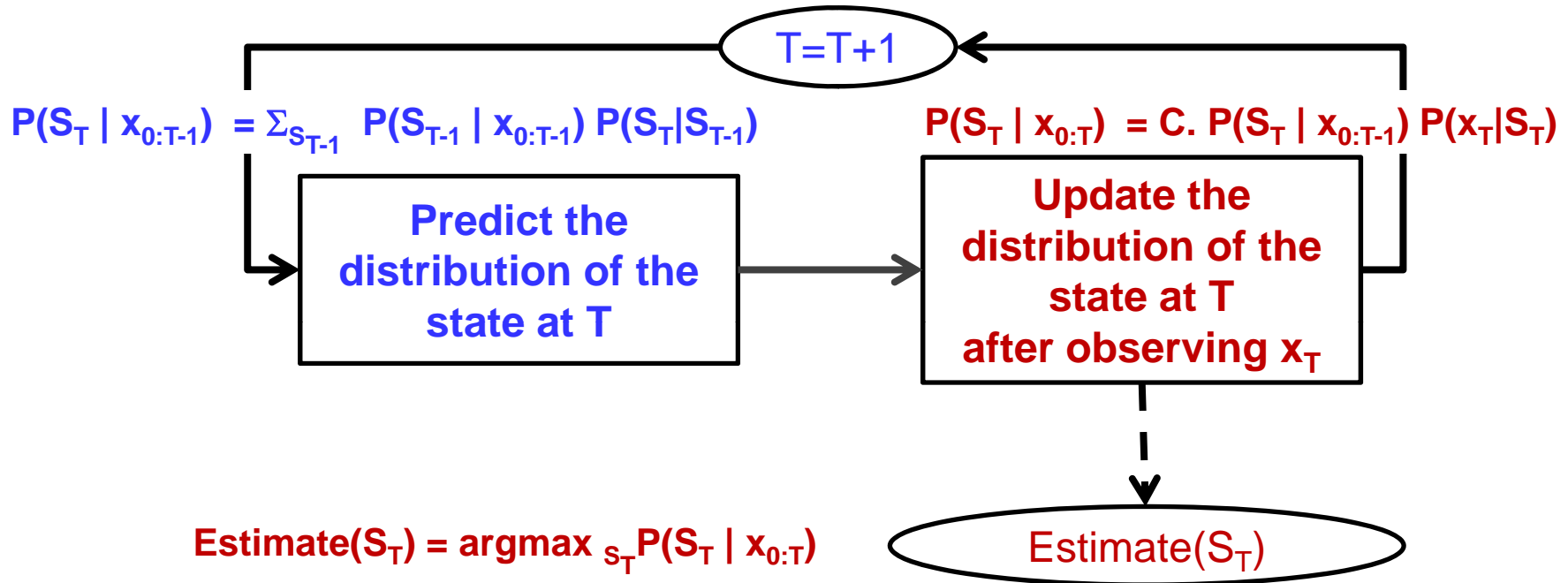
- The state-space model
  - Assuming all transitions from a state are equally probable

# Overall procedure



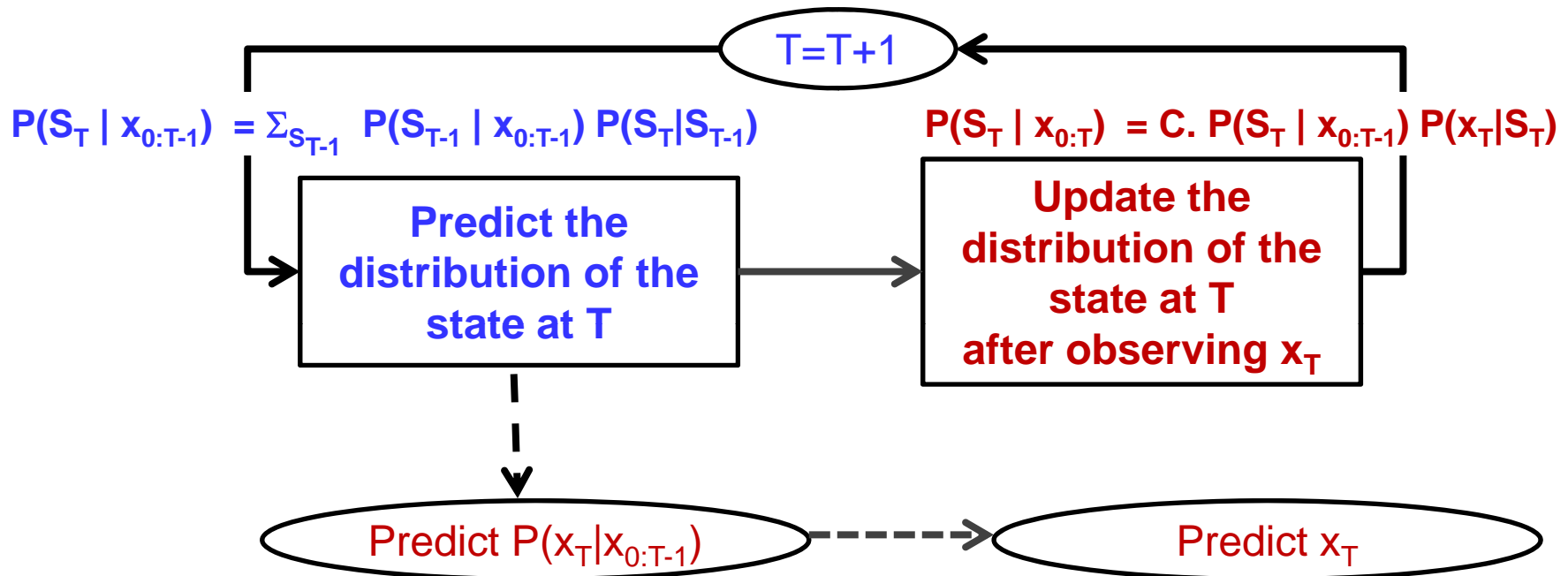
- At  $T=0$  the predicted state distribution is the initial state probability
- At each time  $T$ , the current estimate of the distribution over states considers *all* observations  $x_0 \dots x_T$ 
  - A natural outcome of the Markov nature of the model
- The prediction+update is identical to the forward computation for HMMs to within a normalizing constant

# Estimating the *state*



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

# Predicting the *next observation*



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

# Continuous state system

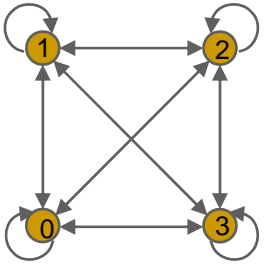


$$s_t = f(s_{t-1}, \varepsilon_t)$$

$$o_t = g(s_t, \gamma_t)$$

- 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)

# Discrete vs. Continuous State Systems



Prediction at time  $t$ :

$$P(s_t | O_{0:t-1}) = \sum_{s_{t-1}} P(s_{t-1} | O_{0:t-1}) P(s_t | s_{t-1})$$

Update after  $O_t$ :

$$P(s_t | O_{0:t}) = CP(s_t | O_{0:t-1}) P(O_t | s_t)$$

$$s_t = f(s_{t-1}, \varepsilon_t)$$

$$o_t = g(s_t, \gamma_t)$$

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

$$P(s_t | O_{0:t}) = CP(s_t | O_{0:t-1}) P(O_t | s_t)$$



# Special case: Linear Gaussian model

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

$$P(\varepsilon) = \frac{1}{\sqrt{(2\pi)^d |\Theta_\varepsilon|}} \exp\left(-0.5(\varepsilon - \mu_\varepsilon)^T \Theta_\varepsilon^{-1} (\varepsilon - \mu_\varepsilon)\right)$$

$$o_t = B_t s_t + \gamma_t$$

$$P(\gamma) = \frac{1}{\sqrt{(2\pi)^d |\Theta_\gamma|}} \exp\left(-0.5(\gamma - \mu_\gamma)^T \Theta_\gamma^{-1} (\gamma - \mu_\gamma)\right)$$

- A *linear* state dynamics equation
  - Probability of state driving term  $\varepsilon$  is Gaussian
  - Sometimes viewed as a driving term  $\mu_\varepsilon$  and additive zero-mean noise
- A *linear* observation equation
  - Probability of observation noise  $\gamma$  is Gaussian
- $A_t$ ,  $B_t$  and Gaussian parameters assumed known
  - May vary with time

# The Linear Gaussian model (KF)

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

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

$$P(o_t | s_t) = \text{Gaussian}(o_t; B_t s_t, \Theta_\gamma)$$

$$P(s_t | o_{0:t-1}) = \text{Gaussian}(s; \bar{s}_t, R_t)$$

$$\begin{aligned}\bar{s}_t &= \mu_\varepsilon + A_t \hat{s}_{t-1} \\ R_t &= \Theta_\varepsilon + A_t \hat{R}_{t-1} A_t^T\end{aligned}$$



$$P(s_t | o_{0:t}) = \text{Gaussian}(s; \hat{s}_t, \hat{R}_t)$$

$$\begin{aligned}\hat{s}_t &= \bar{s}_t + R_t B_t^T (B_t R_t B_t^T + \Theta_\gamma)^{-1} (o - B_t \bar{s}_t) \\ \hat{R}_t &= \left( I - R_t B_t^T (B_t R_t B_t^T + \Theta_\gamma)^{-1} B_t \right) R_t\end{aligned}$$

- Iterative prediction and update

# The Kalman filter

## ■ Prediction

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

$$R_t = \Theta_\varepsilon + A_{t-1} \hat{R}_{t-1} A_{t-1}^T$$

## ■ Update

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

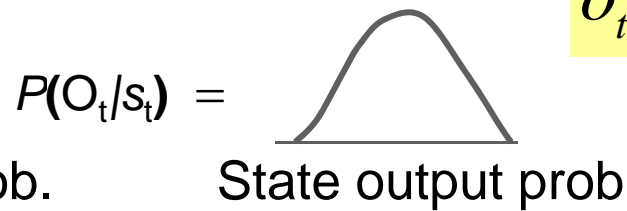
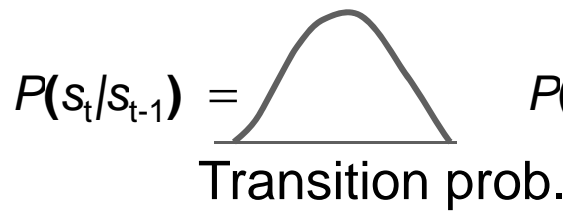
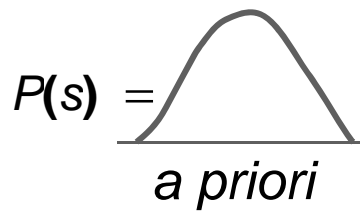
$$\hat{s}_t = \bar{s}_t + K_t (o_t - B_t \bar{s}_t)$$

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

# Linear Gaussian Model

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

$$o_t = B_t s_t + \gamma_t$$



$$P(s_0) = P(s)$$



$$P(s_0 | O_0) = C P(s_0) P(O_0 | s_0)$$



$$P(s_1 | O_0) = \int_{-\infty}^{\infty} P(s_0 | O_0) P(s_1 | s_0) ds_0$$



$$P(s_1 | O_{0:1}) = C P(s_1 | O_0) P(O_1 | s_0)$$



$$P(s_2 | O_{0:1}) = \int_{-\infty}^{\infty} P(s_1 | O_{0:1}) P(s_2 | s_1) ds_1$$



$$P(s_2 | O_{0:2}) = C P(s_2 | O_{0:1}) P(O_2 | s_2)$$

All distributions remain Gaussian

# Problems

$$s_t = f(s_{t-1}, \varepsilon_t)$$

$$o_t = g(s_t, \gamma_t)$$

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

# Problems

$$s_t = f(s_{t-1}, \varepsilon_t)$$

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  - Conventional Kalman update rules are no longer valid
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  - Gaussian based update rules no longer valid

# The problem with non-linear functions

$$s_t = f(s_{t-1}, \varepsilon_t)$$

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

$$o_t = g(s_t, \gamma_t)$$

$$P(s_t | o_{0:t}) = CP(s_t | o_{0:t-1}) P(o_t | s_t)$$

- 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

# The problem with nonlinearity

$$o_t = g(s_t, \gamma_t)$$

- The PDF may not have a closed form

$$P(o_t | s_t) = \sum_{\gamma: g(s_t, \gamma) = o_t} \frac{P_\gamma(\gamma)}{|J_{g(s_t, \gamma)}(o_t)|}$$

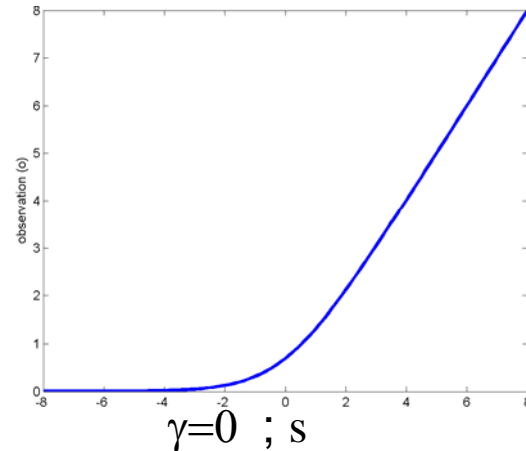
$$|J_{g(s_t, \gamma)}(o_t)| = \begin{vmatrix} \frac{\partial o_t(1)}{\partial \gamma(1)} & \dots & \frac{\partial o_t(1)}{\partial \gamma(n)} \\ \vdots & \ddots & \vdots \\ \frac{\partial o_t(n)}{\partial \gamma(1)} & \dots & \frac{\partial o_t(n)}{\partial \gamma(n)} \end{vmatrix}$$

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



# Example: a simple nonlinearity

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

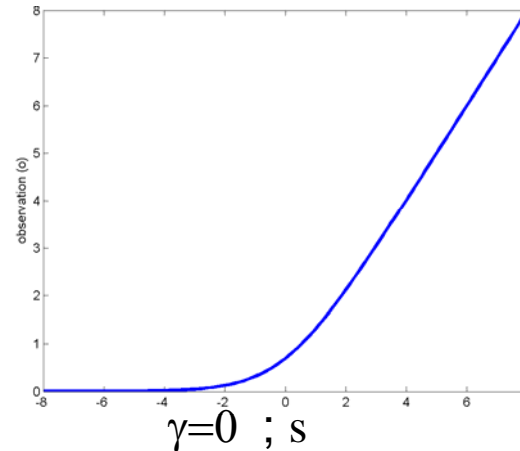


- $P(o|s) = ?$ 
  - Assume  $\gamma$  is Gaussian
  - $P(\gamma) = \text{Gaussian}(\gamma; \mu_\gamma, \Theta_\gamma)$

# Example: a simple nonlinearity

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

■  $P(o|s) = ?$

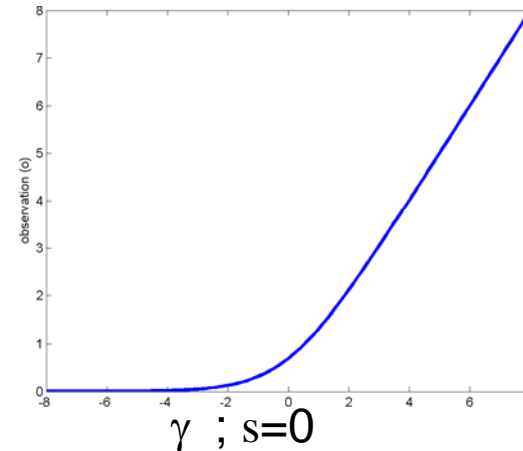


$$P(\gamma) = \text{Gaussian}(\gamma; \mu_\gamma, \Theta_\gamma)$$

$$P(o | s) = \text{Gaussian}(o; \mu_\gamma + \log(1 + \exp(s)), \Theta_\gamma)$$

# Example: At $T=0$ .

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



- Assume initial probability  $P(s)$  is Gaussian

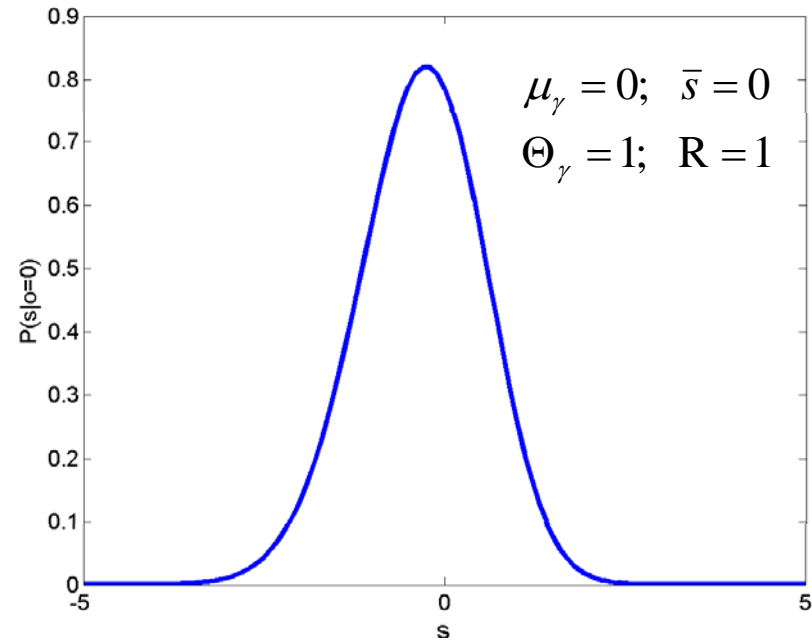
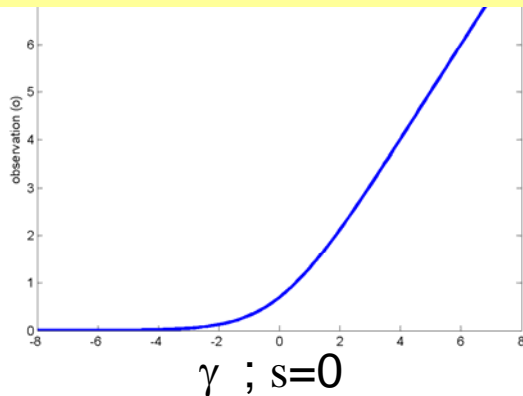
$$P(s_0) = P_0(s) = \text{Gaussian}(s; \bar{s}, R)$$

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

$$P(s_0 | o_0) = C \text{Gaussian}(o; \mu_\gamma + \log(1 + \exp(s_0)), \Theta_\gamma) \text{Gaussian}(s_0; \bar{s}, R)$$

# UPDATE: At T=0.

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



$$P(s_0 | o_0) = C \text{Gaussian}(o; \mu_\gamma + \log(1 + \exp(s_0)), \Theta_\gamma) \text{Gaussian}(s_0; \bar{s}, R)$$

$$P(s_0 | o_0) = C \exp \left( \begin{aligned} & -0.5(\mu_\gamma + \log(1 + \exp(s_0)) - o)^T \Theta_\gamma^{-1} (\mu_\gamma + \log(1 + \exp(s_0)) - o) \\ & - 0.5(s_0 - \bar{s})^T R^{-1} (s_0 - \bar{s}) \end{aligned} \right)$$

■ = Not Gaussian

# Prediction for $T = 1$

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

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

- Trivial, linear state transition equation

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

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

$$P(s_1 | o_0) = \int_{-\infty}^{\infty} C \exp\left(-0.5(\mu_\gamma + \log(1 + \exp(s_0)) - o)^T \Theta_\gamma^{-1} (\mu_\gamma + \log(1 + \exp(s_0)) - o) - 0.5(s_0 - \bar{s})^T R^{-1} (s_0 - \bar{s})\right) \exp\left(-0.5(s_1 - s_0)^T \Theta_\varepsilon^{-1} (s_1 - s_0)\right) ds_0$$

- = intractable

# Update at $T=1$ and later

- Update at  $T=1$

$$P(s_t | o_{0:t}) = CP(s_t | o_{0:t-1})P(o_t | s_t)$$

- Intractable

- Prediction for  $T=2$

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

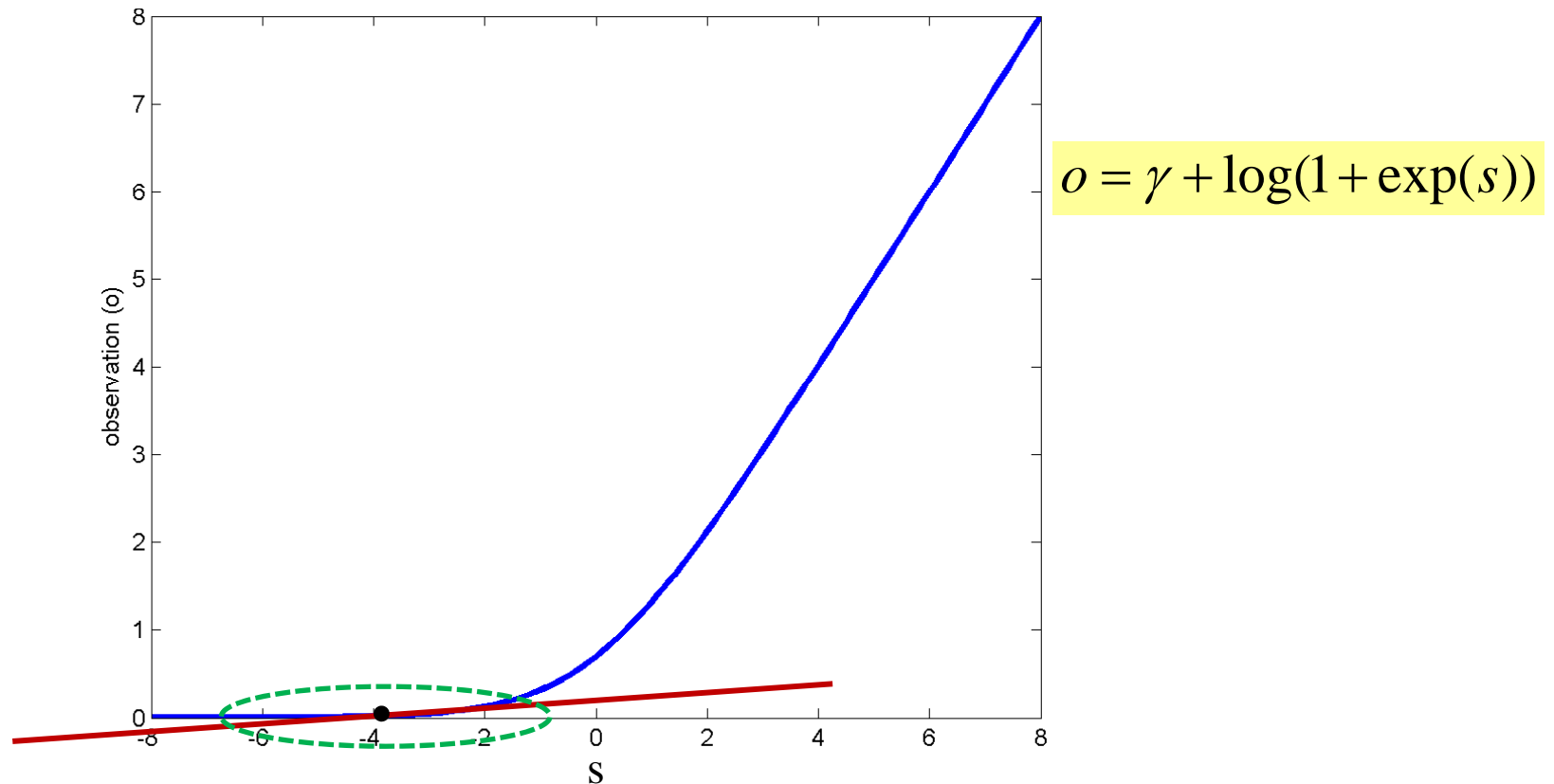
- Intractable

# The State prediction Equation

$$s_t = f(s_{t-1}, \varepsilon_t)$$

- 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

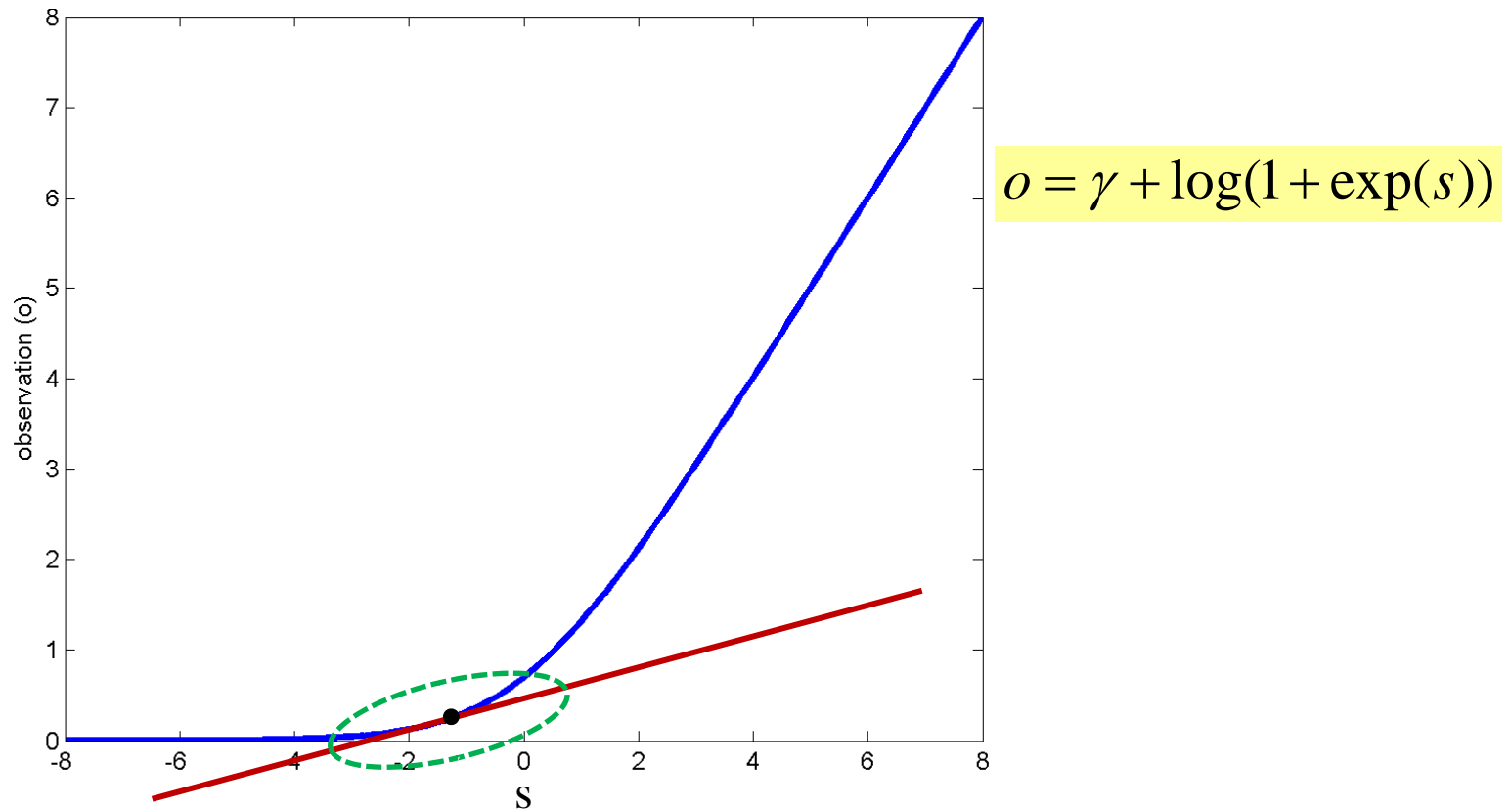
# Simplifying the problem: Linearize



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

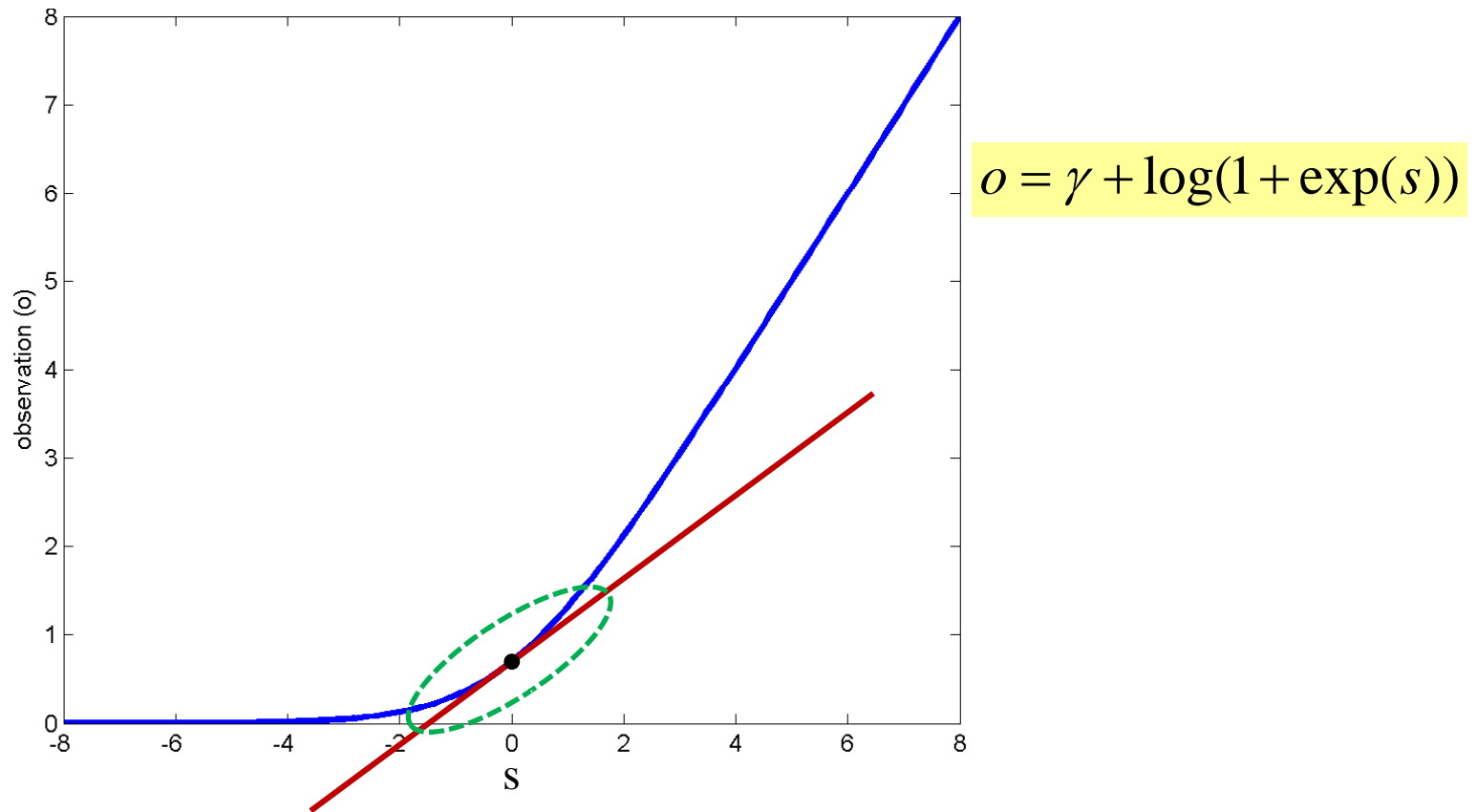


# Simplifying the problem: Linearize



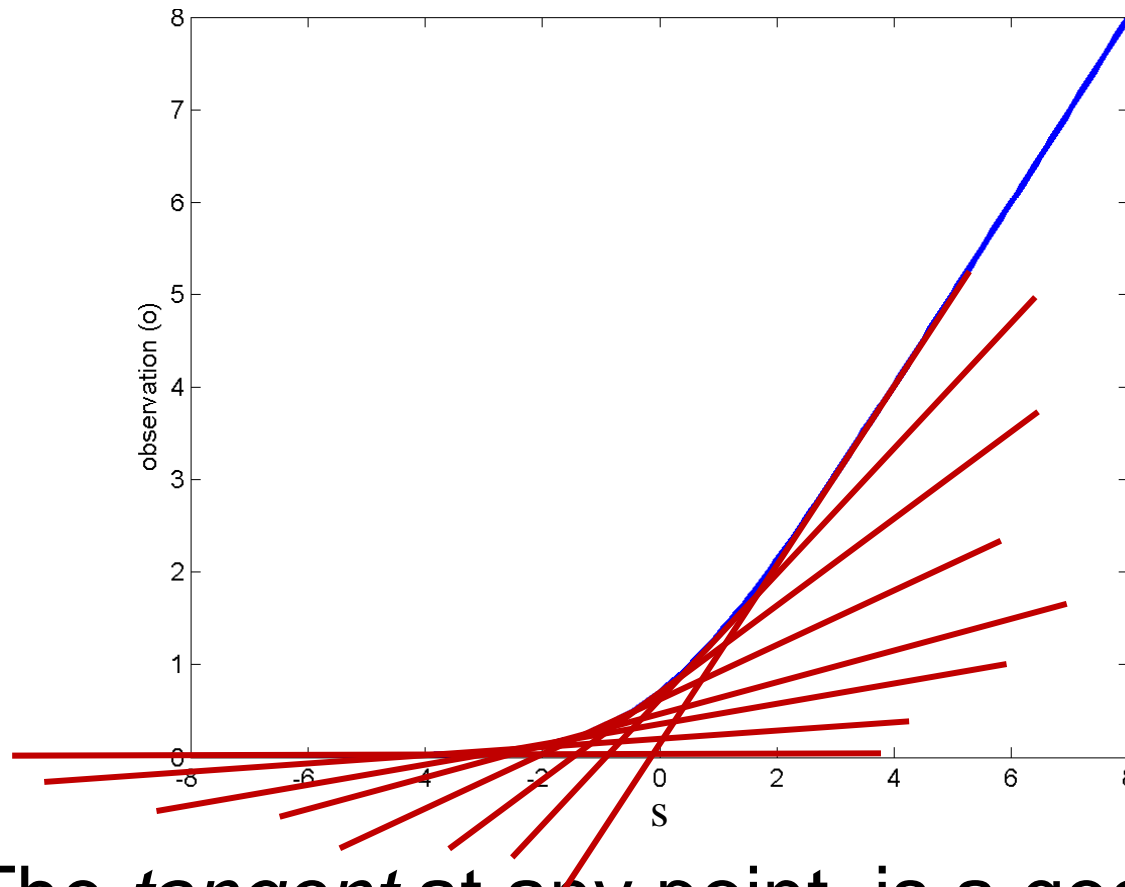
- The *tangent* at any point is a good *local* approximation if the function is sufficiently smooth

# Simplifying the problem: Linearize



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# Simplifying the problem: Linearize

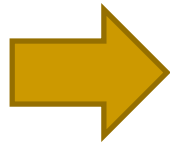


- The *tangent* at any point is a good *local* approximation if the function is sufficiently smooth

# Linearizing the observation function

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

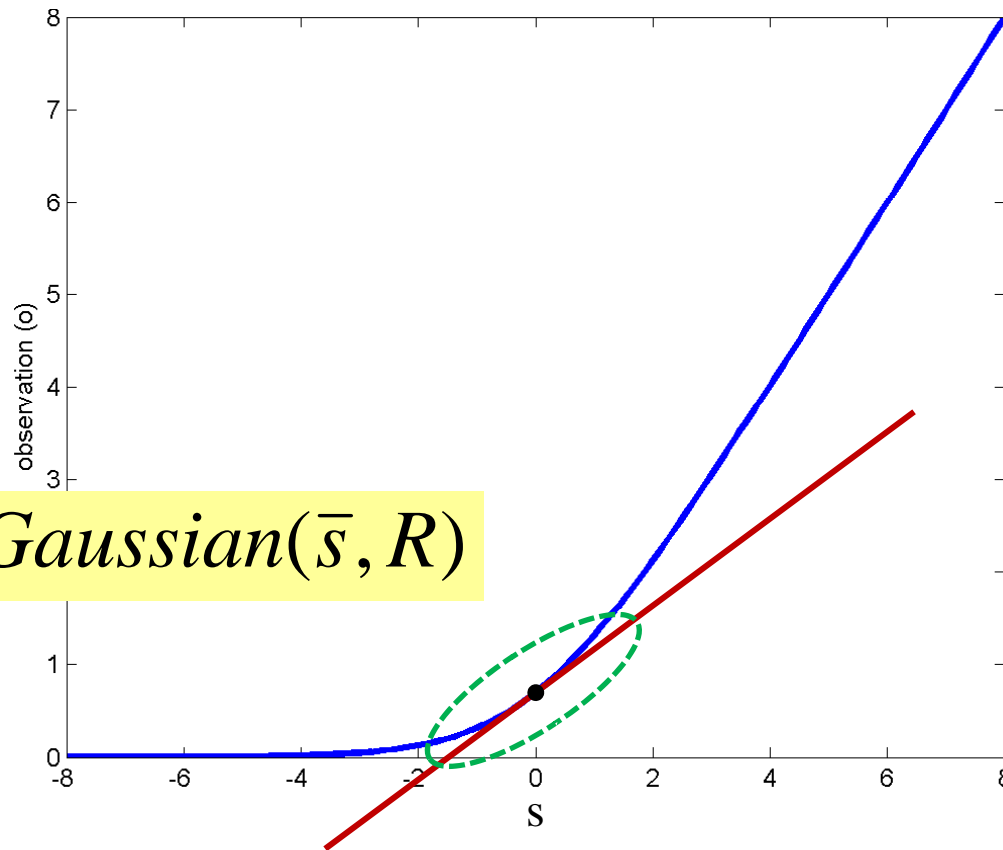
$$o = \gamma + g(s)$$



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

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

# Most probability is in the low-error region



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

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

# Linearizing the observation function

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

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

- Observation PDF is Gaussian

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

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

# UPDATE.

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

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

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

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

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

## ■ Gaussian!!

□ Note: This is actually only an approximation

# Prediction?

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

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

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

$$P(s_{t-1} | o_{0:t-1}) = \text{Gaussian}(s_{t-1}; \hat{s}_{t-1}, \hat{R}_{t-1})$$

$$s_t = f(s_{t-1}) + \varepsilon \quad \longrightarrow \quad s_t \approx \varepsilon + f(\hat{s}_{t-1}) + J_f(\hat{s}_{t-1})(s_{t-1} - \hat{s}_{t-1})$$

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



# Prediction

$$s_t = f(s_{t-1}) + \varepsilon \quad \Rightarrow \quad s_t \approx \varepsilon + f(\hat{s}_{t-1}) + J_f(\hat{s}_{t-1})(s_{t-1} - \hat{s}_{t-1})$$

$$P(s_{t-1} | o_{0:t-1}) = \text{Gaussian}(s_{t-1}; \hat{s}_{t-1}, \hat{R}_{t-1}) \quad P(\varepsilon) = \text{Gaussian}(\varepsilon; 0, \Theta_\varepsilon)$$

- The state transition probability is now:

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

- The predicted state probability is:

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

# Prediction

$$P(s_{t-1} | o_{0:t-1}) = \text{Gaussian}(s_{t-1}; \hat{s}_{t-1}, \hat{R}_{t-1})$$

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

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

$$P(s_t | o_{0:t-1}) = \int_{-\infty}^{\infty} \text{Gaussian}(s_{t-1}; \hat{s}_{t-1}, \hat{R}_{t-1}) \text{Gaussian}(s_t; f(\hat{s}_{t-1}) + J_f(\hat{s}_{t-1})(s_{t-1} - \hat{s}_{t-1}), \Theta_\varepsilon) ds_{t-1}$$

- The predicted state probability is:

$$P(s_t | o_{0:t-1}) = \text{Gaussian}(s_t; \hat{f}(s_{t-1}), J_f(\hat{s}_{t-1})\hat{R}_{t-1}J_f(\hat{s}_{t-1})^T + \Theta_\varepsilon)$$

- **Gaussian!!**

□ This is actually only an approximation

# The linearized prediction/update

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

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

- Given: two non-linear functions for state update and observation generation
- 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

# Linearized Prediction and Update

- Prediction for time t

$$P(s_t | o_{0:t-1}) = \text{Gaussian}(s_t; \bar{s}_t, R_t)$$

$$\bar{s}_t = f(\hat{s}_{t-1}) \quad R_t = J_f(\hat{s}_{t-1})\hat{R}_{t-1}J_f(\hat{s}_{t-1})^T + \Theta_\varepsilon$$

- Update at time t

$$P(s_t | o_{0:t}) = \text{Gaussian}(s_t; \hat{s}_t, \hat{R}_t)$$

$$\hat{s}_t = \bar{s}_t + R_t J_g(\bar{s}_t)^T (J_g(\bar{s}_t)R_t J_g(\bar{s}_t)^T + \Theta_\gamma)^{-1} (o_t - g(\bar{s}_t))$$
$$\hat{R}_t = \left( I - R_t J_g(\bar{s}_t)^T (J_g(\bar{s}_t)R_t J_g(\bar{s}_t)^T + \Theta_\gamma)^{-1} J_g(\bar{s}_t) \right) R_t$$

# Linearized Prediction and Update

- Prediction for time t

$$P(s_t | o_{0:t-1}) = \text{Gaussian}(s_t; \bar{s}_t, R_t)$$

$$\bar{s}_t = f(\hat{s}_{t-1}) \quad R_t = A_t \hat{R}_{t-1} A_t^T + \Theta_\varepsilon$$

$$A_t = J_f(\hat{s}_{t-1})$$

$$B_t = J_g(\bar{s}_t)$$

- Update at time t

$$P(s_t | o_{0:t}) = \text{Gaussian}(s_t; \hat{s}_t, \hat{R}_t)$$

$$\hat{s}_t = \bar{s}_t + R_t B_t^T (B_t R_t B_t^T + \Theta_\gamma)^{-1} (o_t - g(\bar{s}_t))$$

$$\hat{R}_t = \left( I - R_t B_t^T (B_t R_t B_t^T + \Theta_\gamma)^{-1} B_t \right) R_t$$

# The Extended Kalman filter

## ■ Prediction

$$\bar{s}_t = f(\hat{s}_{t-1})$$

$$A_t = J_f(\hat{s}_{t-1})$$

$$B_t = J_g(\bar{s}_t)$$

$$R_t = \Theta_\varepsilon + A_t \hat{R}_{t-1} A_t^T$$

## ■ Update

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

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

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

# The Kalman filter

## ■ Prediction

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

$$R_t = \Theta_\varepsilon + A_t \hat{R}_{t-1} A_t^T$$

## ■ Update

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

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

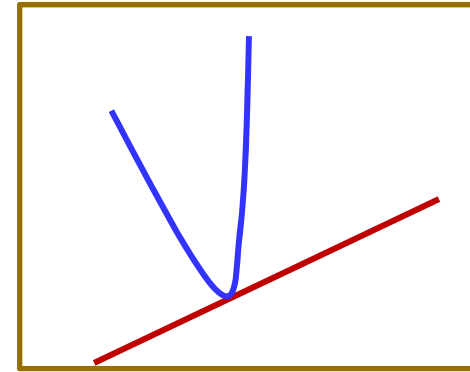
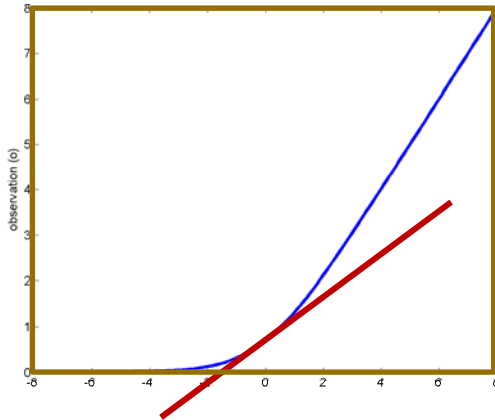
$$\hat{R}_t = (I - K_t B_t) R_t$$

# EKF<sub>s</sub>

- 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..



# 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:
  - Invariant extended Kalman filters (IEKF)
  - Unscented Kalman filters (UKF)

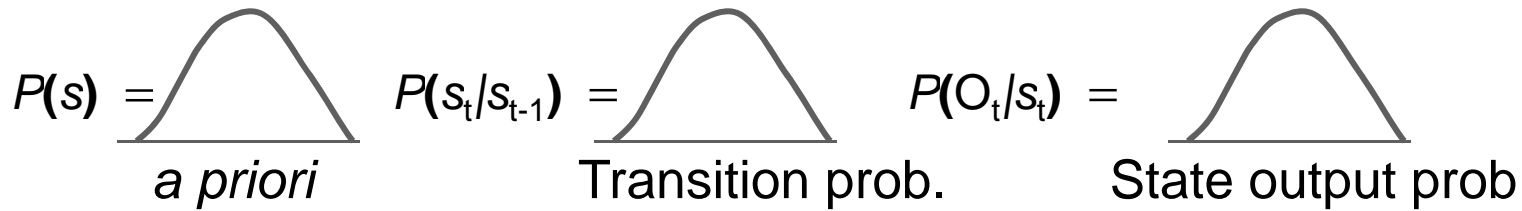
# A different problem: Non-Gaussian PDFs

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

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

- 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

# Linear Gaussian Model



$$P(s_0) = P(s)$$



$$P(s_0 | O_0) = C P(s_0) P(O_0 | s_0)$$



$$P(s_1 | O_0) = \int_{-\infty}^{\infty} P(s_0 | O_0) P(s_1 | s_0) ds_0$$



$$P(s_1 | O_{0:1}) = C P(s_1 | O_0) P(O_1 | s_0)$$



$$P(s_2 | O_{0:1}) = \int_{-\infty}^{\infty} P(s_1 | O_{0:1}) P(s_2 | s_1) ds_1$$



$$P(s_2 | O_{0:2}) = C P(s_2 | O_{0:1}) P(O_2 | s_2)$$

All distributions remain Gaussian

# 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

# A simple case



$$o_t = Bs_t + \gamma$$

$$P(\gamma) = \sum_{i=0}^1 w_i \text{Gaussian}(\gamma; \mu_i, \Theta_i)$$

- $P(\gamma)$  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!

$$P(o_t | s_t) = P(\gamma = o_t - Bs_t) = \sum_{i=0}^1 w_i \text{Gaussian}(o; \mu_i + Bs_t, \Theta_i)$$

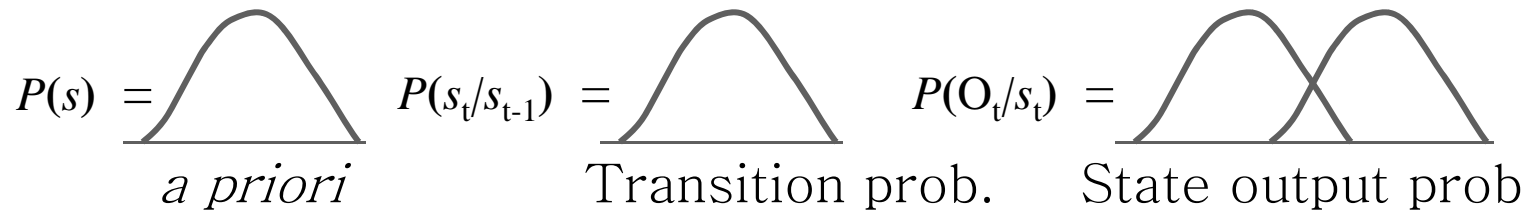


$P(\gamma)$



$P(o_t | s_t)$

# When distributions are not Gaussian



$$P(s_0) = P(s)$$



# When distributions are not Gaussian

$$P(s) = \text{[Gaussian curve]}$$

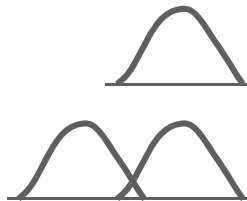
*a priori*

$$P(s_t/s_{t-1}) = \text{[Gaussian curve]}$$

Transition prob.

$$P(O_t/s_t) = \text{[Two overlapping Gaussian curves]}$$

State output prob

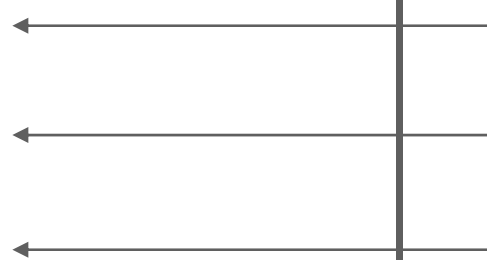
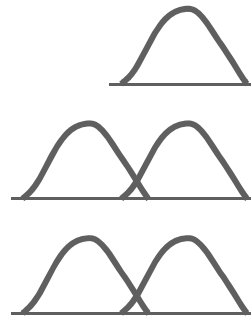
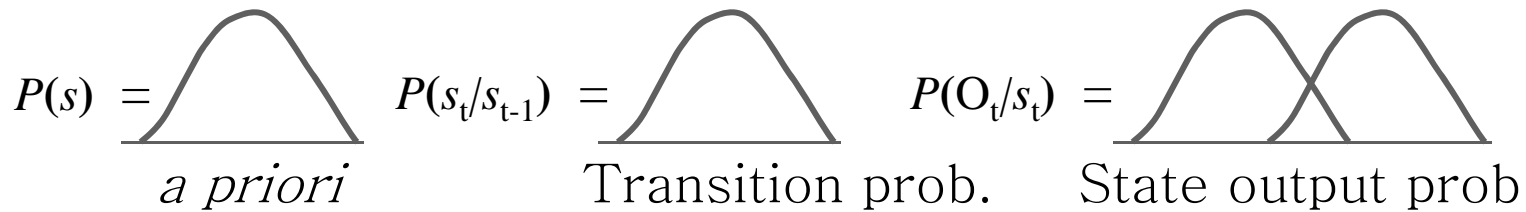


$$P(s_0) = P(s)$$



$$P(s_0 | O_0) = C P(s_0) P(O_0 | s_0)$$

# When distributions are not Gaussian



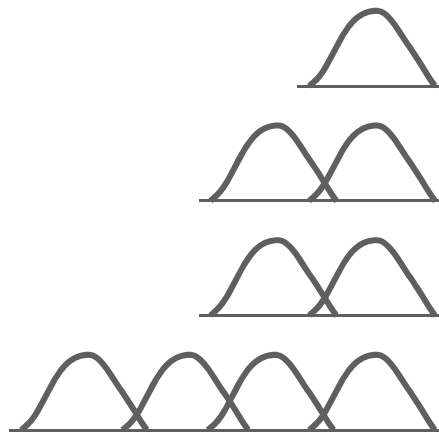
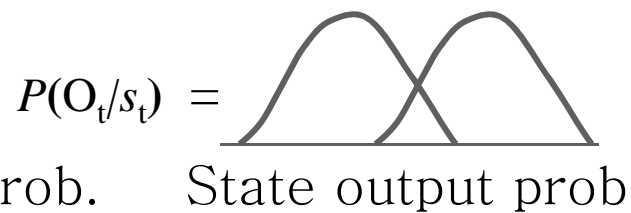
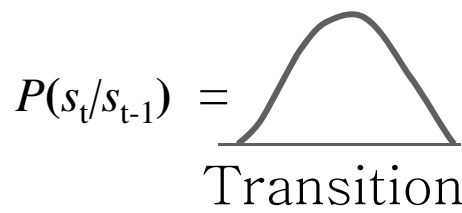
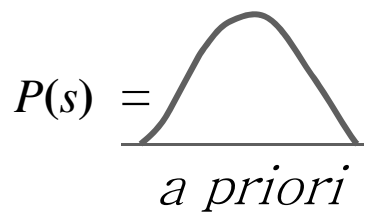
$$P(s_0) = P(s)$$

$$P(s_0 | O_0) = C P(s_0) P(O_0 | s_0)$$

$$P(s_1 | O_0) = \int_{-\infty}^{\infty} P(s_0 | O_0) P(s_1 | s_0) ds_0$$



# When distributions are not Gaussian



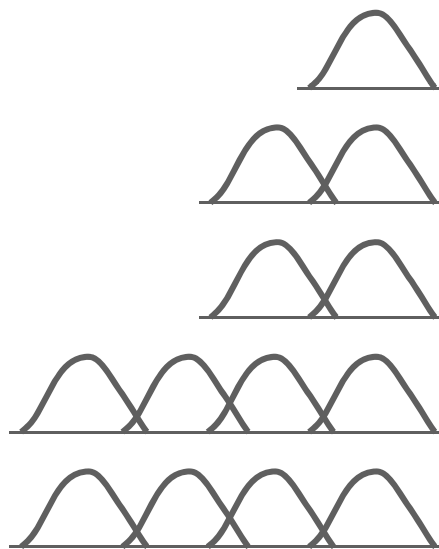
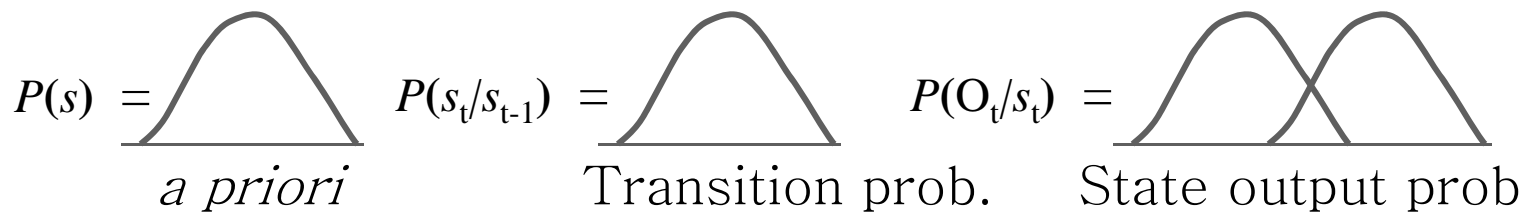
$$P(s_0) = P(s)$$

$$P(s_0 | O_0) = C P(s_0) P(O_0 | s_0)$$

$$P(s_1 | O_0) = \int_{-\infty}^{\infty} P(s_0 | O_0) P(s_1 | s_0) ds_0$$

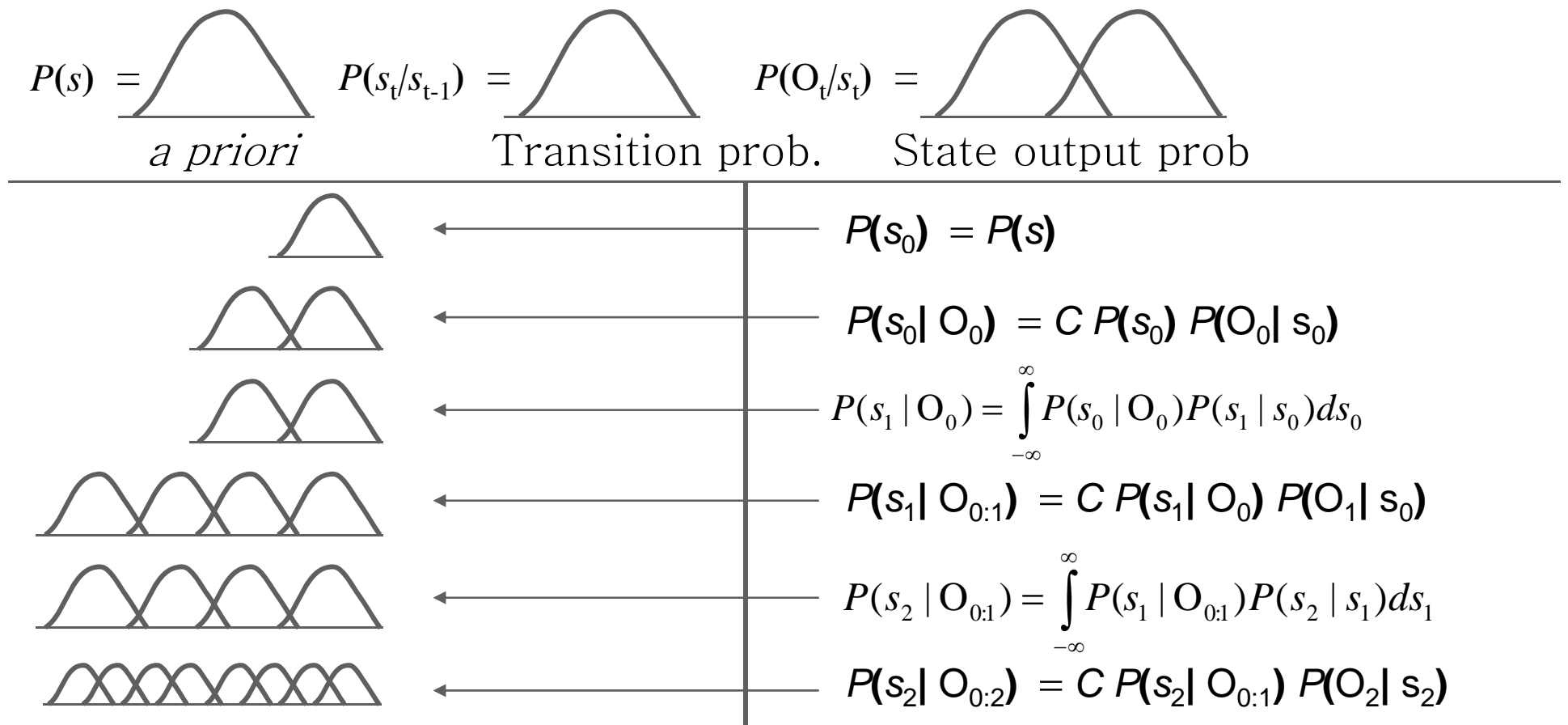
$$P(s_1 | O_{0:1}) = C P(s_1 | O_0) P(O_1 | s_0)$$

# When distributions are not Gaussian



←	$P(s_0) = P(s)$
←	$P(s_0   O_0) = C P(s_0) P(O_0   s_0)$
←	$P(s_1   O_0) = \int_{-\infty}^{\infty} P(s_0   O_0) P(s_1   s_0) ds_0$
←	$P(s_1   O_{0:1}) = C P(s_1   O_0) P(O_1   s_0)$
←	$P(s_2   O_{0:1}) = \int_{-\infty}^{\infty} P(s_1   O_{0:1}) P(s_2   s_1) ds_1$

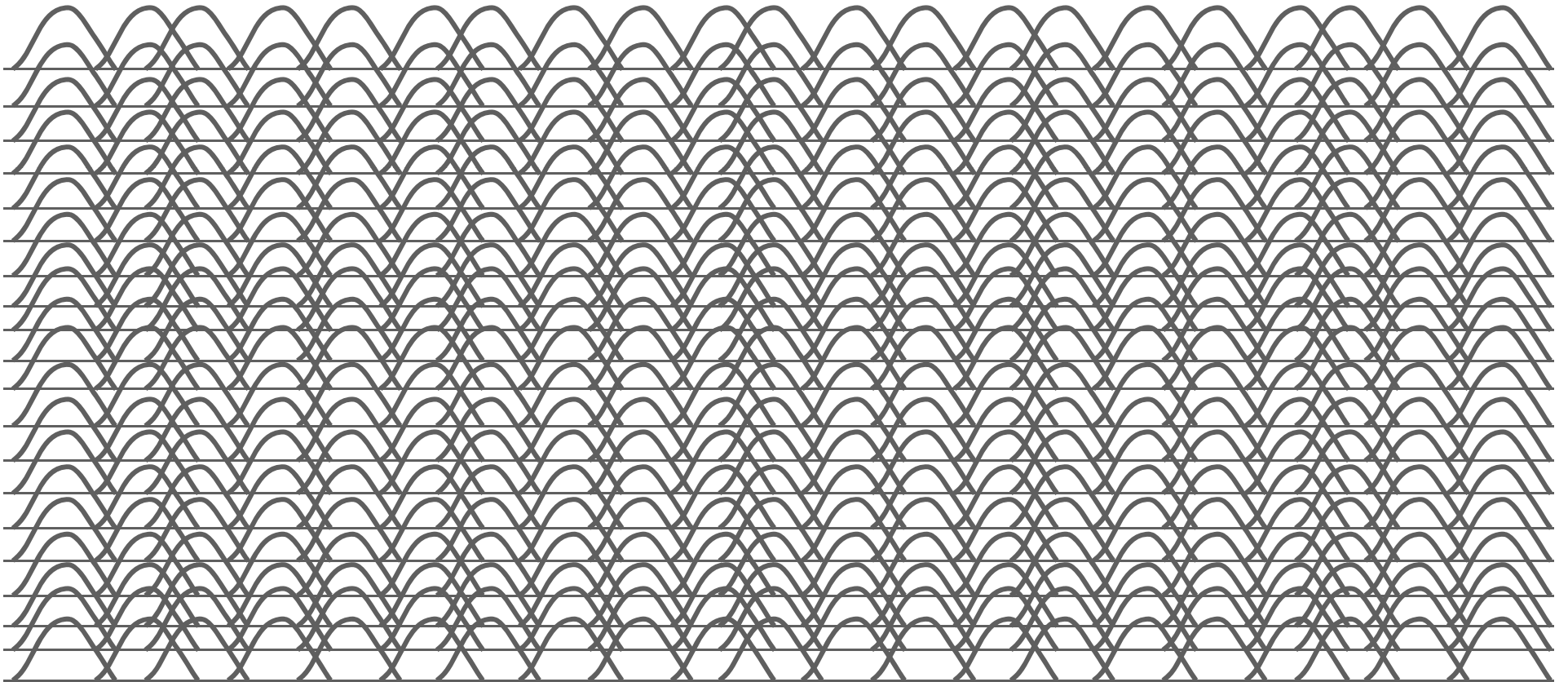
# When distributions are not Gaussian



When  $P(O_t/s_t)$  has more than one Gaussian, after only a few time steps...

# When distributions are not Gaussian

$$P(s_t | O_{0:t}) =$$



We have too many Gaussians for comfort..

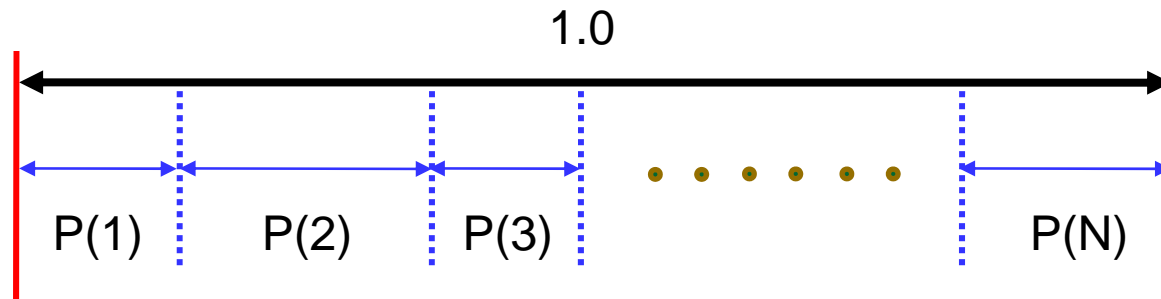
# Related Topic: How to sample from a Distribution?

- “Sampling from a Distribution  $P(x; \Gamma)$  with parameters  $\Gamma$ ”
- Generate random numbers such that
  - The distribution of a large number of generated numbers is  $P(x; \Gamma)$
  - The parameters of the distribution are  $\Gamma$
- 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

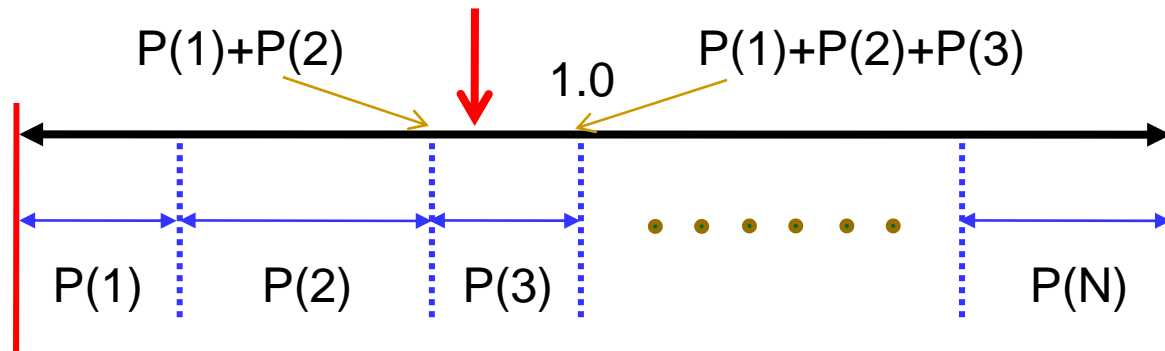
- Given a multinomial over  $N$  symbols, with probability of  $i^{\text{th}}$  symbol =  $P(i)$
- Randomly generate symbols from this distribution
- Can be done by sampling from a uniform distribution

# Sampling a multinomial



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

# 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  $i^{\text{th}}$  segment, select the  $i^{\text{th}}$  symbol



## 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  $\mu$ , std dev  $\sigma$ :

$$x = \sum_{i=1}^{12} r_i - 6$$

- Matlab :  $x = \mu + \text{randn} * \sigma$

- “randn” draws from a Gaussian of mean=0, variance=1

## Related Topic: Sampling from a Gaussian

- Multivariate (d-dimensional) Gaussian with mean  $\mu$  and covariance  $\Theta$ 
  - Compute eigen value matrix  $\Lambda$  and eigenvector matrix  $E$  for  $\Theta$ 
    - $\Theta = E \Lambda E^T$
  - Generate d 0-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  $\Lambda$  and  $E$ , add  $\mu$

$$Y = \mu + E \text{sqrt}(\Lambda) X$$

# Sampling from a Gaussian Mixture

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

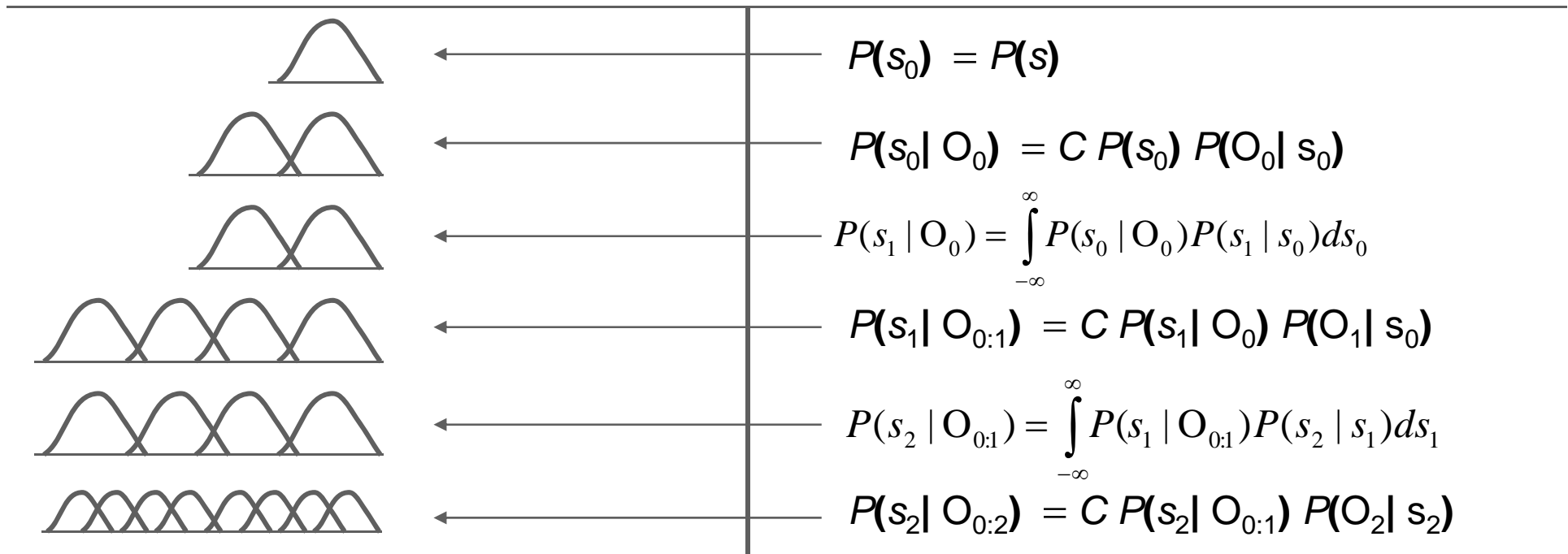
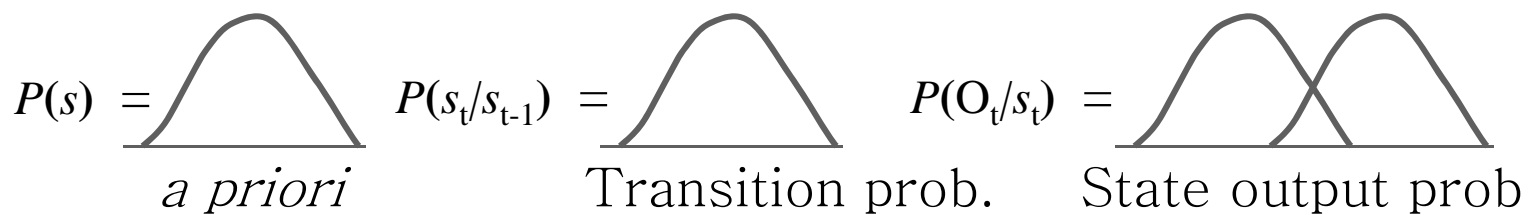
- Select a Gaussian by sampling the multinomial distribution of weights:

$$j \sim \text{multinomial}(w_1, w_2, \dots)$$

- Sample from the selected Gaussian

$$\text{Gaussian}(X; \mu_j, \Theta_j)$$

# Returning to our problem:

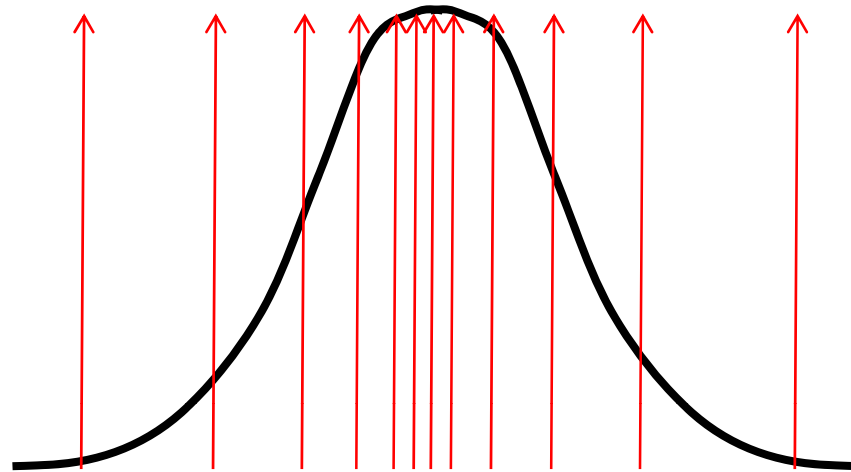


When  $P(O_t/s_t)$  has more than one Gaussian, after only a few time steps...

# The problem of the exploding distribution

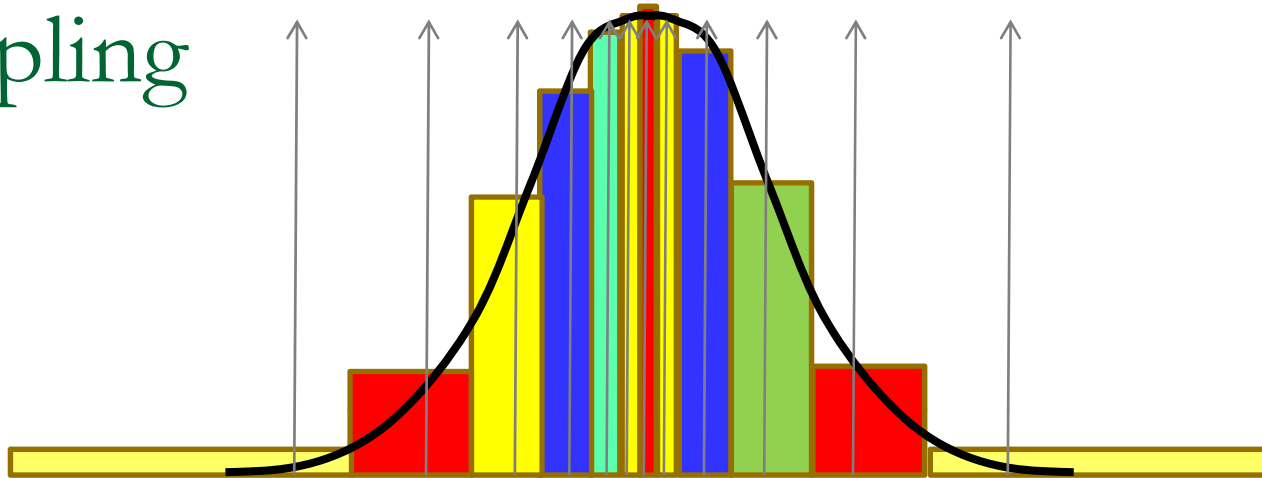
- The complexity of the distribution increases exponentially with time
- This is a consequence of having a *continuous* state space
  - Only Gaussian PDFs propagate without increase of complexity
- *Discrete-state* systems do not have this problem
  - The number of states in an HMM stays fixed
  - However,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

# Discrete approximation to a distribution



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

# Discrete approximation: Random sampling



- 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)

$$P(x) \approx \frac{1}{M} \sum_{i=0}^{M-1} \delta(x - x_i)$$

## Note: Properties of a discrete distribution

$$P(x) \approx \frac{1}{M} \sum_{i=0}^{M-1} \delta(x - x_i)$$

$$P(x)P(y|x) \propto \sum_{i=0}^{M-1} P(y|x_i)\delta(x - x_i)$$

- The product of a discrete distribution with another distribution is simply a weighted discrete probability

$$P(x) \approx \sum_{i=0}^{M-1} w_i \delta(x - x_i)$$

$$\int_{-\infty}^{\infty} P(x)P(y|x)dx = \sum_{i=0}^{M-1} w_i P(y|x_i)$$

- The integral of the product is a mixture distribution



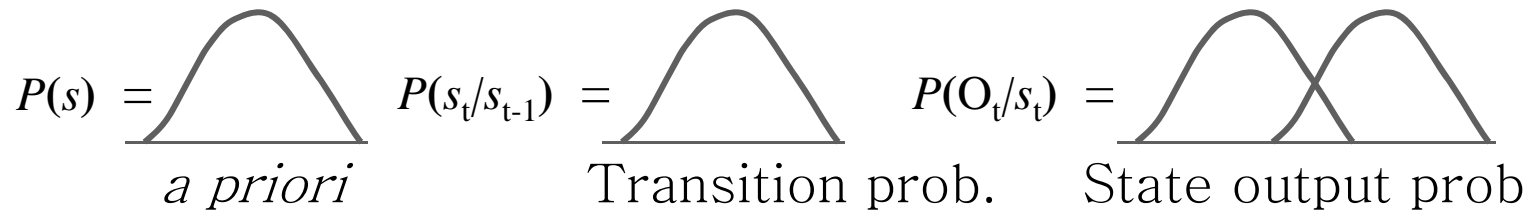
# Discretizing the state space

- At each time, discretize the predicted state space

$$P(s_t | o_{0:t}) \approx \frac{1}{M} \sum_{i=0}^{M-1} \delta(s_t - s_i)$$

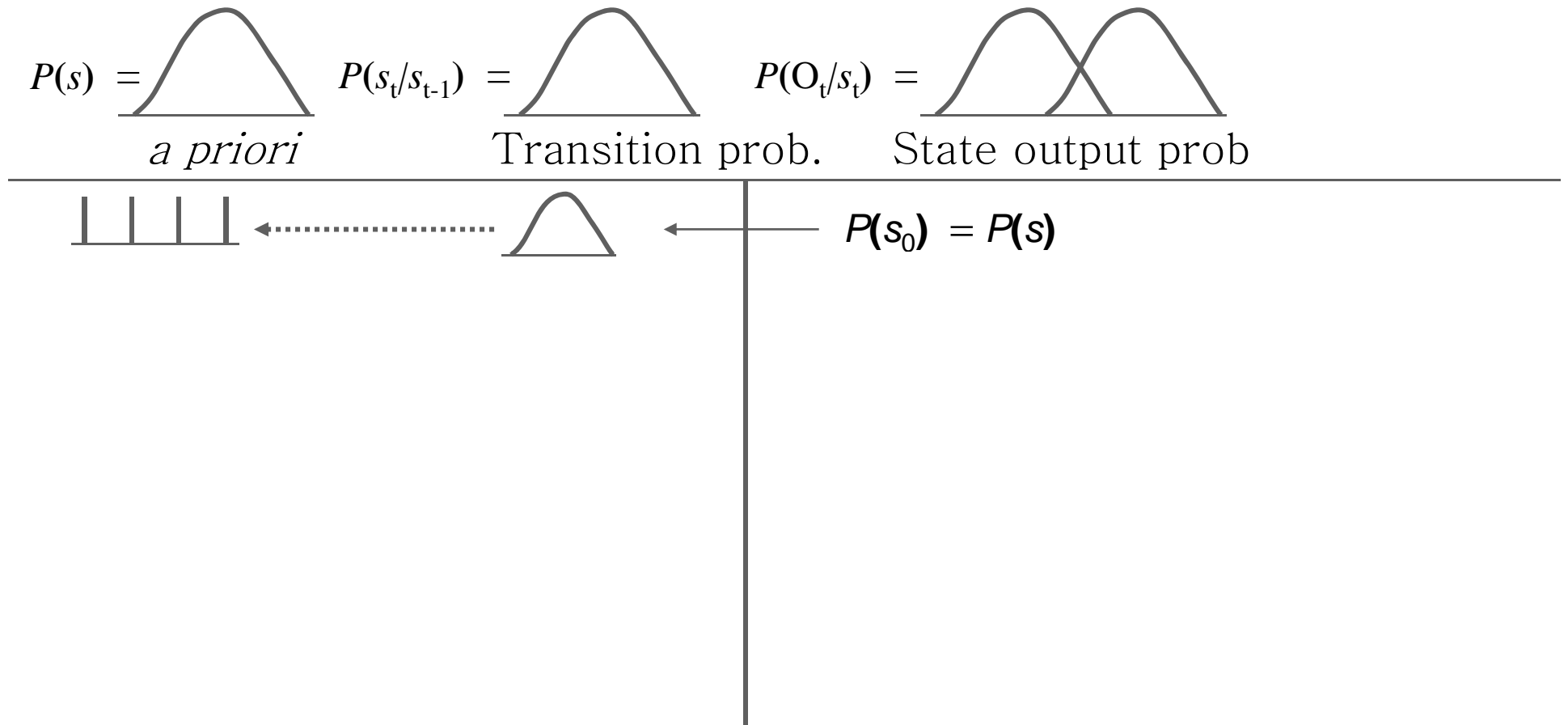
- $s_i$  are randomly drawn samples from  $P(s_t | o_{0:t})$
- Propagate the discretized distribution

# Particle Filtering



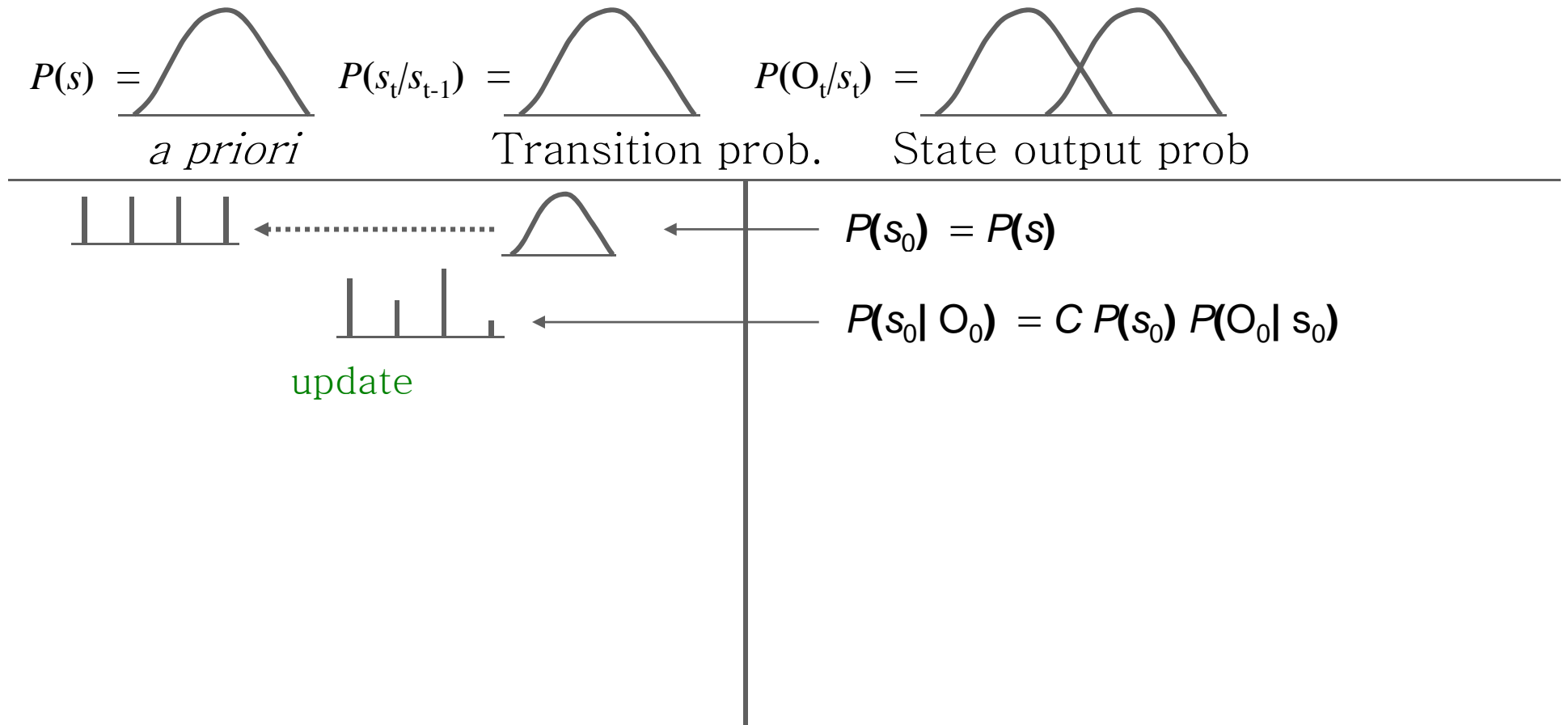
Assuming that we only generate *FOUR* samples from the predicted distributions

# Particle Filtering



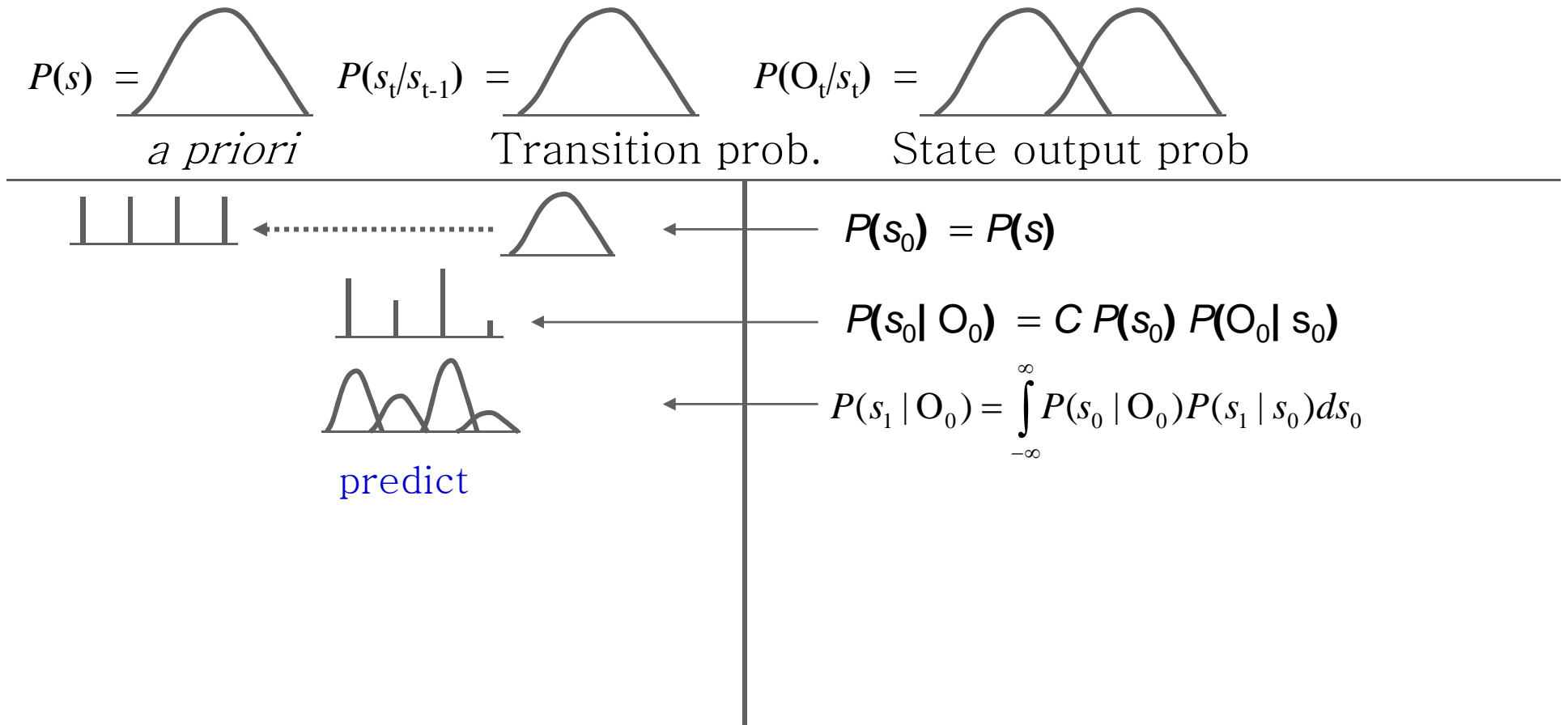
Assuming that we only generate *FOUR* samples from the predicted distributions

# Particle Filtering



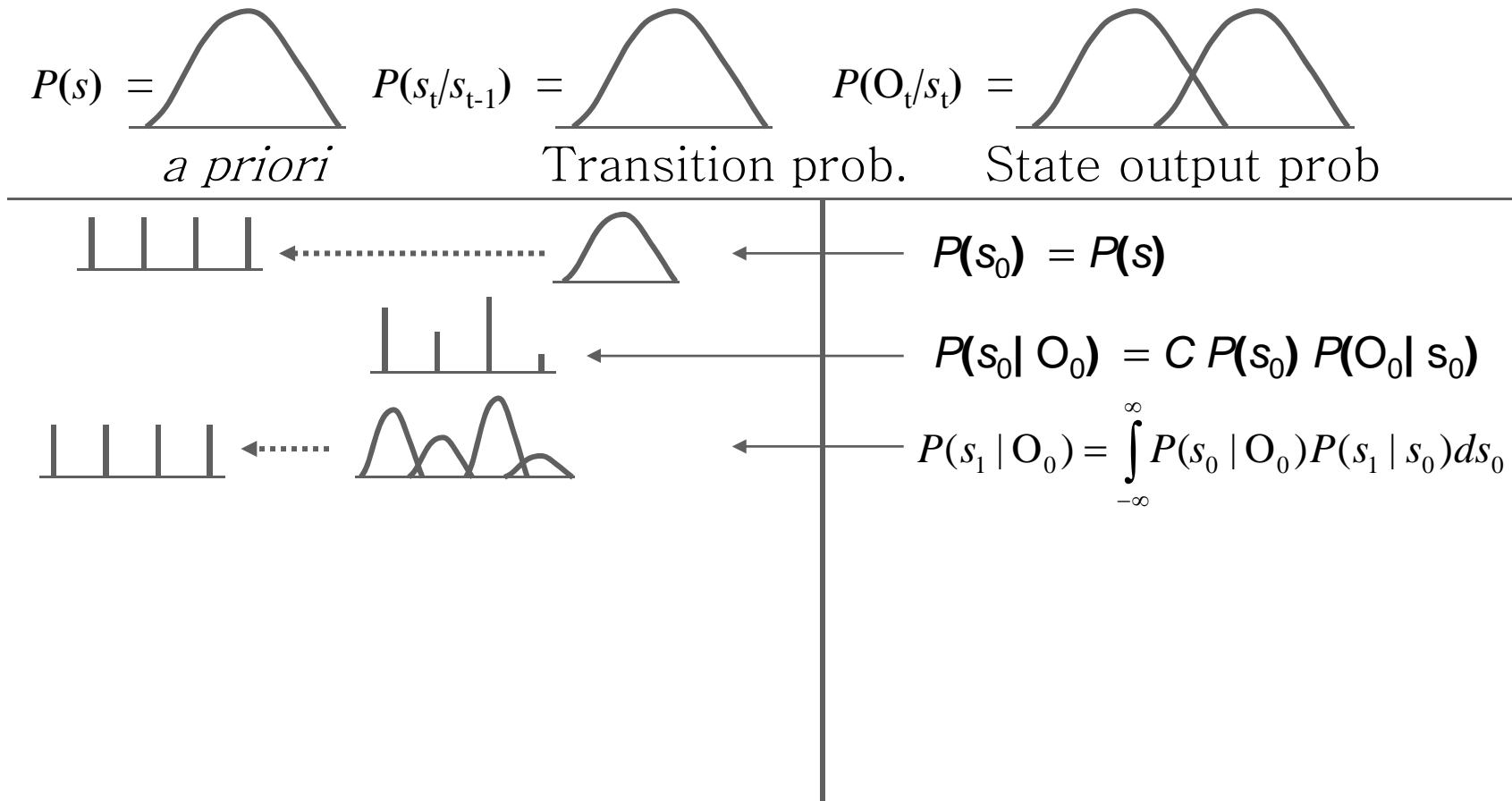
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# Particle Filtering



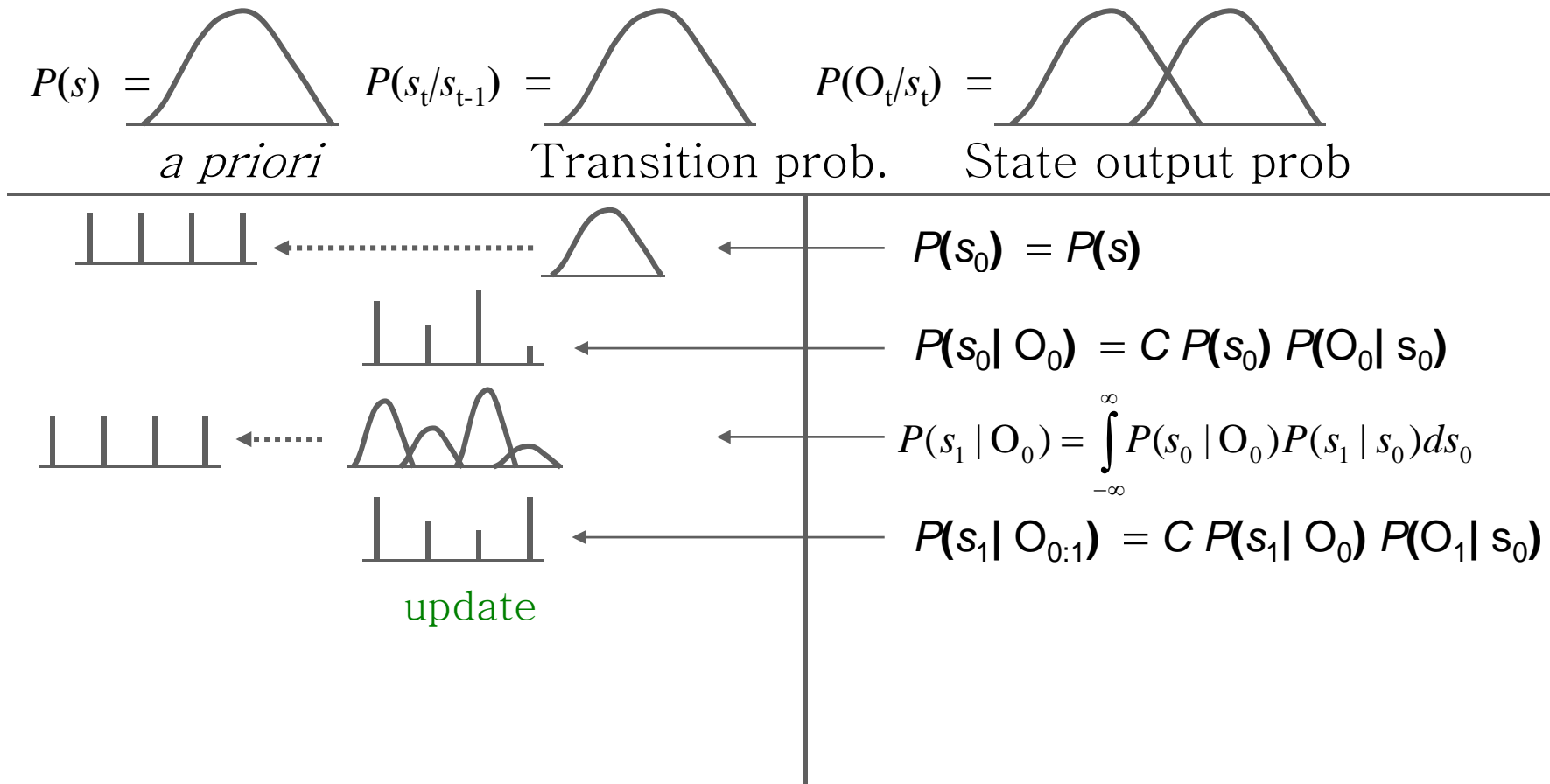
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# Particle Filtering



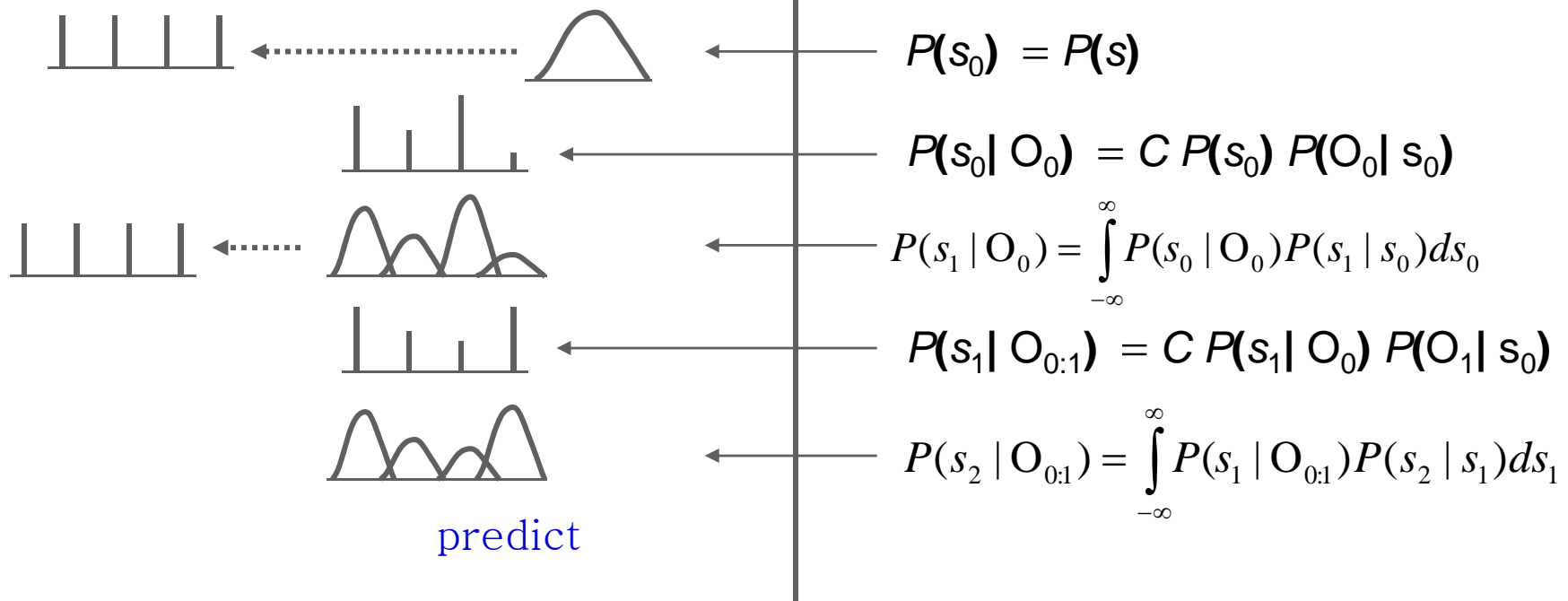
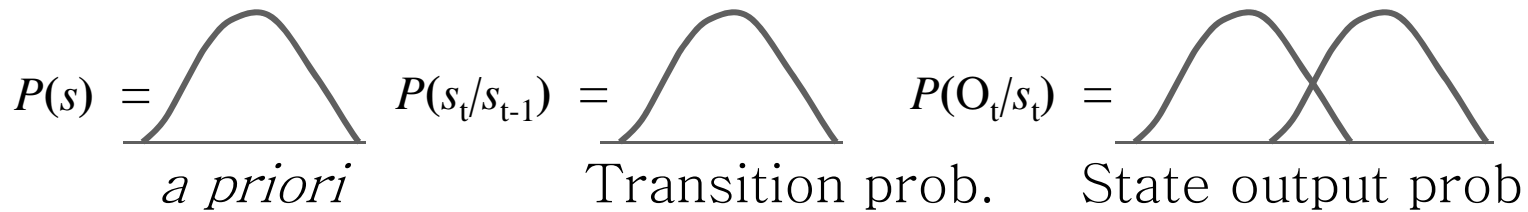
Assuming that we only generate *FOUR* samples from the predicted distributions

# Particle Filtering



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

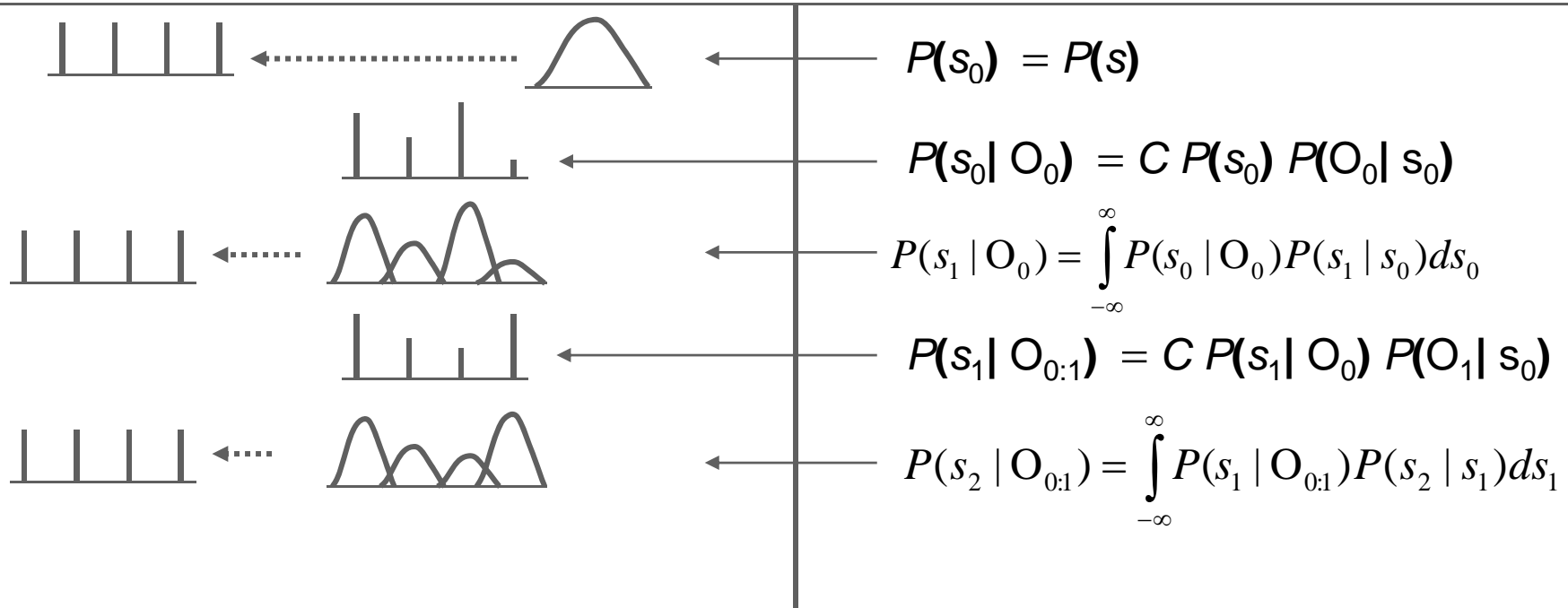
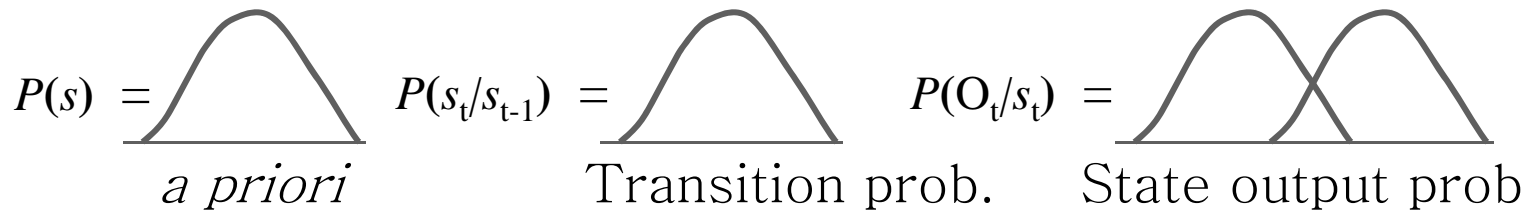
# Particle Filtering



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

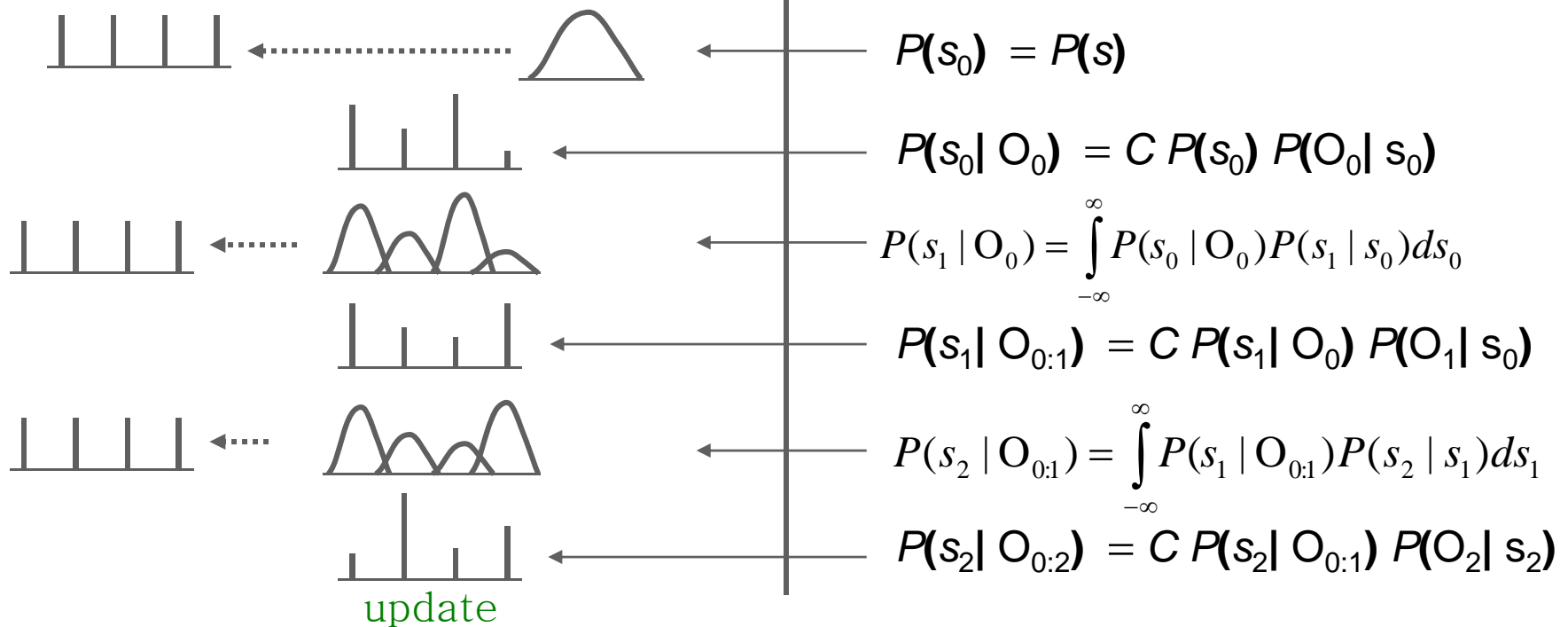
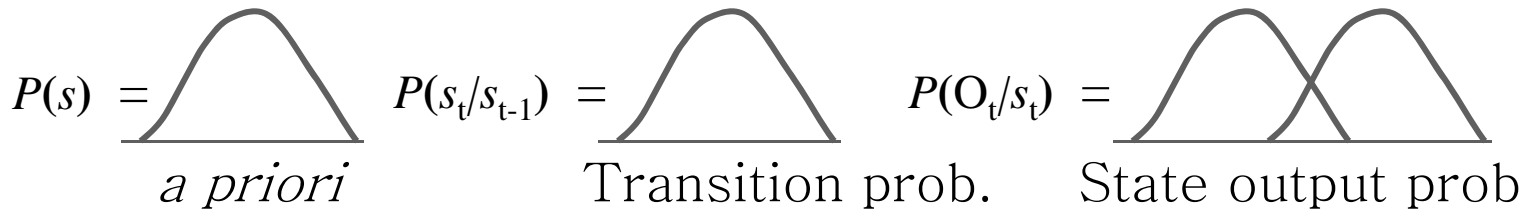


# Particle Filtering



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

# Particle Filtering

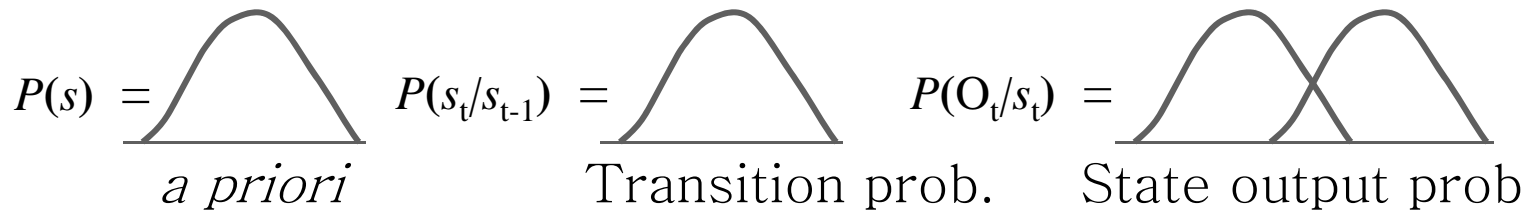


Assuming that we only generate **FOUR** samples from the predicted distributions

# 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



predict



update

Prediction at time t:

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

Update at time t:

$$P(s_t | O_{0:t}) = CP(s_t | O_{0:t-1}) P(O_t | s_t)$$

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

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

$$P_\gamma(\gamma)$$

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

$$P_\varepsilon(\varepsilon)$$

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

$$P(s_0 | o_{-1}) = P(s_0) \approx \frac{1}{M} \sum_{i=0}^{M-1} \delta(s_0 - \bar{s}_i^0) \text{ where } \bar{s}_i^0 \leftarrow P_0(s)$$

- Update the state distribution with the observation

$$P(s_t | o_{0:t}) = C \sum_{i=0}^{M-1} P_\gamma(o_t - g(\bar{s}_i^t)) \delta(s_t - \bar{s}_i^t)$$

$$C = \frac{1}{\sum_{i=0}^{M-1} P_\gamma(o_t - g(\bar{s}_i^t))}$$

# Particle Filtering

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

$$P_\gamma(\gamma)$$

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

$$P_\varepsilon(\varepsilon)$$

- Predict the state distribution at the next time

$$P(s_t | o_{0:t-1}) = C \sum_{i=0}^{M-1} P_\gamma(o_{t-1} - g(\bar{s}_i^{t-1})) P_\varepsilon(s_t - f(\bar{s}_i^{t-1}))$$

- Sample the predicted state distribution

$$P(s_t | o_{0:t-1}) \approx \frac{1}{M} \sum_{i=0}^{M-1} \delta(s_t - \bar{s}_i^t) \text{ where } \bar{s}_i^t \leftarrow P(s_t | o_{0:t-1})$$

# Particle Filtering

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

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

- Predict the state distribution at t

$$P(s_t | o_{0:t-1}) = C \sum_{i=0}^{M-1} P_\gamma(o_{t-1} - g(\bar{s}_i^{t-1})) P_\varepsilon(s_t - f(\bar{s}_i^{t-1}))$$

- Sample the predicted state distribution at t

$$P(s_t | o_{0:t-1}) \approx \frac{1}{M} \sum_{i=0}^{M-1} \delta(s_t - \bar{s}_i^t) \quad \text{where } \bar{s}_i^t \leftarrow P(s_t | o_{0:t-1})$$

- Update the state distribution at t

$$P(s_t | o_{0:t}) = C \sum_{i=0}^{M-1} P_\gamma(o_t - g(\bar{s}_i^t)) \delta(s_t - \bar{s}_i^t)$$

$$C = \frac{1}{\sum_{i=0}^{M-1} P_\gamma(o_t - g(\bar{s}_i^t))}$$

# Estimating a state

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

$$P(s_t | o_{0:t}) = C \sum_{i=0}^{M-1} P_\gamma(o_t - g(\bar{s}_i^t)) \delta(s_t - \bar{s}_i^t)$$

- The actual state can be estimated as the mean of this distribution

$$\hat{s}_t = C \sum_{i=0}^{M-1} \bar{s}_i^t P_\gamma(o_t - g(\bar{s}_i^t))$$

- Alternately, it can be the most likely sample

$$\hat{s}_t = \bar{s}_j^t : j = \arg \max_i P_\gamma(o_t - g(\bar{s}_i^t))$$



# Simulations with a Linear Model

$$S_t = S_{t-1} + \varepsilon_t$$

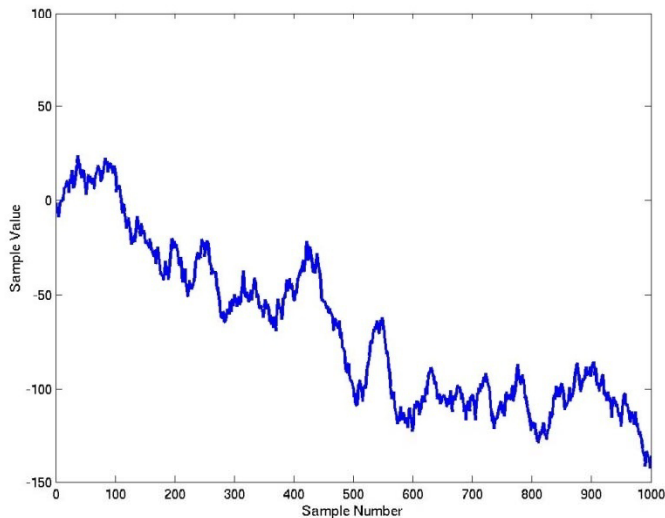
$$O_t = S_t + X_t$$

- $\varepsilon_t$  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  $O_t$

# Simulation: Synthesizing data

Generate state sequence according to:  
 $\varepsilon_t$  is Gaussian with mean 0 and variance 10

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



# Simulation: Synthesizing data

Generate state sequence according to:  
 $\varepsilon_t$  is Gaussian with mean 0 and variance 10

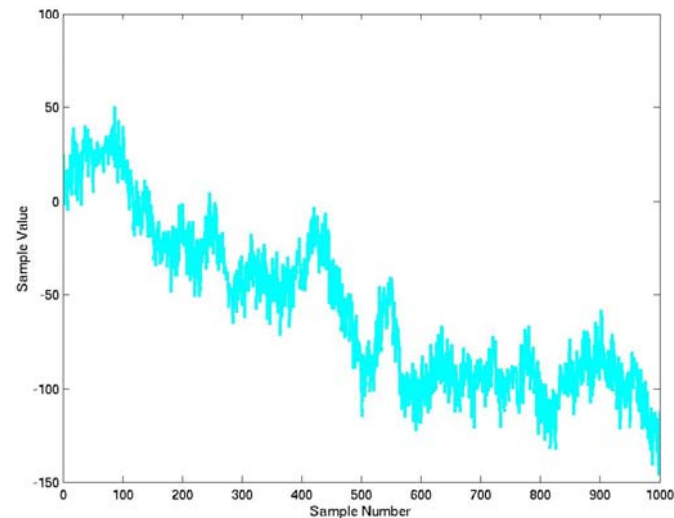
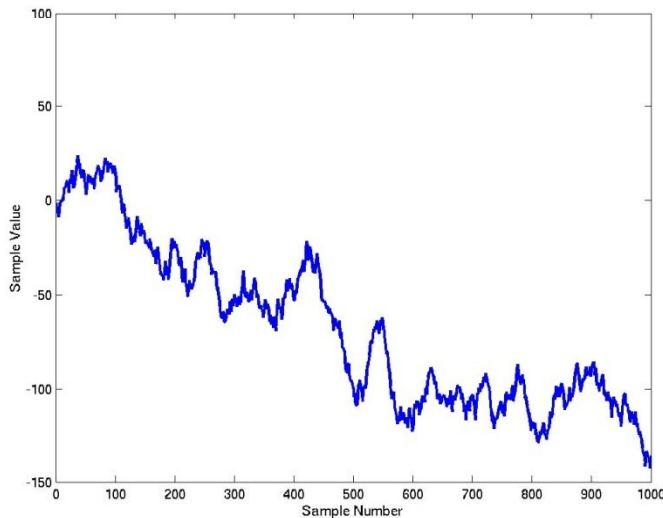
$$s_t = s_{t-1} + \varepsilon_t$$

Generate observation sequence from state sequence according to:  $o_t = s_t + x_t$   
 $x_t$  is mixture Gaussian with parameters:

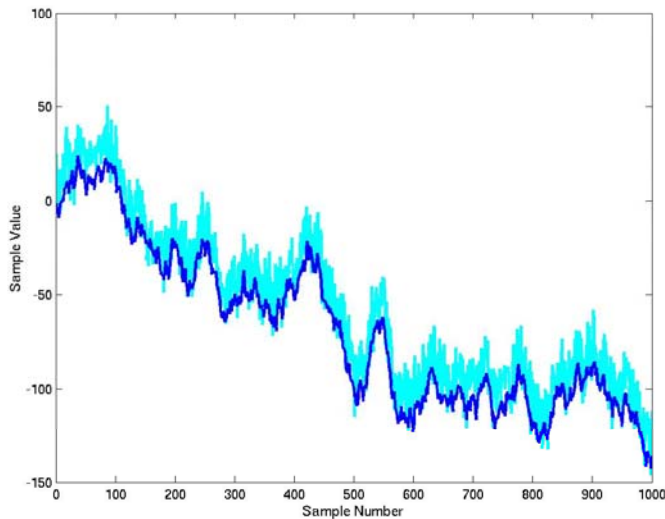
Means = [-4, 0, 4, 8, 12, 16, 18, 20]

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]

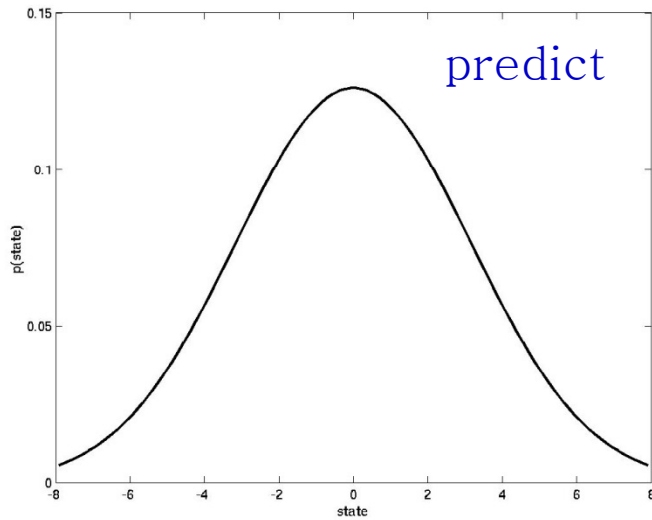


# Simulation: Synthesizing data

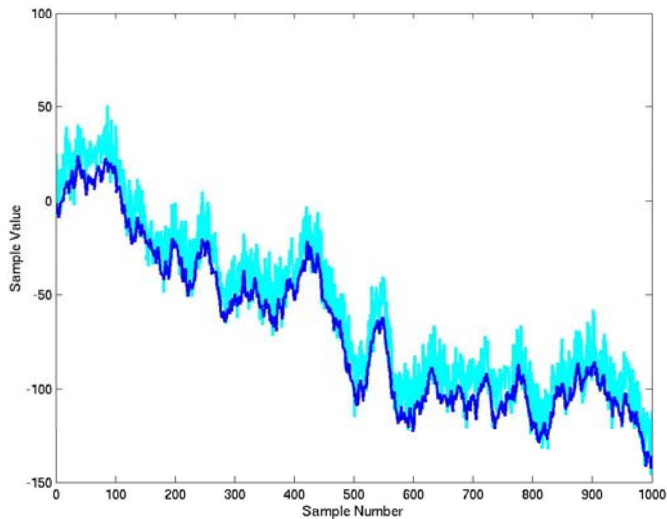


Combined figure for more compact representation

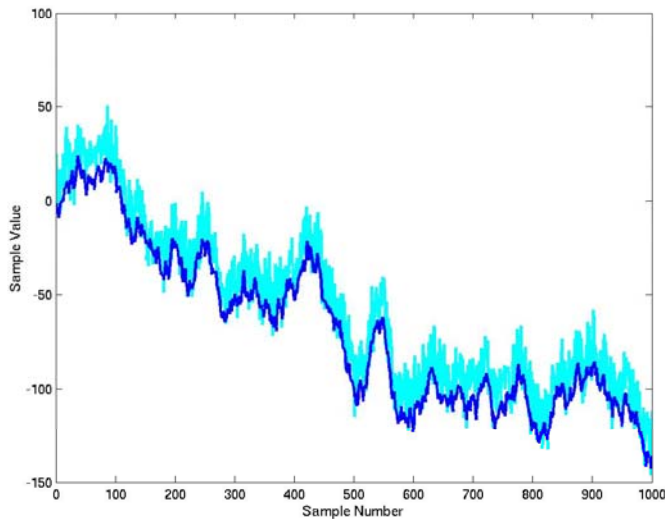
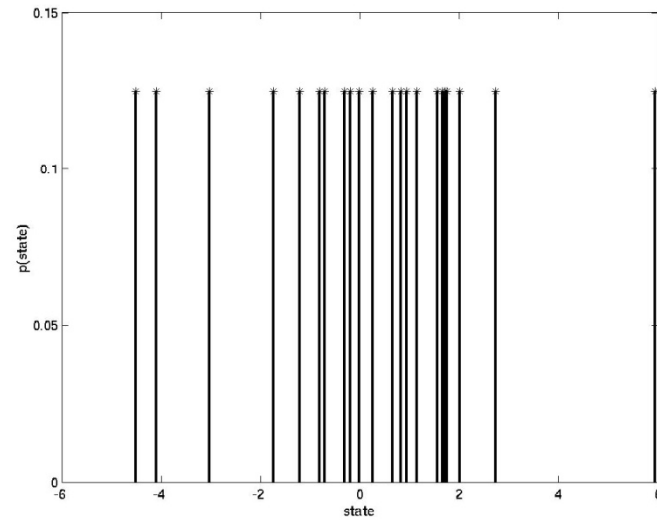
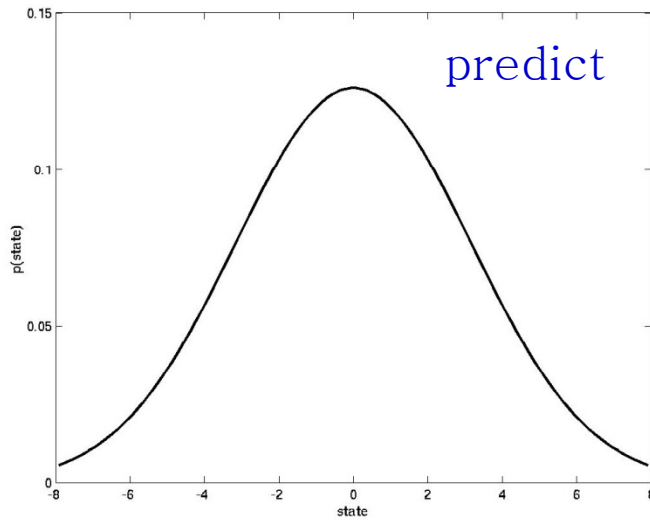
# SIMULATION: TIME = 1



PREDICTED STATE DISTRIBUTION  
AT TIME = 1

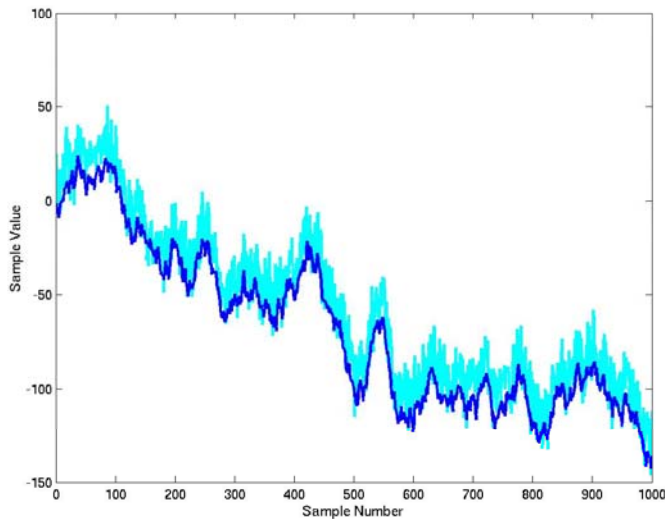
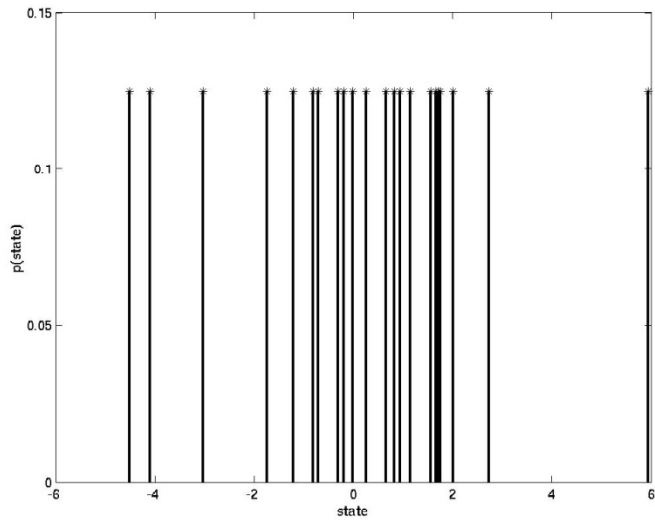


# SIMULATION: TIME = 1



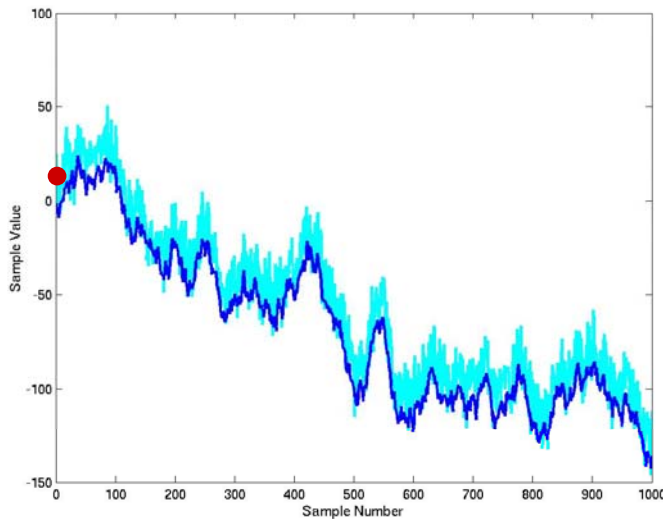
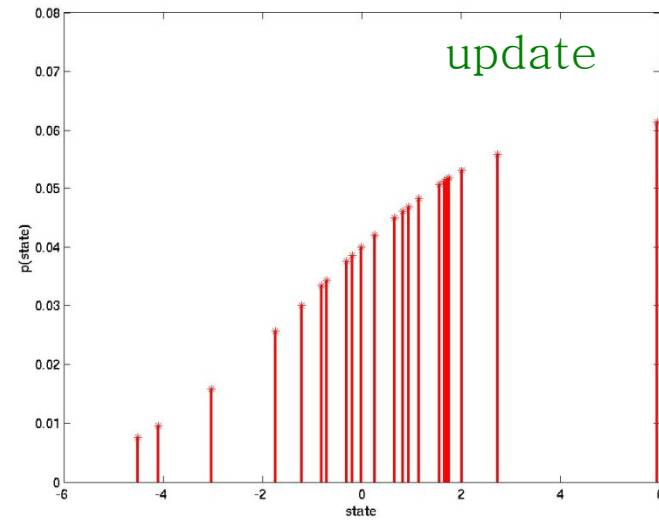
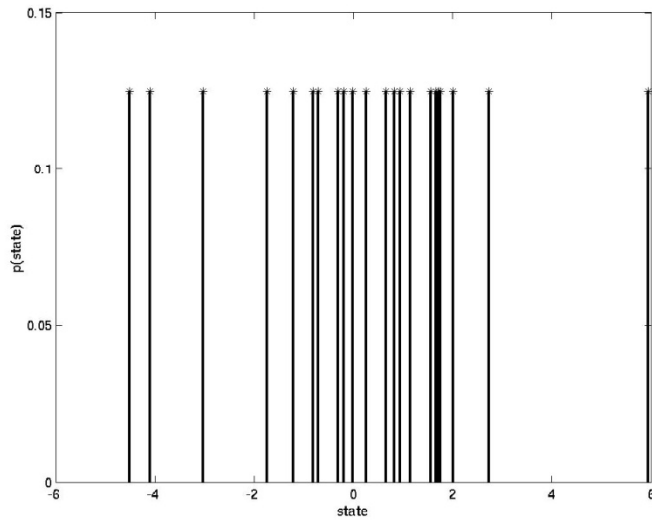
SAMPLED VERSION OF  
PREDICTED STATE DISTRIBUTION  
AT TIME = 1

# SIMULATION: TIME = 1



SAMPLED VERSION OF  
PREDICTED STATE DISTRIBUTION  
AT TIME = 1

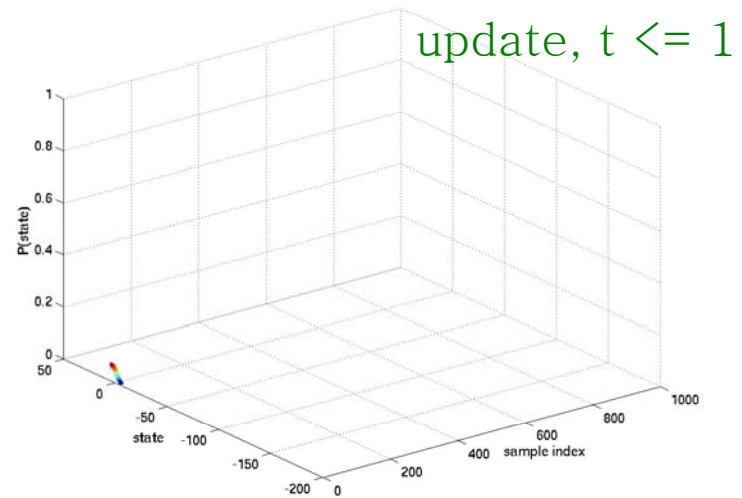
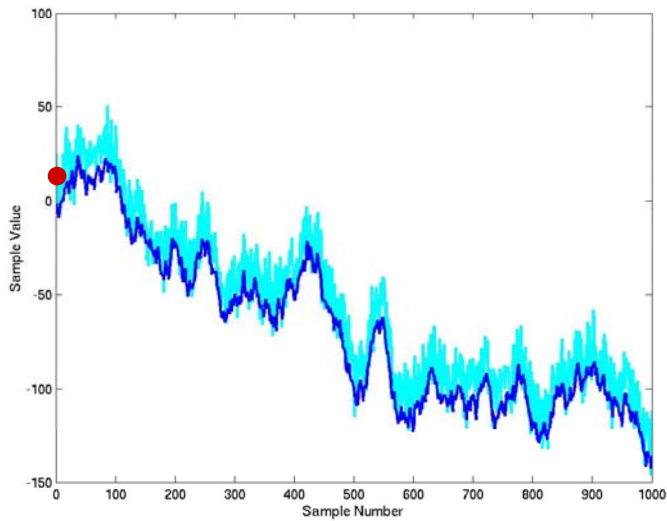
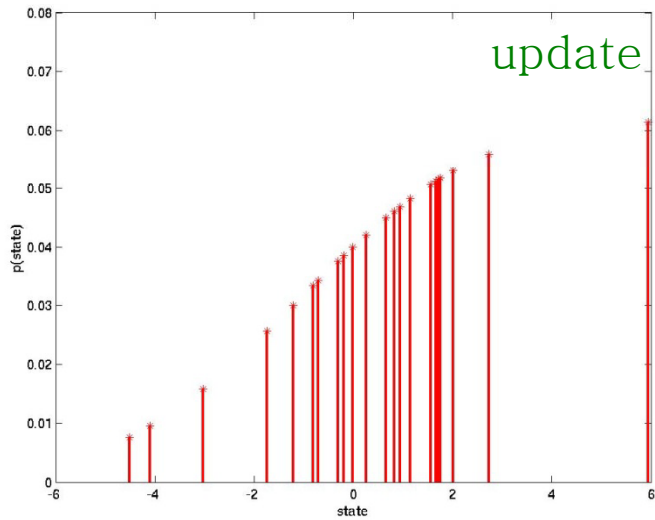
# SIMULATION: TIME = 1



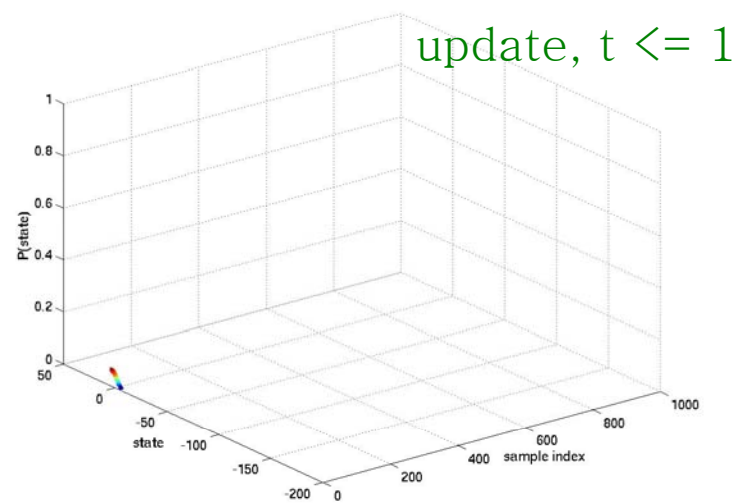
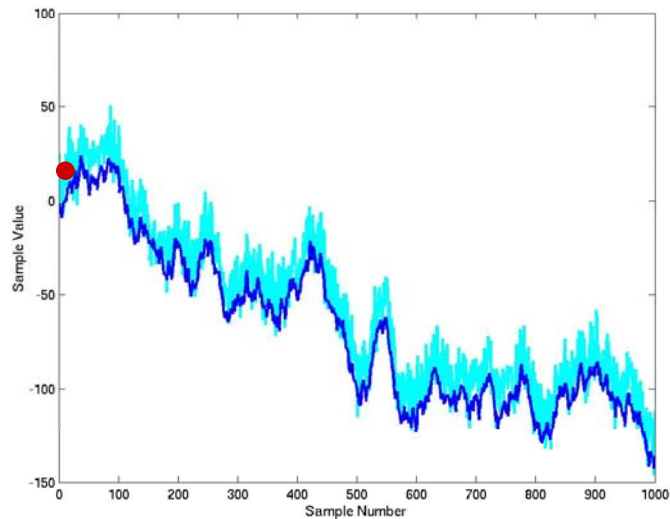
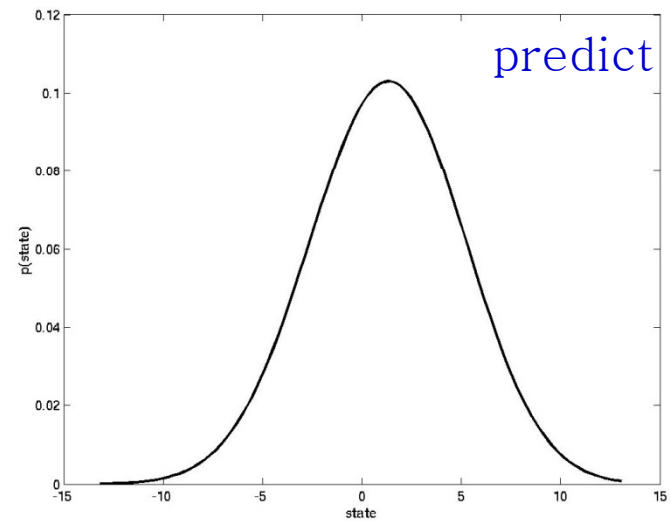
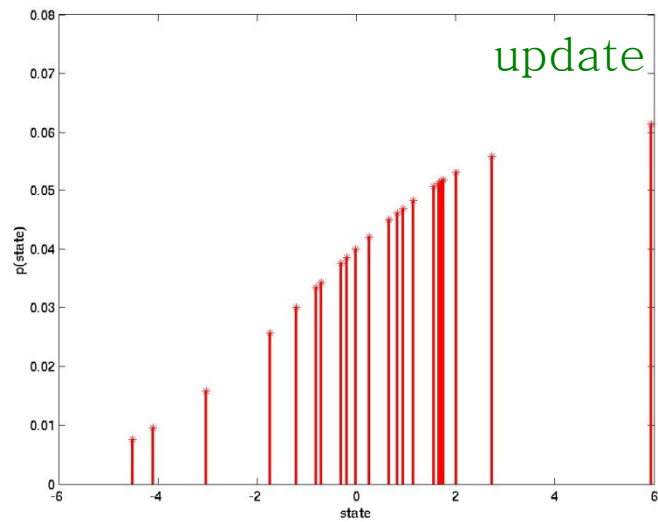
UPDATED VERSION OF  
SAMPLED VERSION OF  
PREDICTED STATE DISTRIBUTION  
AT TIME = 1  
AFTER SEEING FIRST OBSERVATION



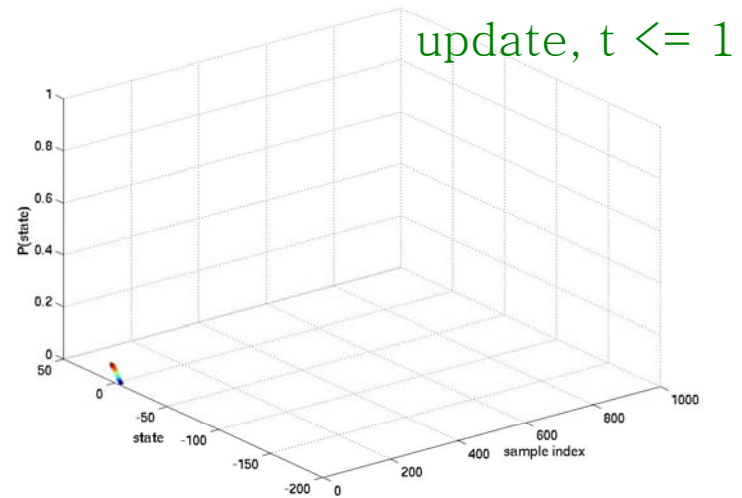
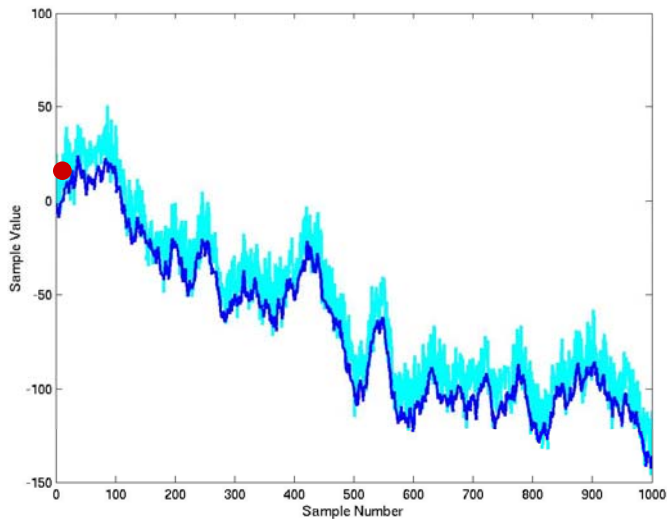
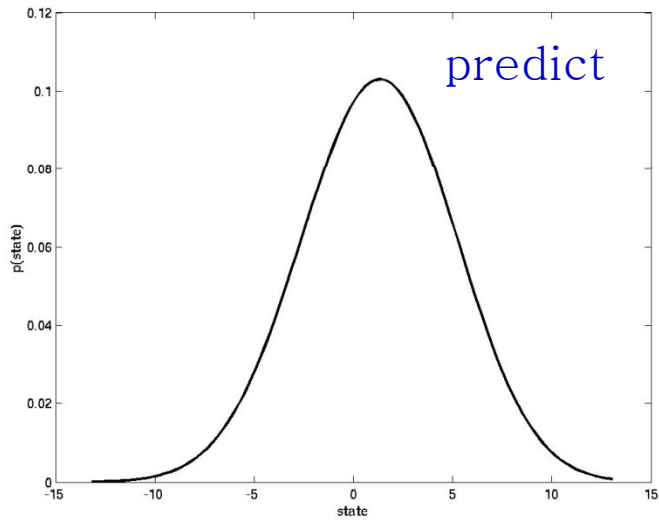
# SIMULATION: TIME = 1



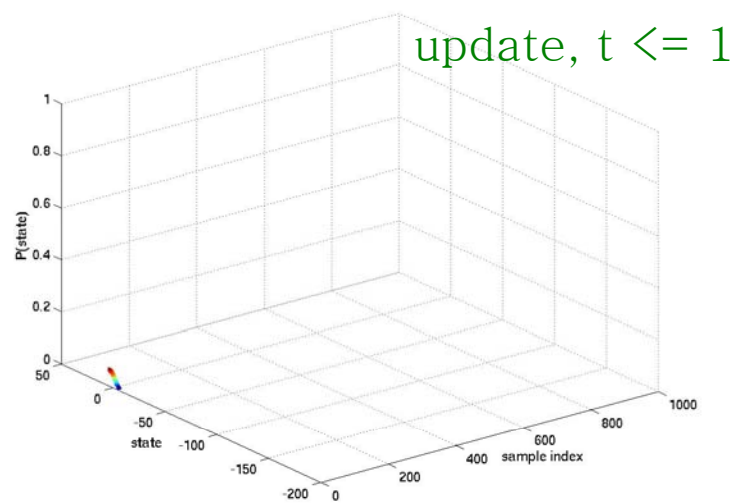
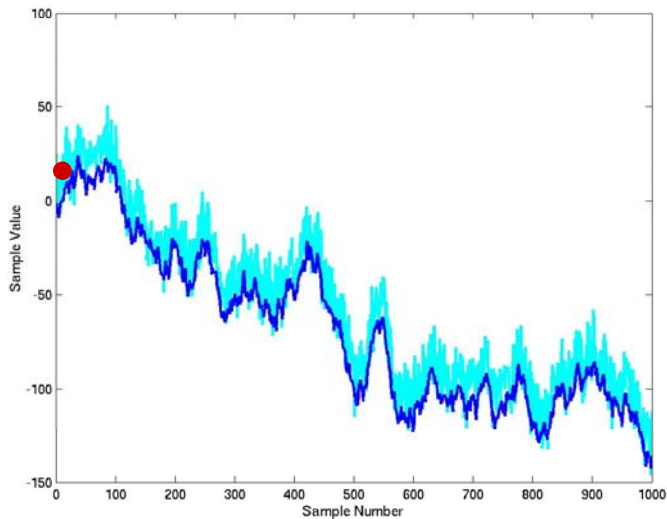
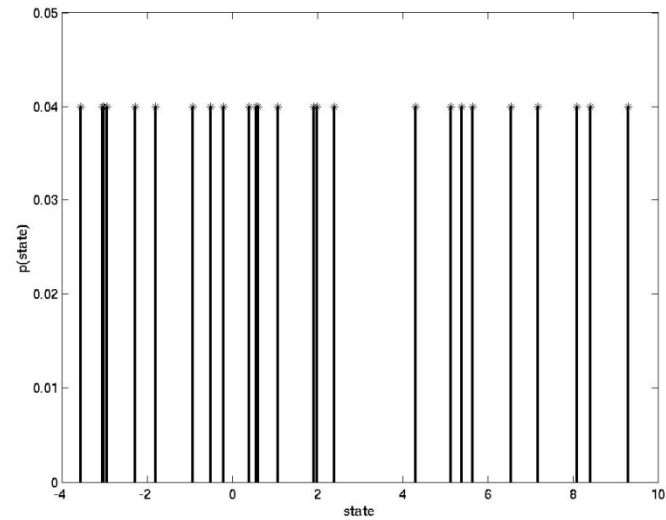
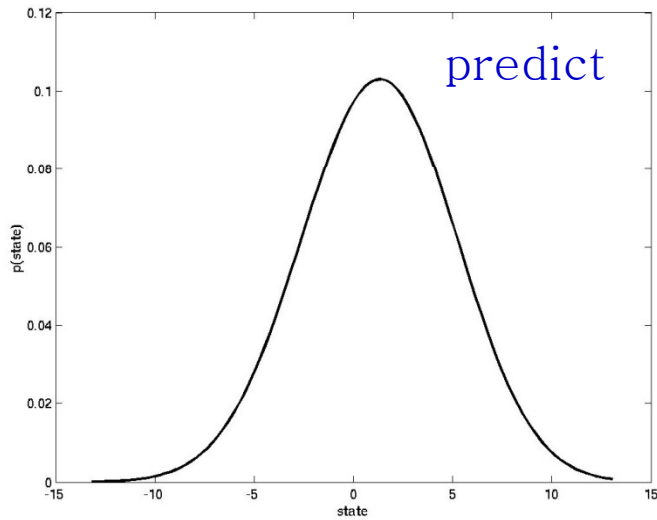
# SIMULATION: TIME = 2



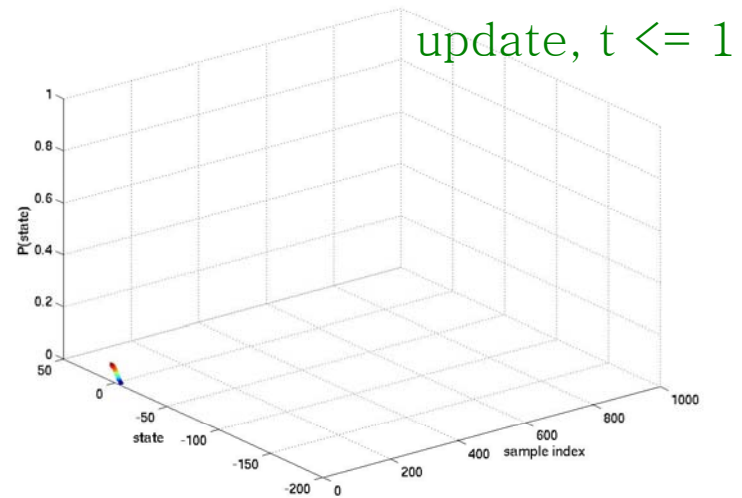
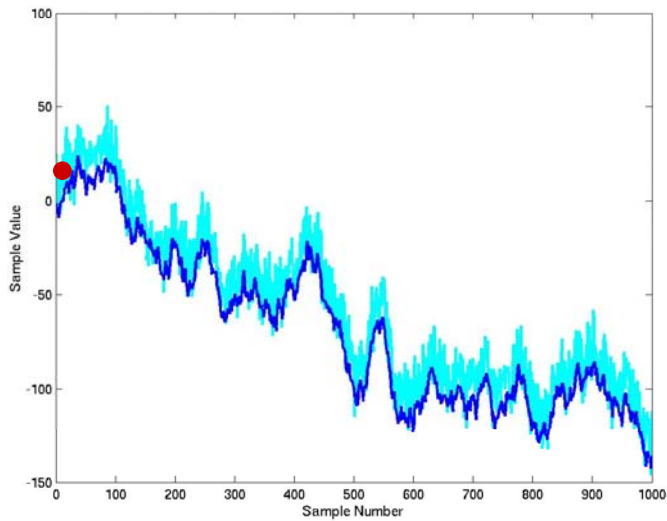
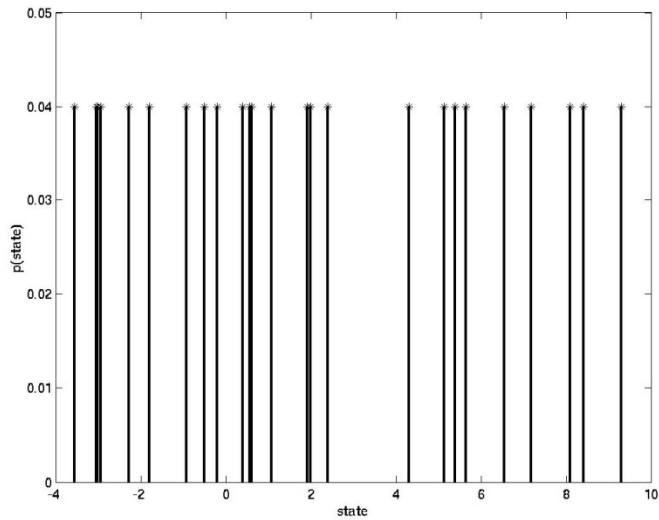
# SIMULATION: TIME = 2



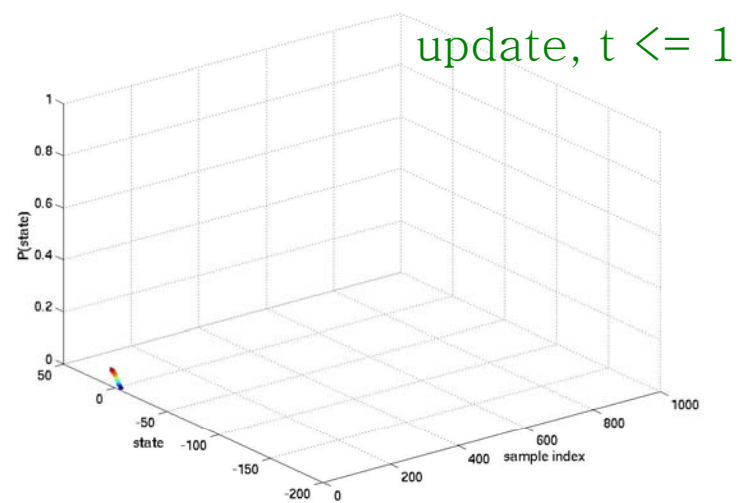
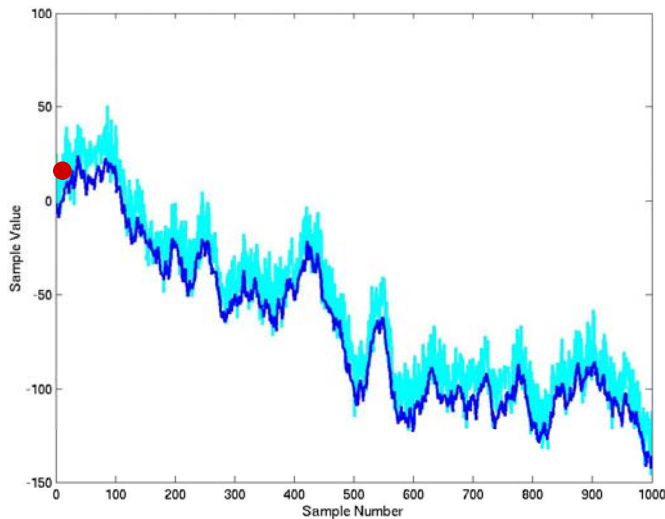
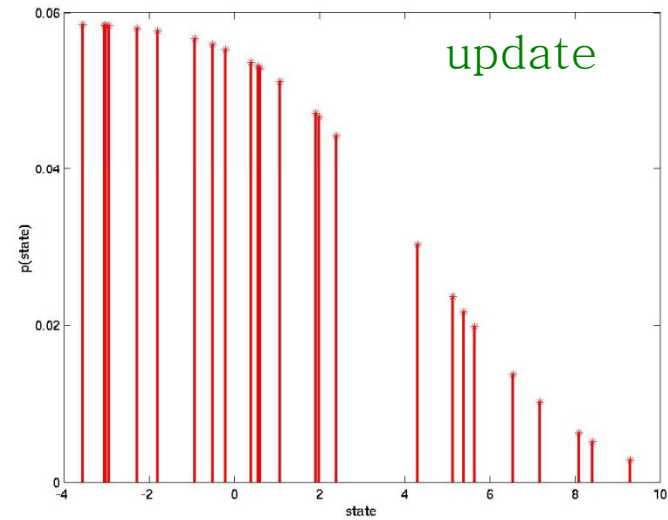
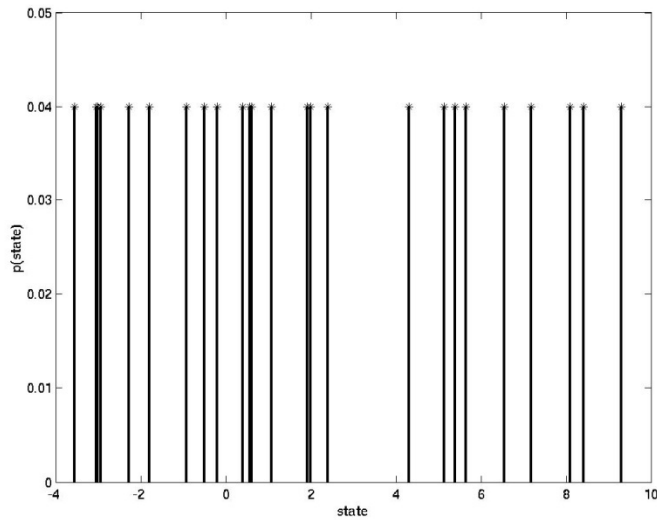
# SIMULATION: TIME = 2



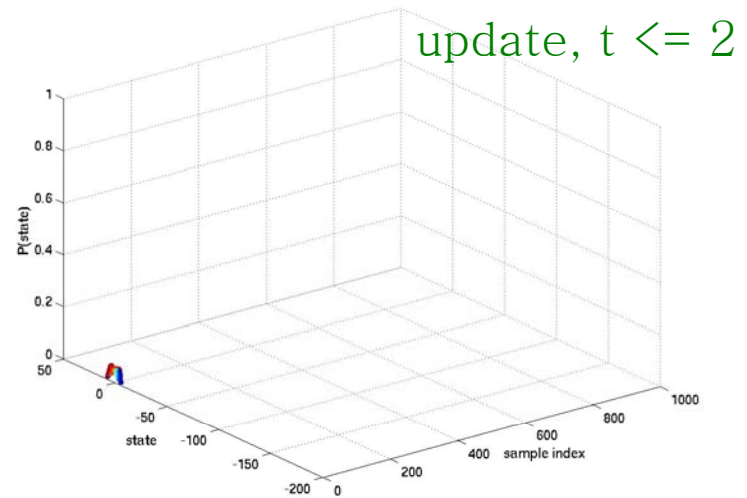
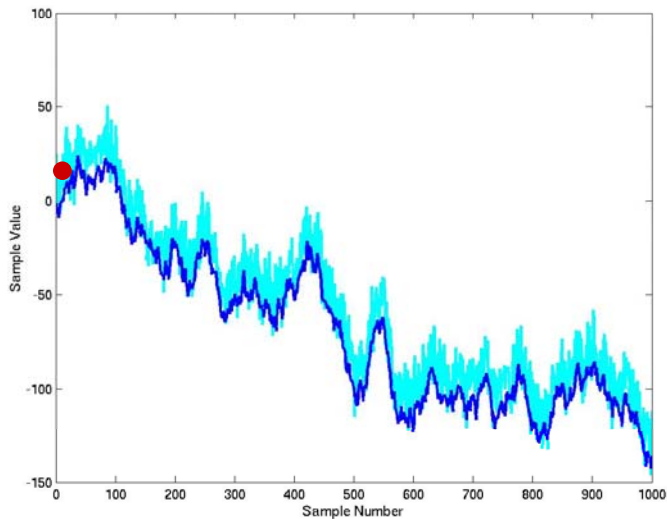
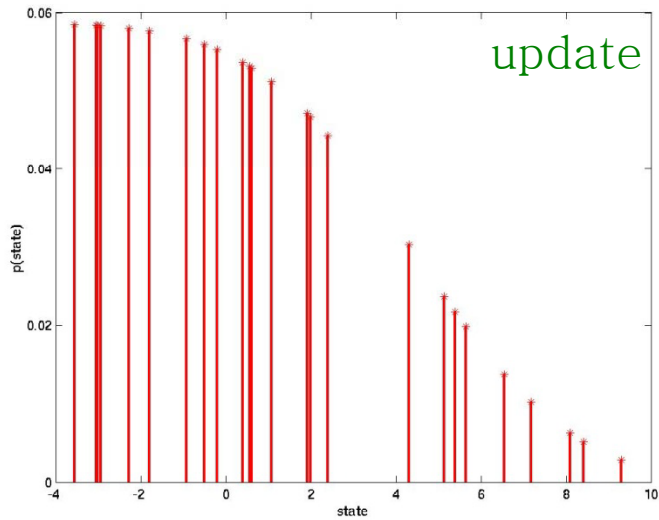
# SIMULATION: TIME = 2



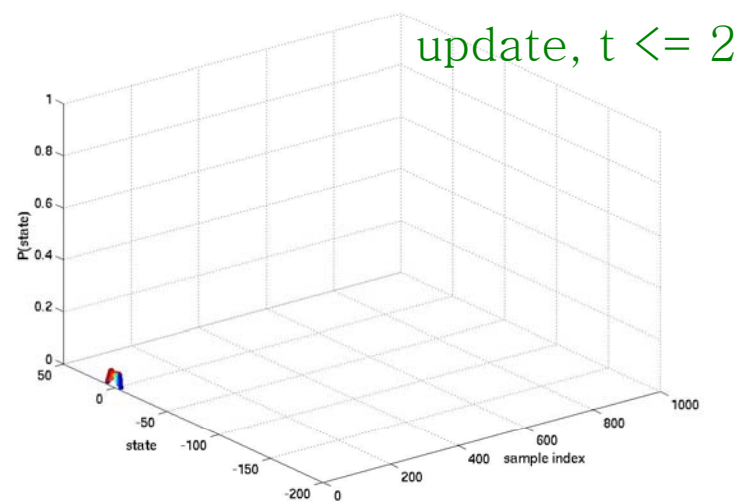
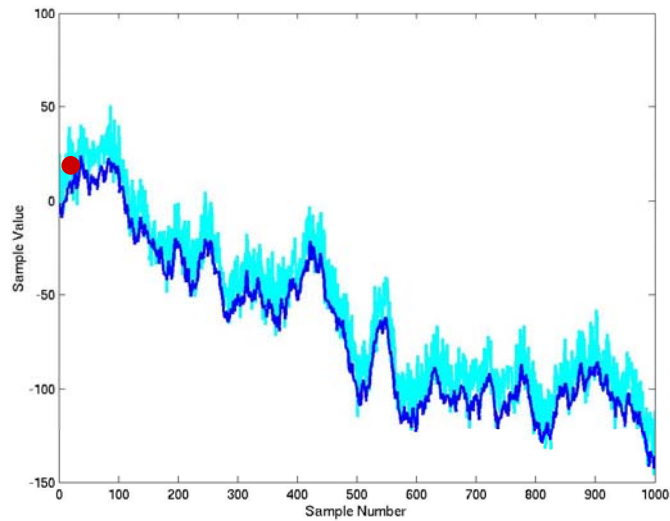
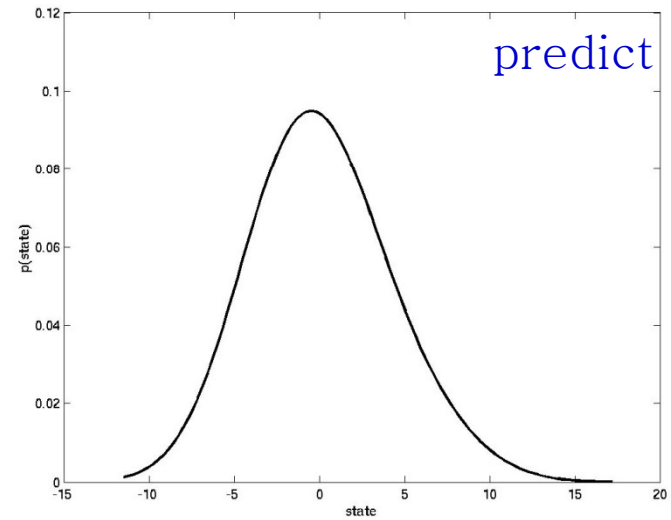
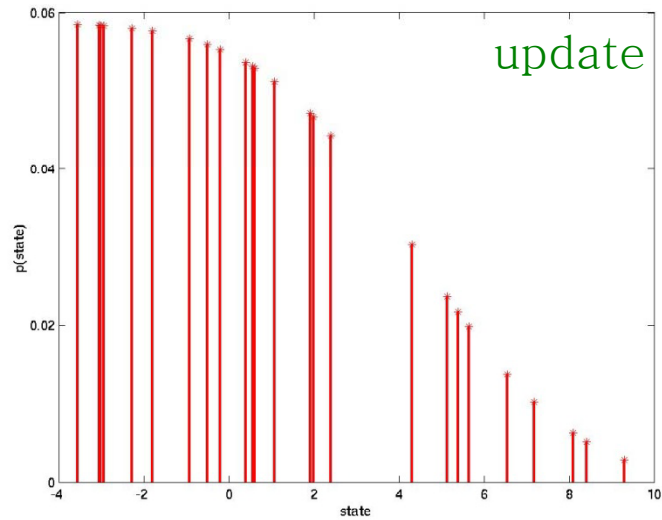
# SIMULATION: TIME = 2



# SIMULATION: TIME = 2

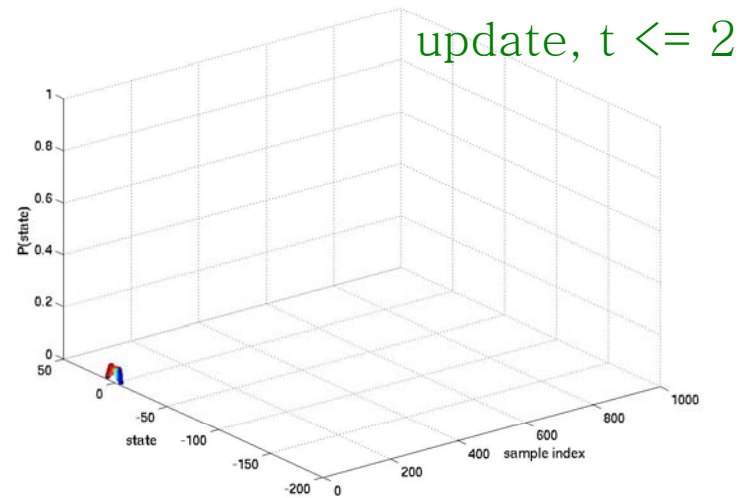
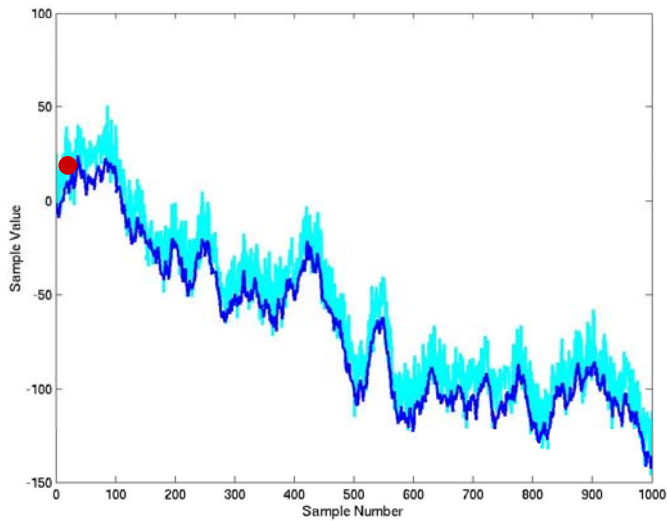
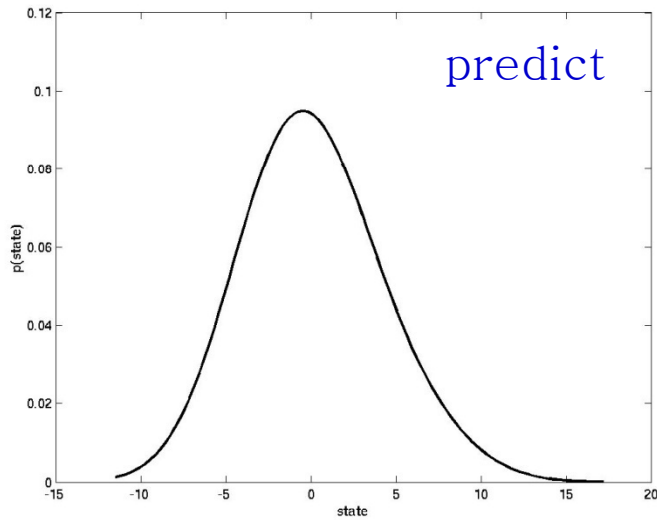


# SIMULATION: TIME = 3

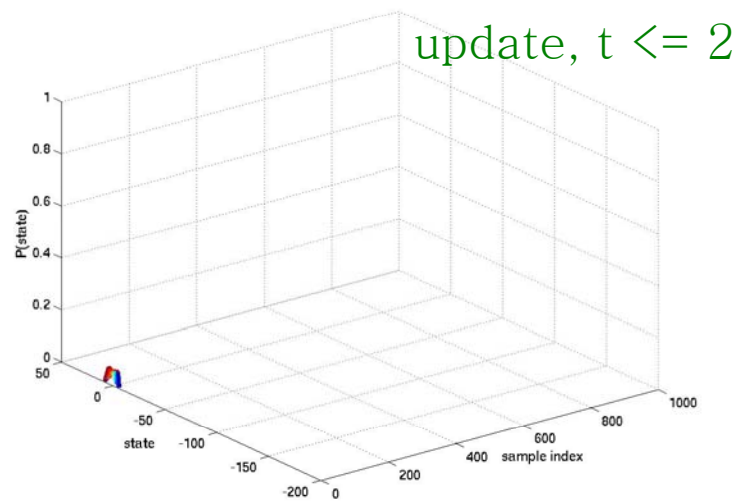
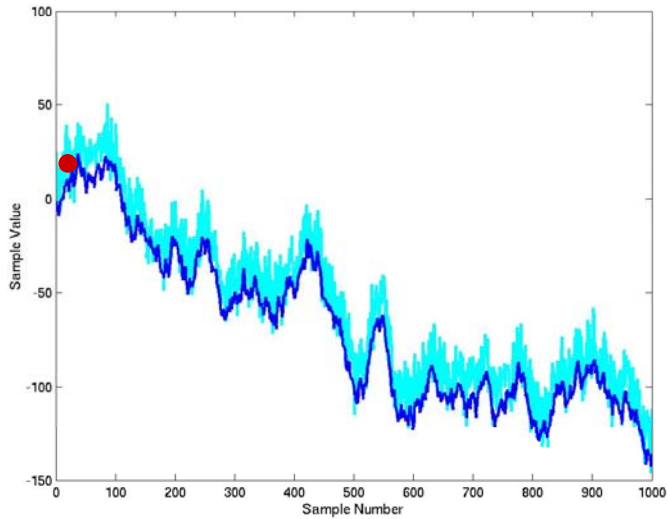
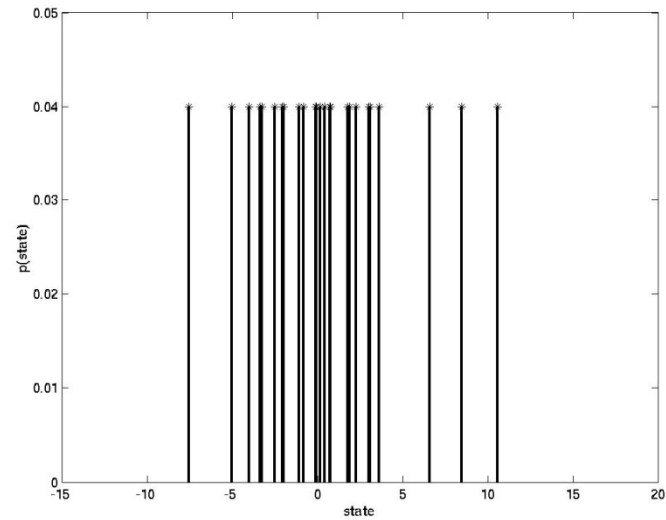
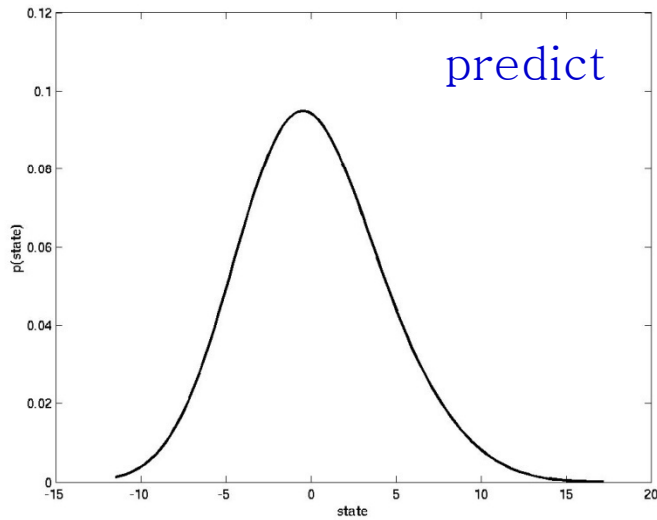




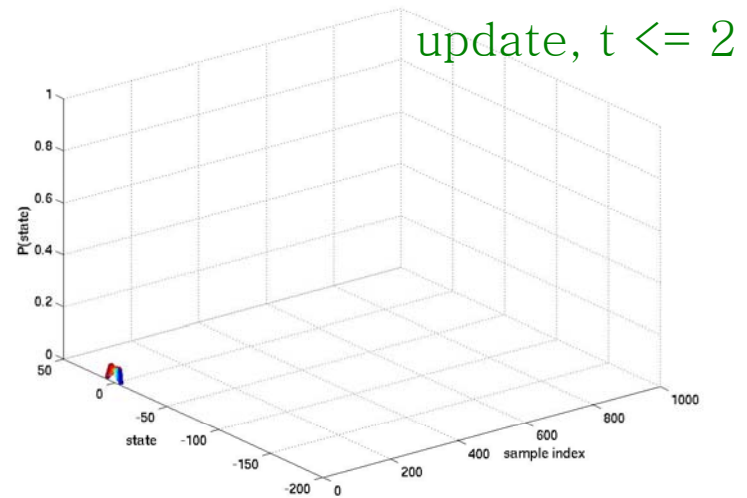
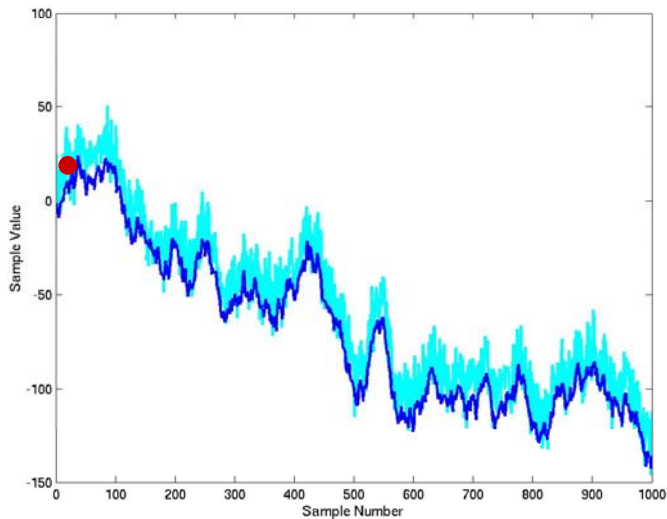
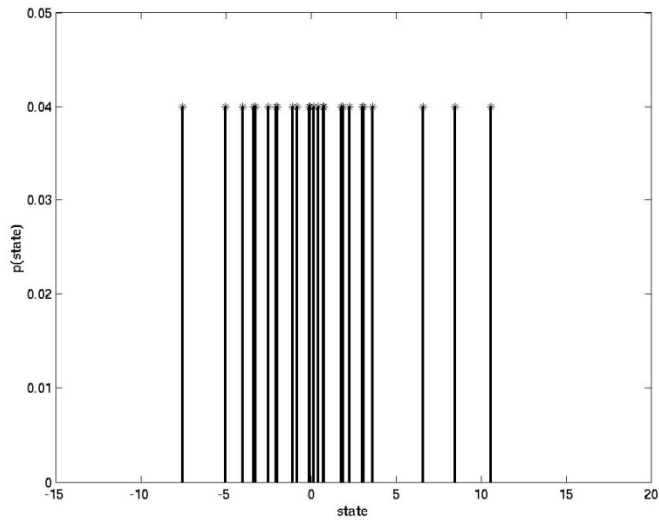
# SIMULATION: TIME = 3



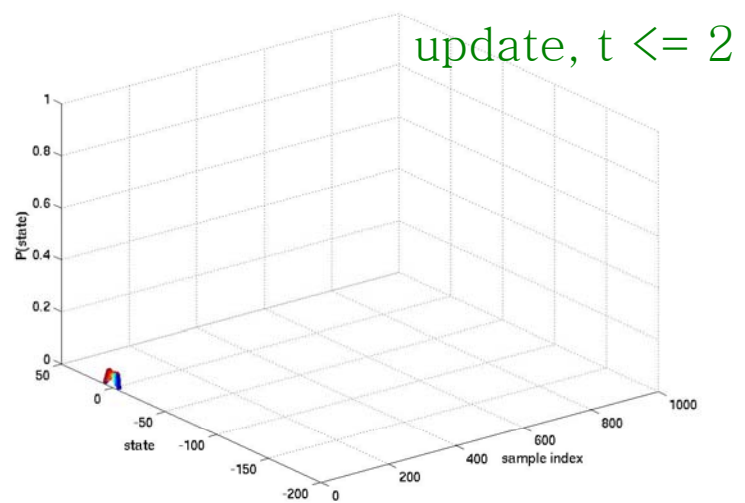
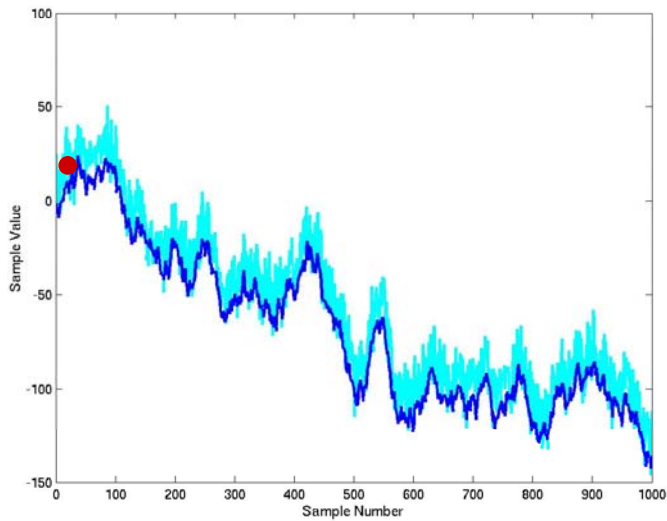
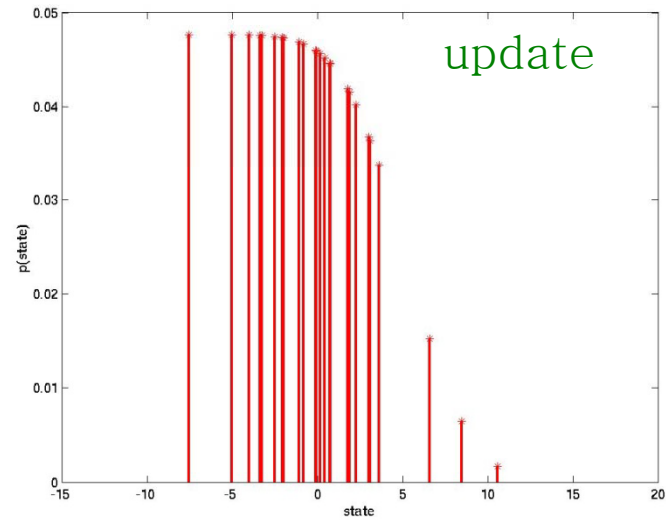
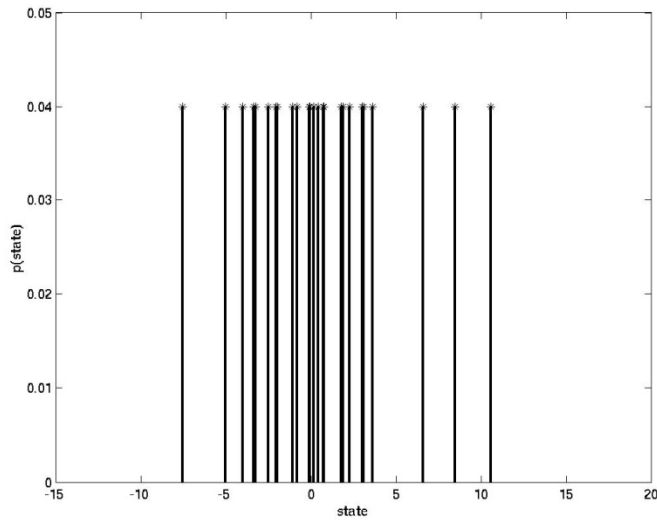
# SIMULATION: TIME = 3



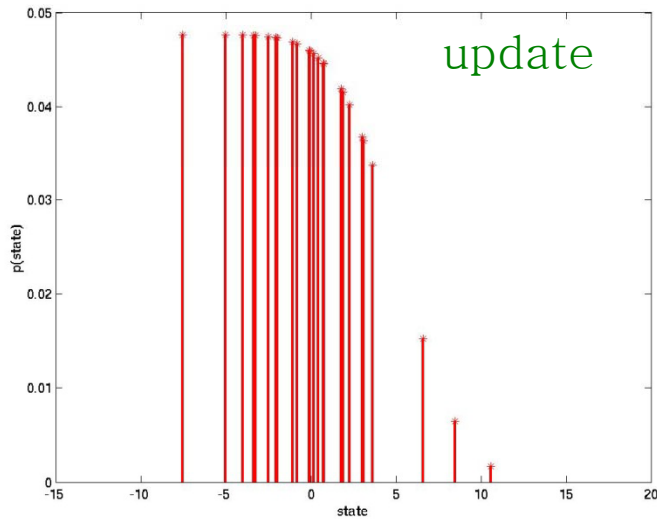
# SIMULATION: TIME = 3



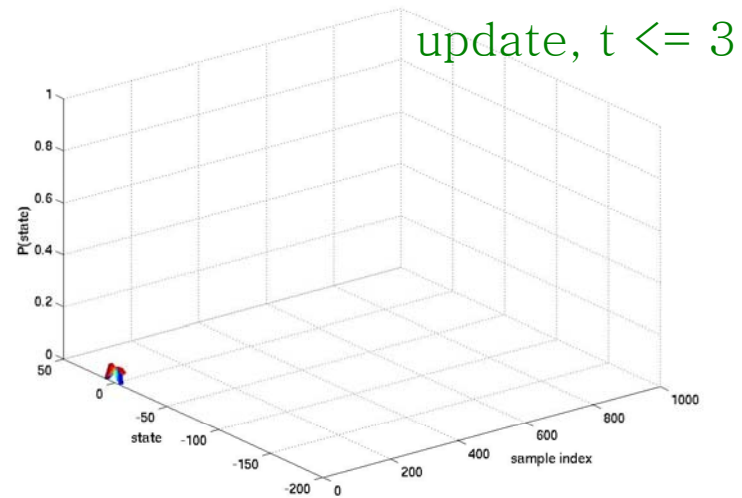
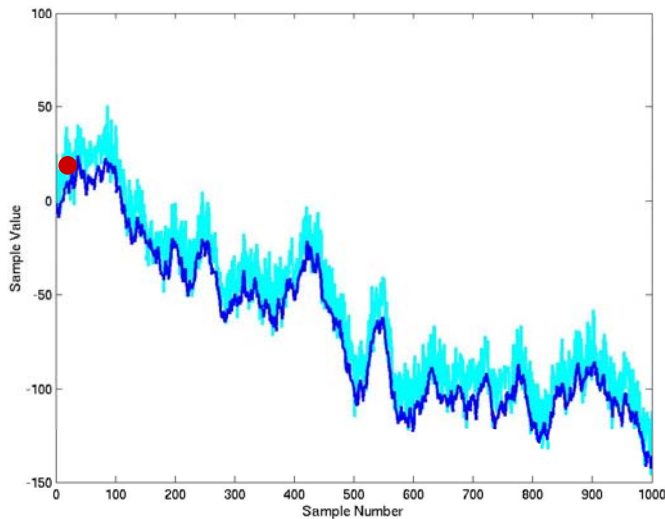
# SIMULATION: TIME = 3



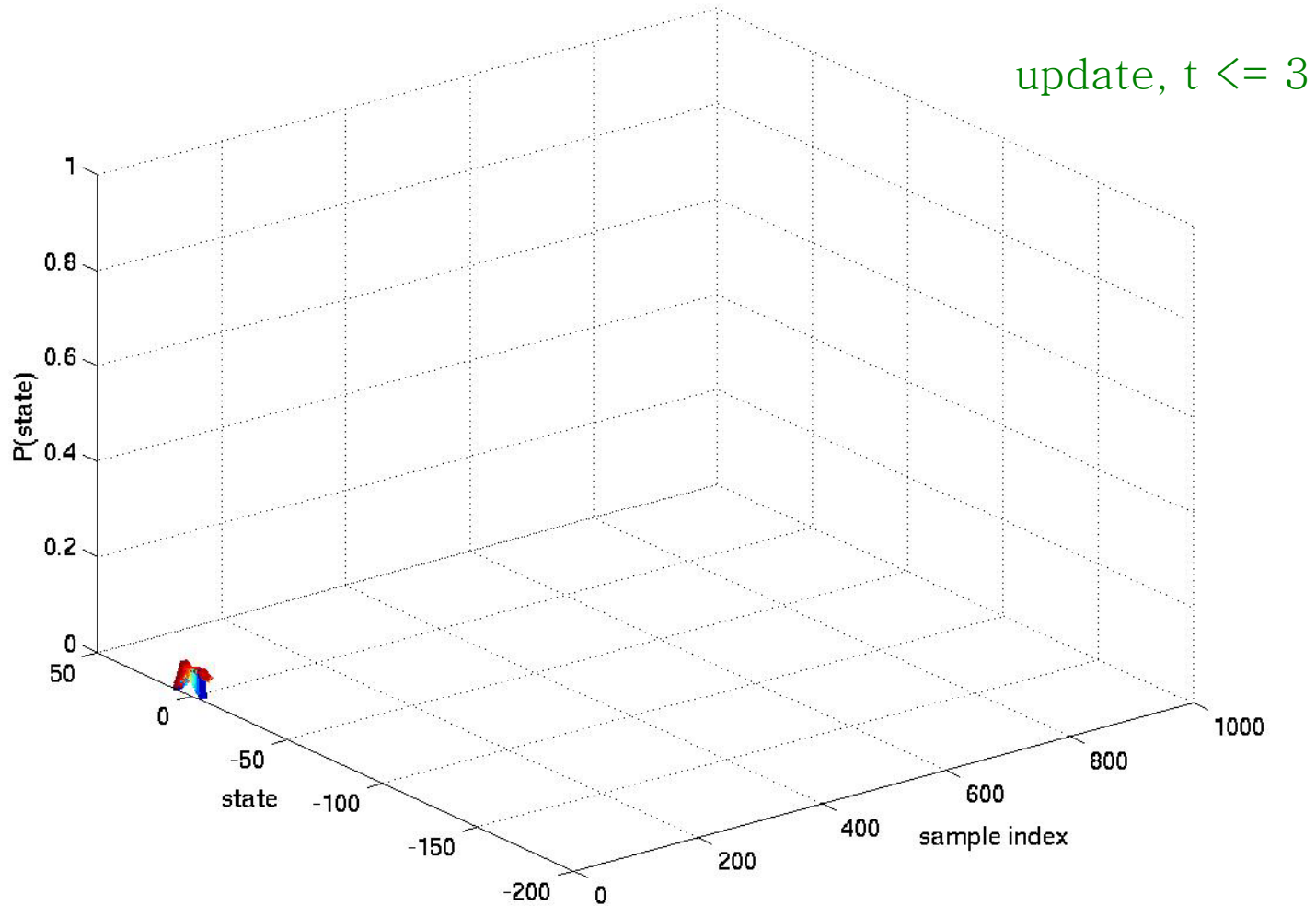
# SIMULATION: TIME = 3



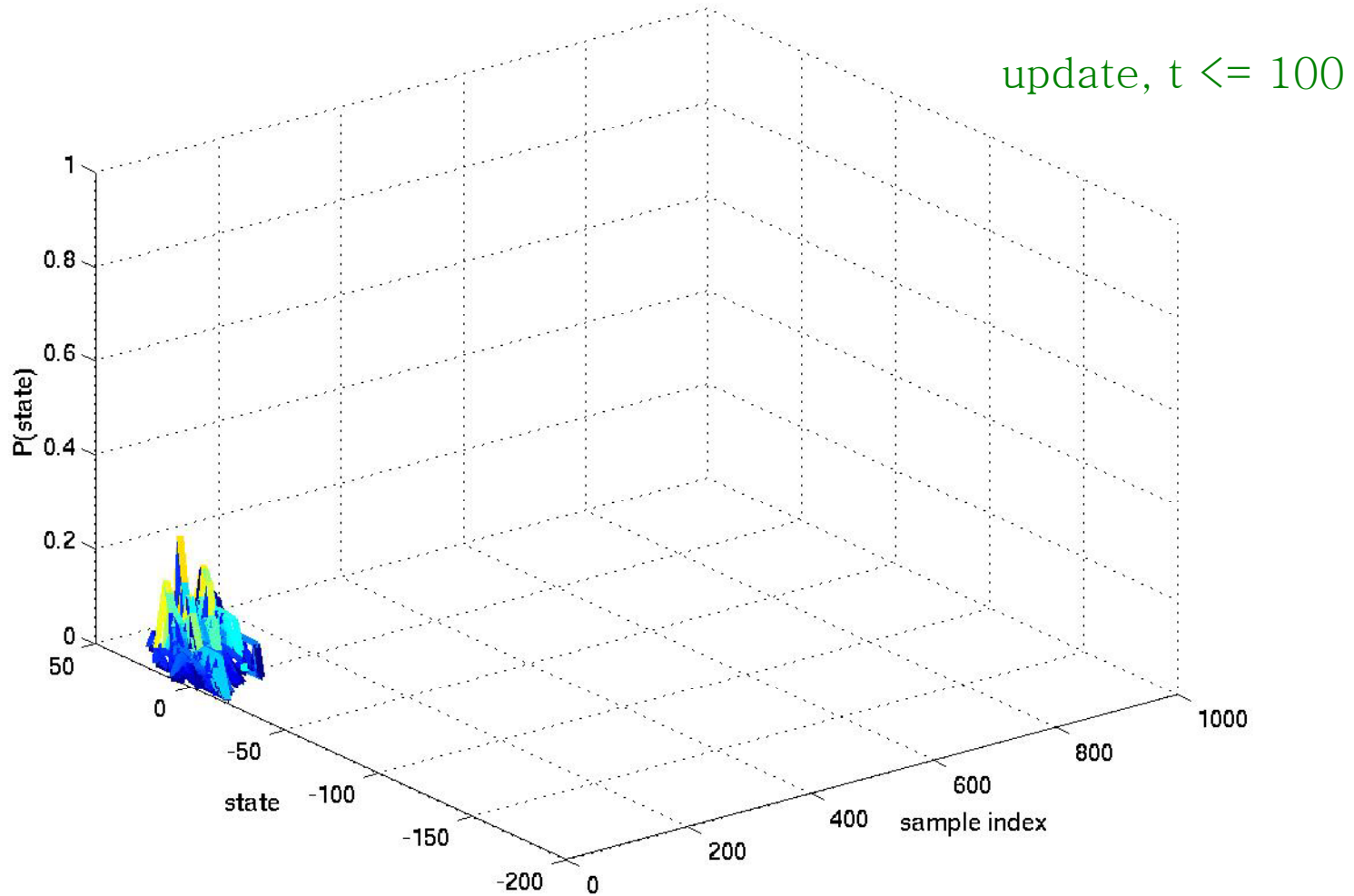
The figure below shows the contour of the updated state probabilities for all time instants until the current instant



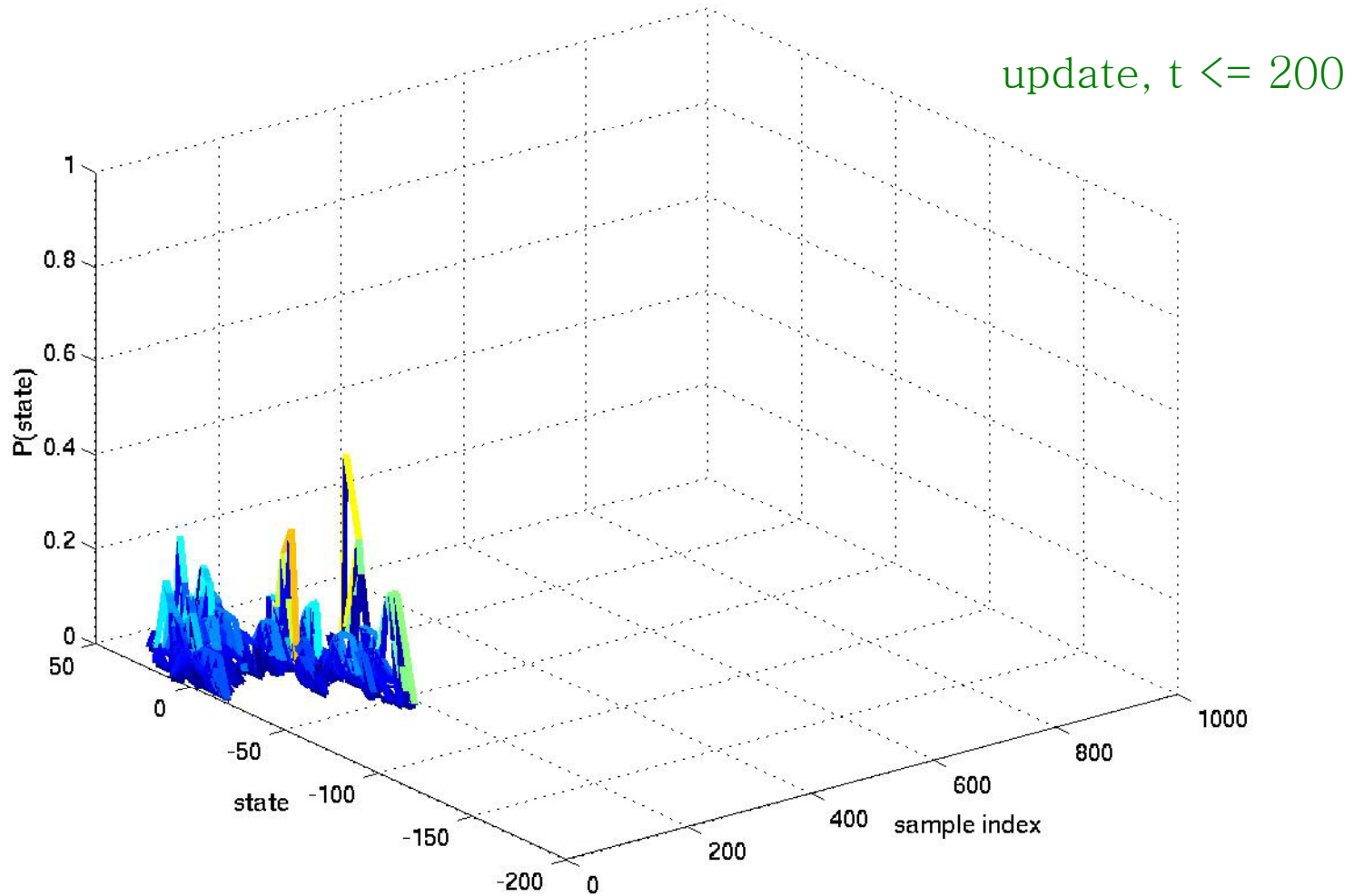
# Simulation: Updated Probs Until $T=3$



# Simulation: Updated Probs Until $T=100$

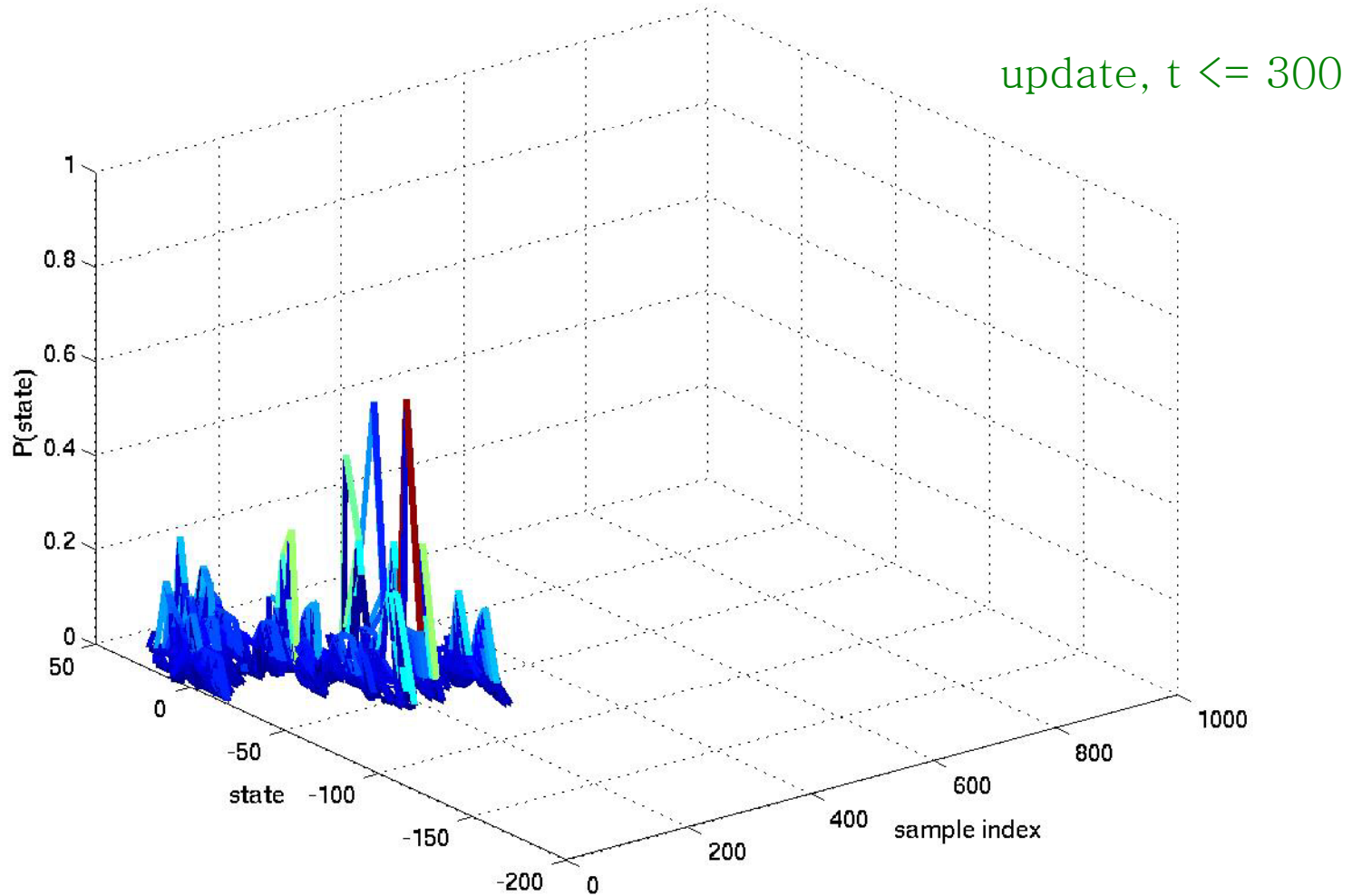


# Simulation: Updated Probs Until $T=200$

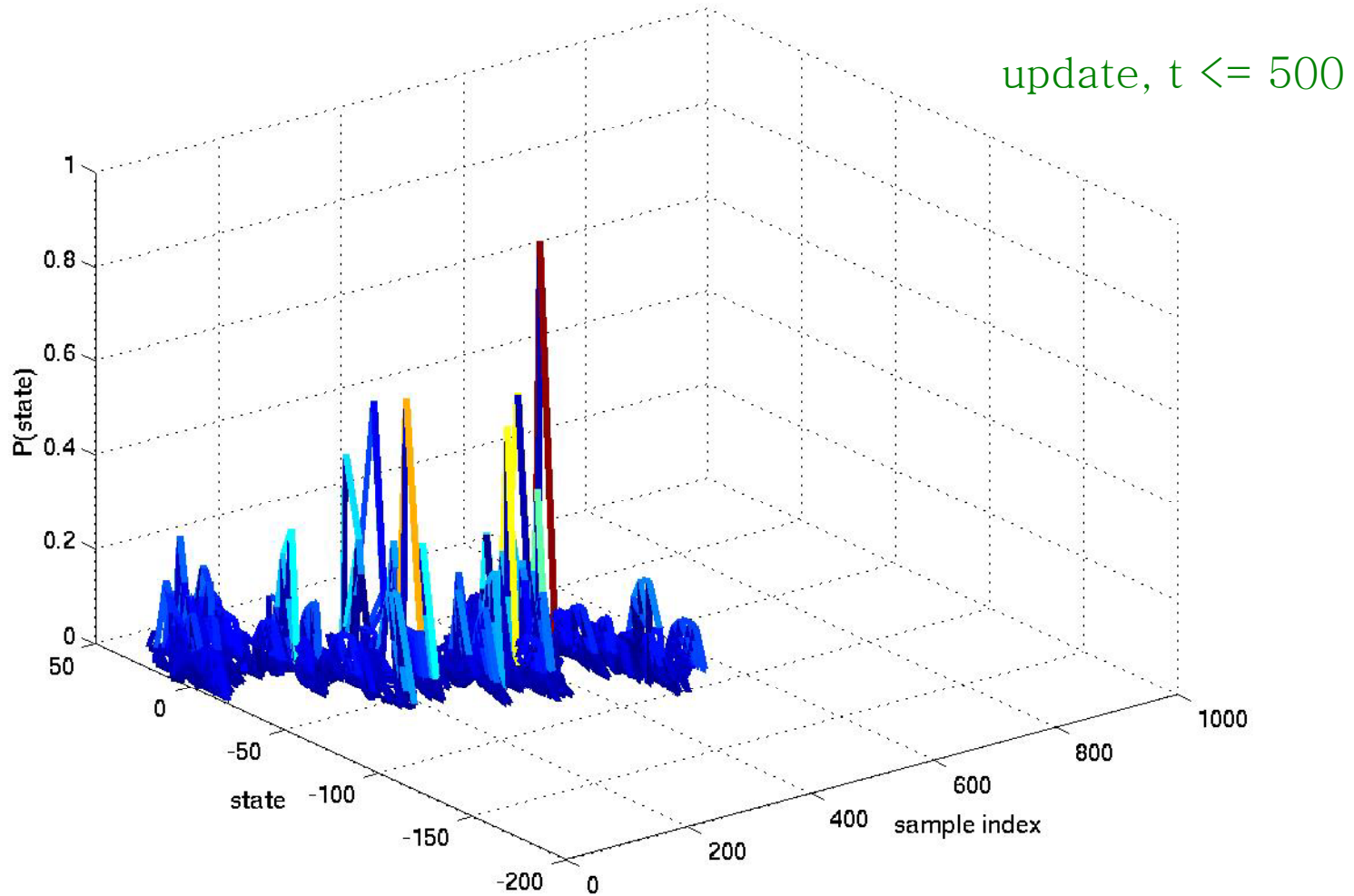




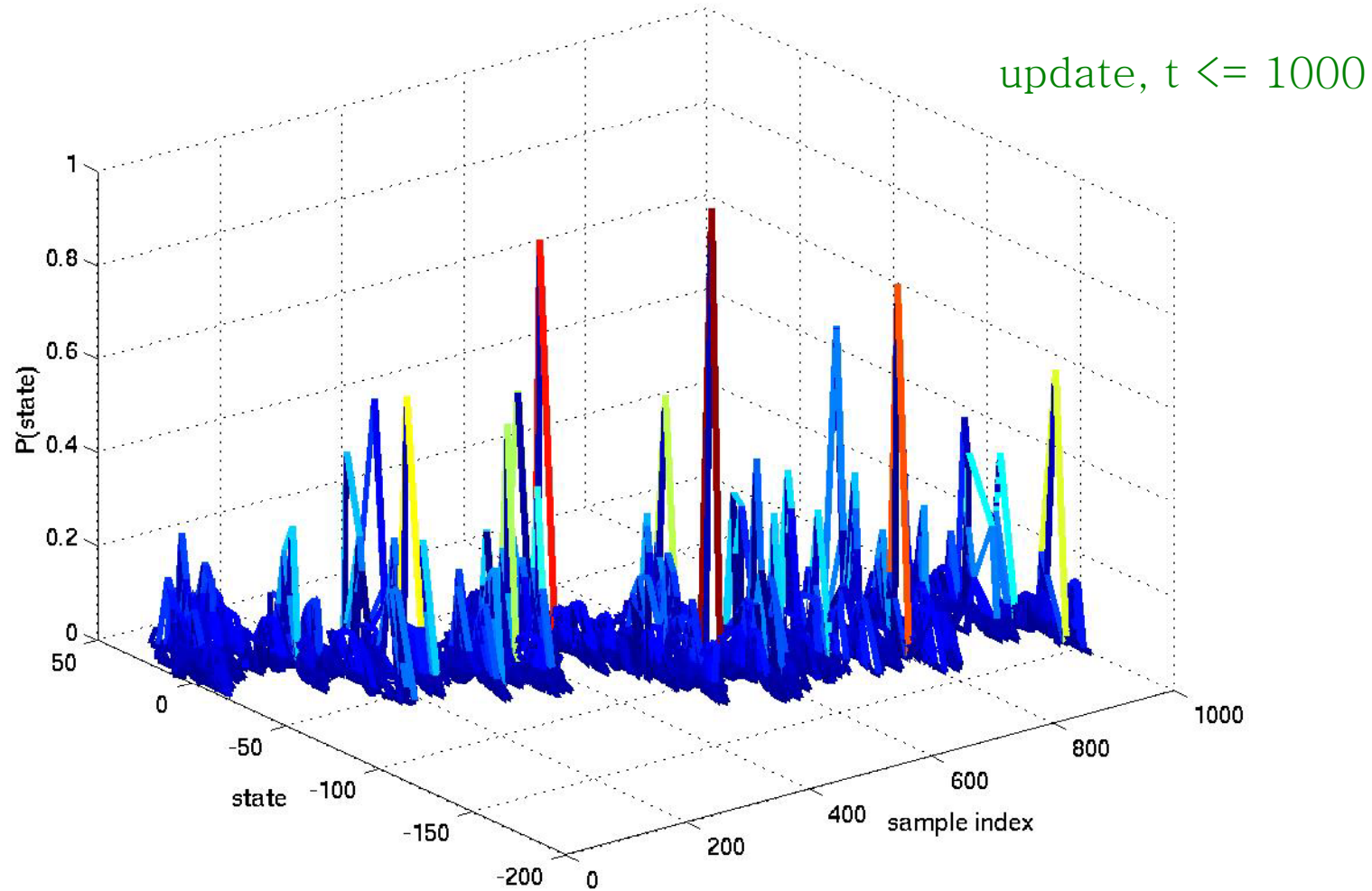
# Simulation: Updated Probs Until $T=300$



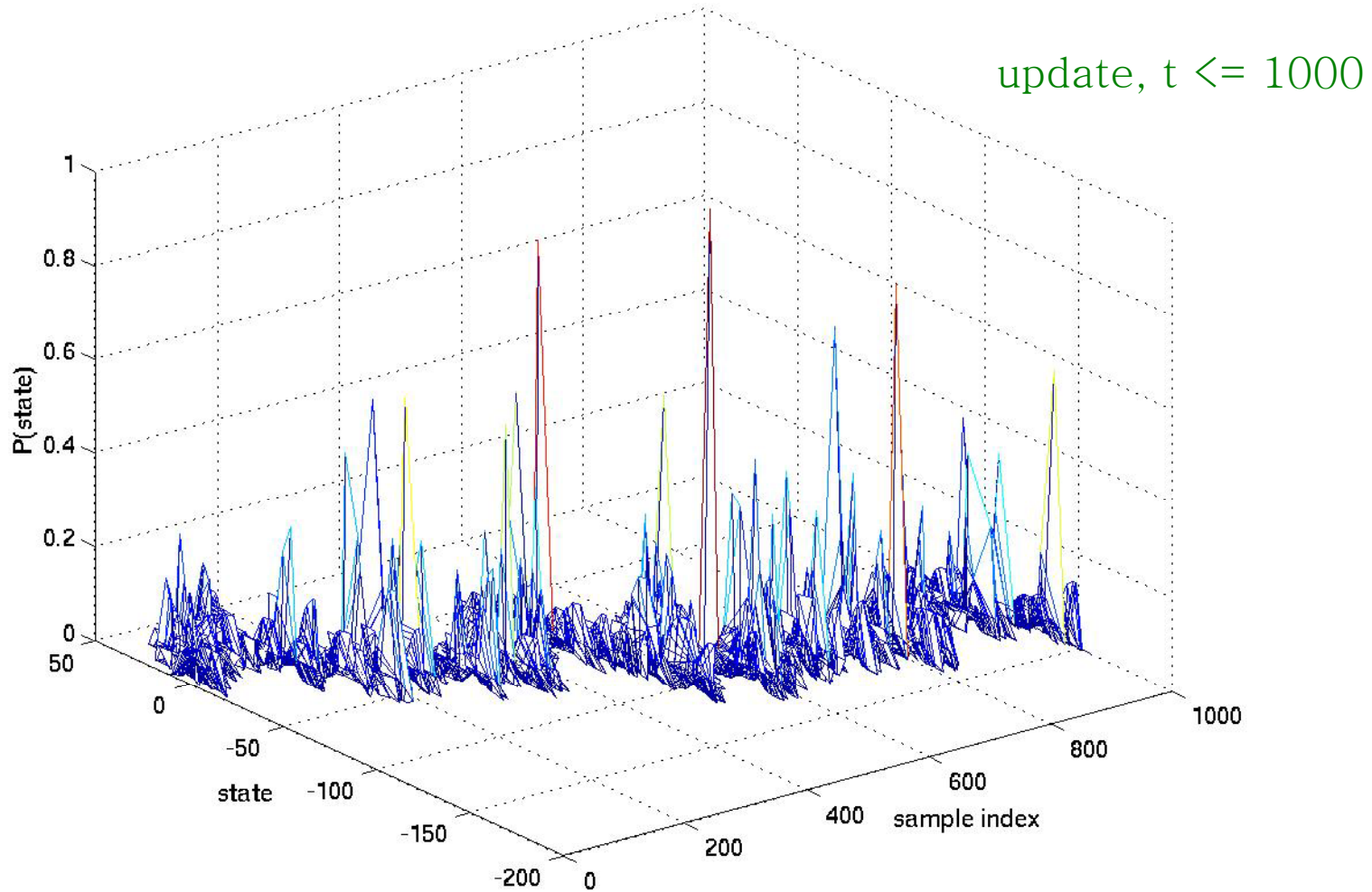
# Simulation: Updated Probs Until $T=500$



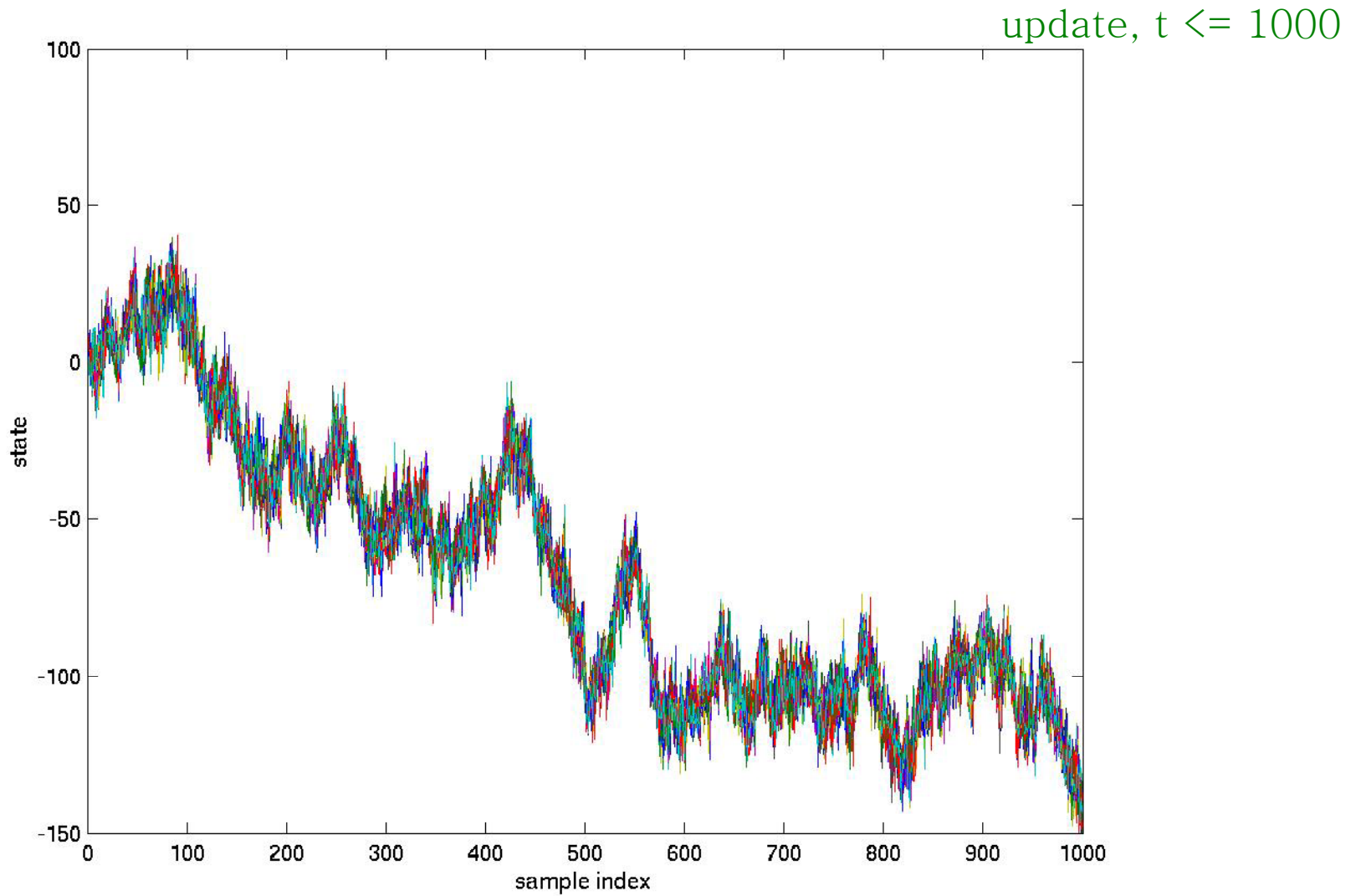
# Simulation: Updated Probs Until $T=1000$



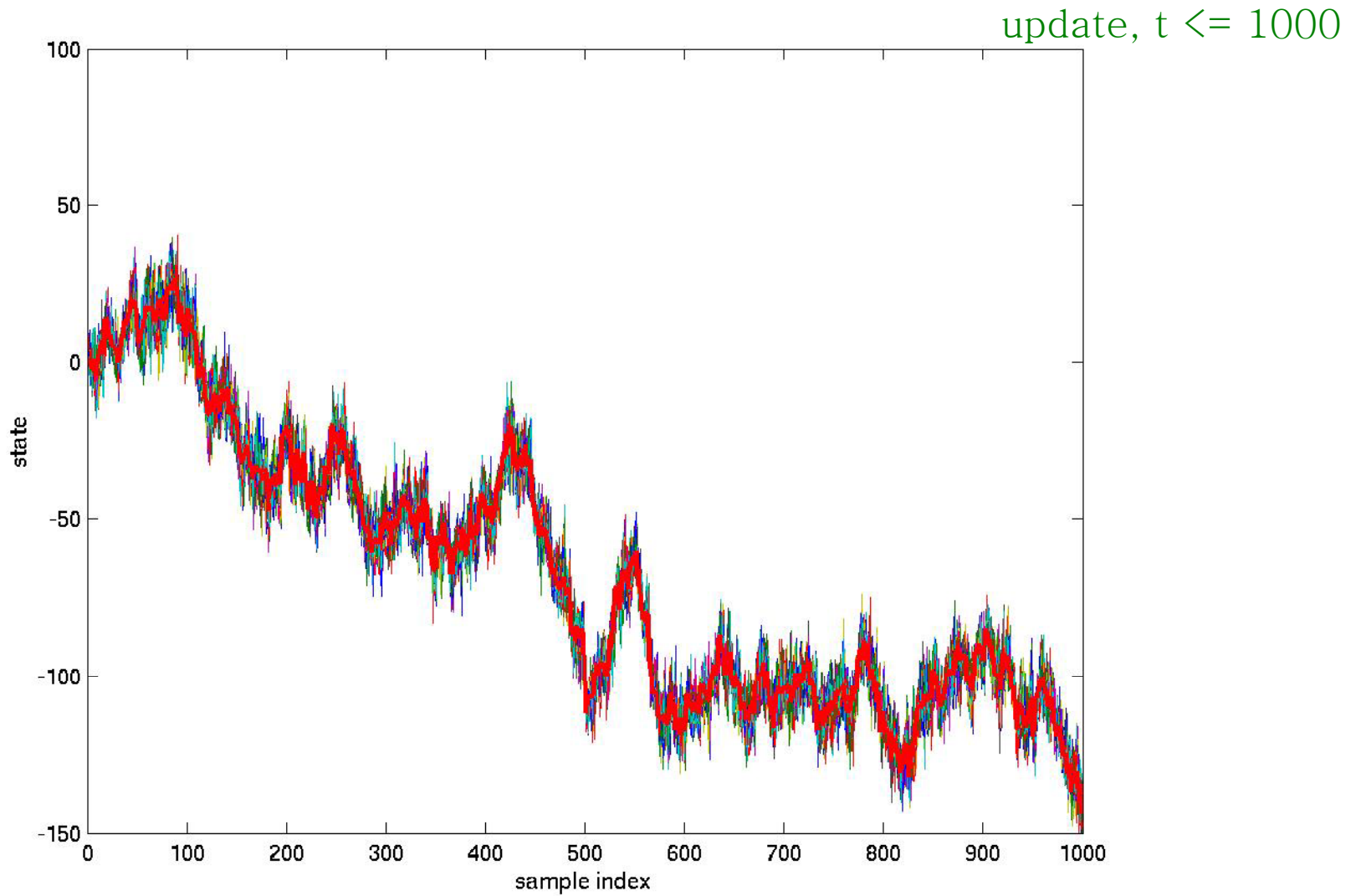
# Updated Probs Until $T = 1000$



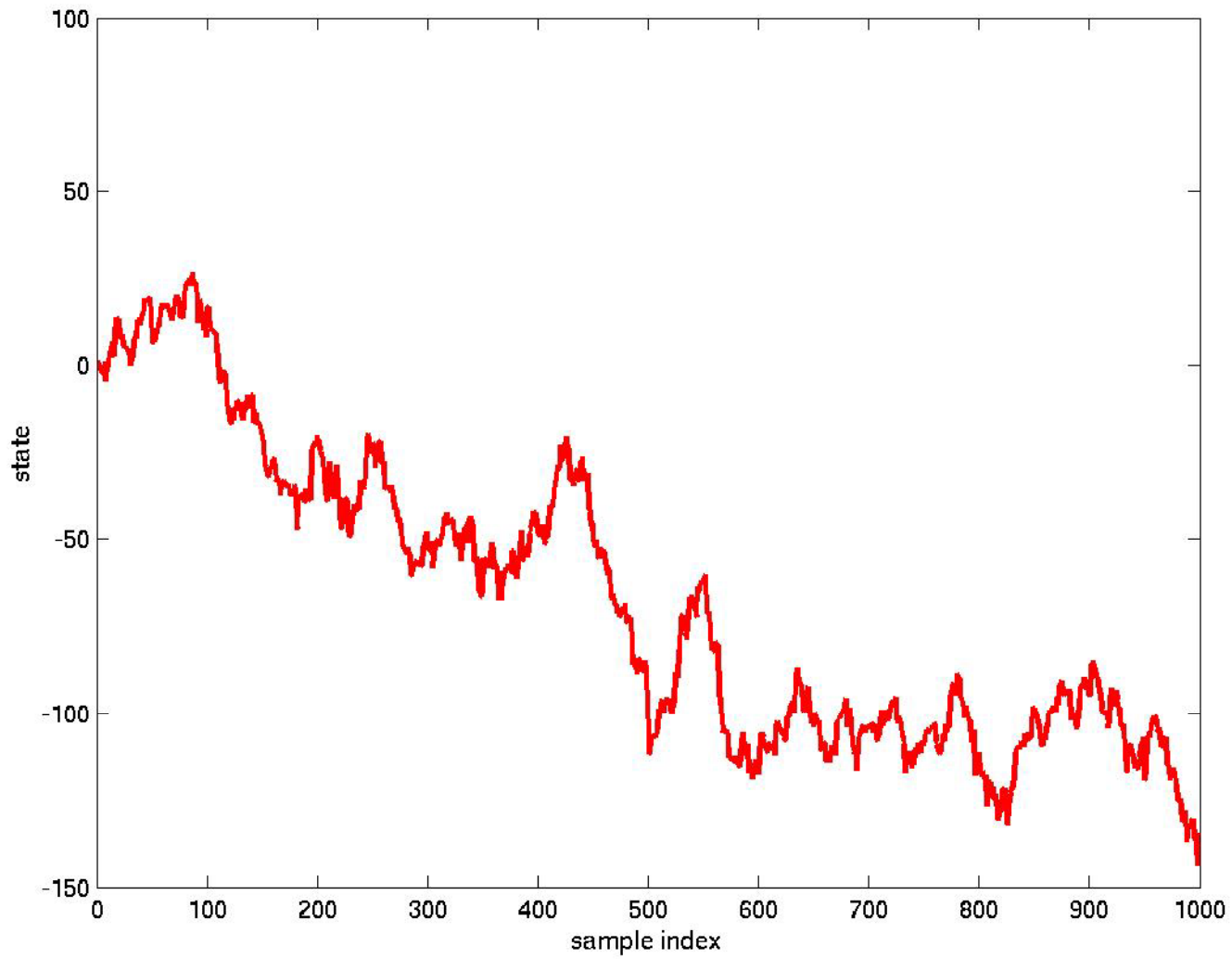
# Updated Probs Until $T = 1000$



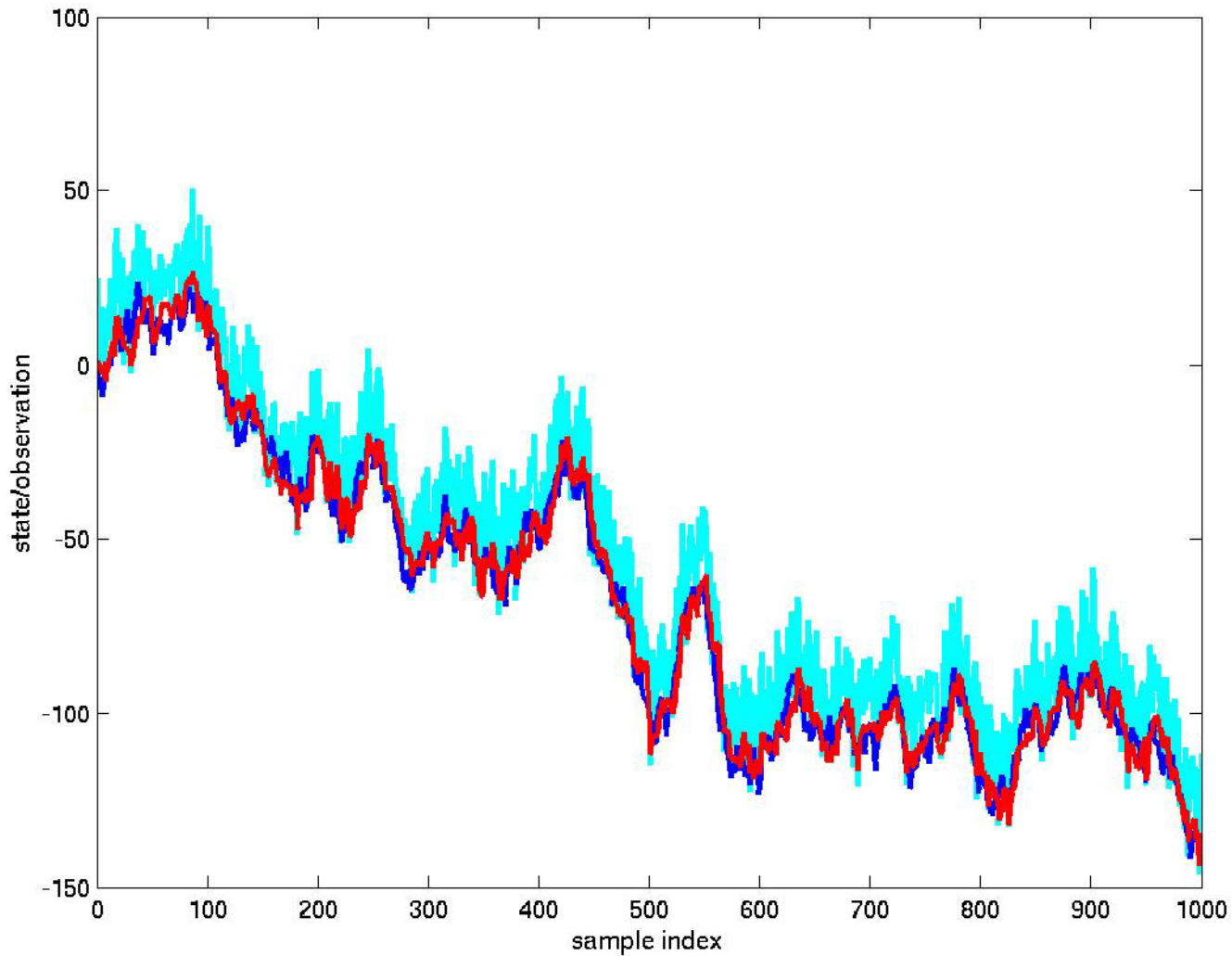
# Updated Probs: Top View



# ESTIMATED STATE



# Observation, True States, Estimate





# 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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# Additional Topics

- NMF / ICA.
- Class on 30 Nov?