## 4.0 More about Hidden Markov Models

Reference: 1. 6.1-6.6, Rabiner and Juang

2. 4.4.1 of Huang

# **Markov Model**

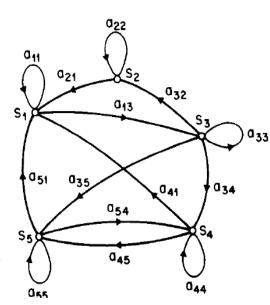
## Markov Model (Markov Chain)

- First-order Markov chain of N states is a triplet  $(S, A, \pi)$ 
  - S is a set of N states
  - A is the  $N \times N$  matrix of state transition probabilities

$$P(q_t=j|q_{t-1}=i, q_{t-2}=k, \ldots)=P(q_t=j|q_{t-1}=i) \equiv \mathbf{a}_{ij}$$

•  $\pi$  is the vector of initial state probabilities  $\pi_i = P(q_0 = j)$ 

- The output for any given state is an observable event (deterministic)
- The output of the process is a sequence of observable events



A Markov chain with 5 states (labeled  $S_1$  to  $S_5$ ) with state transitions.

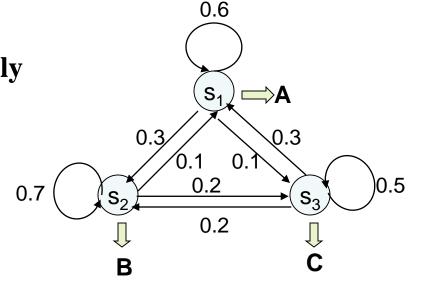
# **Markov Model**

## An example: a 3-state Markov Chain λ

State 1 generates symbol A *only*,
 State 2 generates symbol B **only**,
 and State 3 generates symbol C **only**

$$\mathbf{A} = \begin{bmatrix} 0.6 & 0.3 & 0.1 \\ 0.1 & 0.7 & 0.2 \\ 0.3 & 0.2 & 0.5 \end{bmatrix}$$

$$\pi = \begin{bmatrix} 0.4 & 0.5 & 0.1 \end{bmatrix}$$



- Given a sequence of observed symbols  $O=\{CABBCABC\}$ , the **only one** corresponding state sequence is  $\{S_3S_1S_2S_2S_3S_1S_2S_3\}$ , and the corresponding probability is

$$P(\mathbf{O}|\lambda) = P(q_0 = S_3)$$
  
 $P(S_1/S_3)P(S_2/S_1)P(S_2/S_2)P(S_3/S_2)P(S_1/S_3)P(S_2/S_1)P(S_3/S_2)$   
 $= 0.1 \times 0.3 \times 0.3 \times 0.7 \times 0.2 \times 0.3 \times 0.3 \times 0.2 = 0.00002268$ 

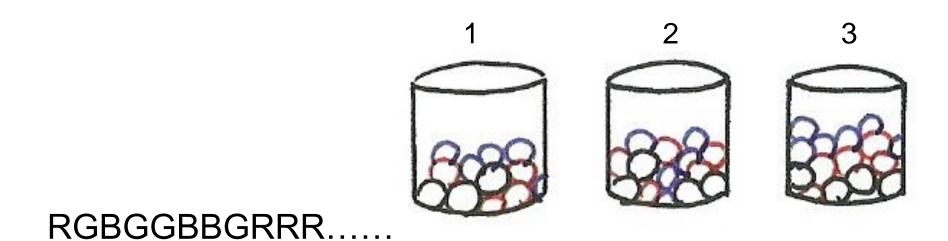
## HMM, an extended version of Markov Model

- The observation is a probabilistic function (discrete or continuous)
   of a state instead of an one-to-one correspondence of a state
- The model is a doubly embedded stochastic process with an underlying stochastic process that is not directly observable (hidden)
  - What is hidden? *The State Sequence*According to the observation sequence, we never know which state sequence generates it

## • Elements of an HMM $\{S,A,B,\pi\}$

- S is a set of N states
- A is the  $N \times N$  matrix of state transition probabilities
- B is a set of N probability functions, each describing the observation probability with respect to a state
- $-\pi$  is the vector of initial state probabilities

# Simplified HMM



## Two types of HMM's according to the observation functions

## Discrete and finite observations:

- The observations that all distinct states generate are finite in number  $\mathbf{V} = {\mathbf{v}_1, \mathbf{v}_2, \mathbf{v}_3, \ldots, \mathbf{v}_M}, \mathbf{v}_k \in \mathbf{R}^D$
- the set of observation probability distributions  $B = \{b_j(\mathbf{v}_k)\}\$  is defined as  $b_j(\mathbf{v}_k) = P(\mathbf{o}_t = \mathbf{v}_k | \mathbf{q}_t = j), 1 \le k \le M, 1 \le j \le N$ 
  - $\mathbf{o}_{t}$ : observation at time t,  $\mathbf{q}_{t}$ : state at time t
  - $\Rightarrow$  for state j,  $b_j(\mathbf{v}_k)$  consists of only M probability values

#### Continuous and infinite observations:

- The observations that all distinct states generate are infinite and continuous,  $\mathbf{V} = \{ \mathbf{v} | \mathbf{v} \in \mathbf{R}^D \}$
- the set of observation probability distributions  $B=\{b_j(\mathbf{v})\}$  is defined as  $b_i(\mathbf{v})=P(\mathbf{o}_t=\mathbf{v}|\mathbf{q}_t=j), 1 \le j \le N$ 
  - $\Rightarrow b_j(\mathbf{v})$  is a continuous probability density function and is often assumed to be a mixture of Gaussian distributions

$$b_{j}(\mathbf{v}) = \sum_{k=1}^{M} c_{jk} \left( \frac{1}{\left(\sqrt{2\pi}\right)^{D} \left|\Sigma_{jk}\right|^{\frac{1}{2}}} \exp\left(-\frac{1}{2} \left( \left(\mathbf{v} - \boldsymbol{\mu}_{jk}\right)^{T} \Sigma_{jk}^{-1} \left(\mathbf{v} - \boldsymbol{\mu}_{jk}\right) \right) \right) \right) = \sum_{k=1}^{M} c_{jk} b_{jk}(V)$$

• An example : a 3-state discrete HMM  $\lambda$ 

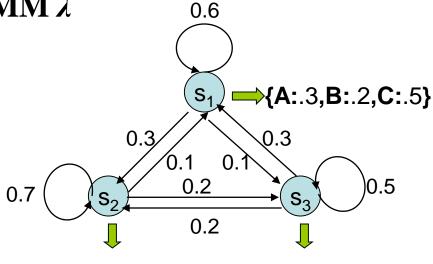
$$\mathbf{A} = \begin{bmatrix} 0.6 & 0.3 & 0.1 \\ 0.1 & 0.7 & 0.2 \\ 0.3 & 0.2 & 0.5 \end{bmatrix}$$

$$b_1(\mathbf{A}) = 0.3, b_1(\mathbf{B}) = 0.2, b_1(\mathbf{C}) = 0.5$$

$$b_2(\mathbf{A}) = 0.7, b_2(\mathbf{B}) = 0.1, b_2(\mathbf{C}) = 0.2$$

$$b_3(\mathbf{A}) = 0.3, b_3(\mathbf{B}) = 0.6, b_3(\mathbf{C}) = 0.1$$

$$\pi = \begin{bmatrix} 0.4 & 0.5 & 0.1 \end{bmatrix}$$



**{A:**.7,**B:**.1,**C:**.2**} {A:**.3,**B:**.6,**C:**.1**}** 

Given a sequence of observations O={ABC}, there are 27 possible corresponding state sequences, and therefore the corresponding probability is

$$\begin{split} P(\overline{\mathbf{O}}|\lambda) &= \sum_{i=1}^{27} P(\overline{\mathbf{O}}, \mathbf{q}_i | \lambda) = \sum_{i=1}^{27} P(\overline{\mathbf{O}}|\mathbf{q}_i, \lambda) P(\mathbf{q}_i | \lambda), \quad \mathbf{q}_i : \text{state sequence} \\ e.g. \text{ when } \mathbf{q}_i &= \left\{ S_2 S_2 S_3 \right\}, P(\overline{\mathbf{O}}|\mathbf{q}_i, \lambda) = P(\mathbf{A}|S_2) P(\mathbf{B}|S_2) P(\mathbf{C}|S_3) = 0.7*0.1*0.1 = 0.007 \\ P(\mathbf{q}_i | \lambda) &= P(q_0 = S_2) P(S_2 | S_2) P(S_3 | S_2) = 0.5*0.7*0.2 = 0.07 \end{split}$$

```
    Three Basic Problems for HMMs

 Given an observation sequence \overline{O}=(o_1,o_2,\ldots,o_T), and an HMM
\lambda = (A,B,\pi)
     – Problem 1 :
         How to efficiently compute P(\overline{O}|\lambda)?
         \Rightarrow Evaluation problem
     - Problem 2:
         How to choose an optimal state sequence \mathbf{q} = (q_1, q_2, \dots, q_T)?
          \Rightarrow Decoding Problem
     - Problem 3:
         Given some observations \overline{O} for the HMM \lambda, how to adjust the
         model parameter \lambda = (A,B,\pi) to maximize P(\overline{O}|\lambda)?
         ⇒ Learning /Training Problem
```

$$\begin{aligned} O &= o_1 o_2 o_3 ..... o_t ..... o_T & observation sequence \\ \overline{q} &= q_1 q_2 q_3 ..... q_t ..... q_T & state sequence \end{aligned}$$

- **Problem 1:** Given  $\lambda$  and O, find  $P(O|\lambda)=Prob[observing O given <math>\lambda]$
- Direct Evaluation: considering all possible state sequence  $\overline{q}$

$$\begin{split} P(\overline{O}|\lambda) &= \sum_{all \ \bar{q}} P(\overline{O}, \overline{q}|\lambda) = \sum_{all \ \bar{q}} P(\overline{O}|\overline{q}, \lambda) P(\overline{q}|\lambda) \\ &\qquad P(\overline{O}|\bar{q}, \lambda) \\ P(\overline{O}|\lambda) &= \sum_{all \ \bar{q}} ([b_{q_1}(o_1) \bullet b_{q_2}(o_2) \bullet \ldots ... .b_{q_T}(o_T)] \bullet \\ &\qquad [\pi_{q_1} \bullet a_{q_1q_2} \bullet a_{q_2q_3} \bullet \ldots ... a_{q_{T-1}q_T}]) \\ &\qquad \qquad \square \\ P(\bar{q}|\lambda) \end{split}$$

total number of different  $\overline{q}: N^T$  huge computation requirements

• Forward Algorithm: defining a forward variable  $\alpha_t(i)$ 

$$\alpha_{t}(i) = P(o_{1}o_{2}...o_{t}, q_{t} = i|\lambda)$$

$$= Prob[observing o_{1}o_{2}...o_{t}, state i at time t|\lambda]$$

- Initialization

$$\alpha_1(i) = \pi_i b_i(o_1) , \quad 1 \le i \le N$$

- Induction

$$\alpha_{t+1}(j) = \left[\sum_{i=1}^{N} \alpha_{t}(i) a_{ij}\right] b_{j}(o_{t+1})$$

$$1 \le j \le N$$

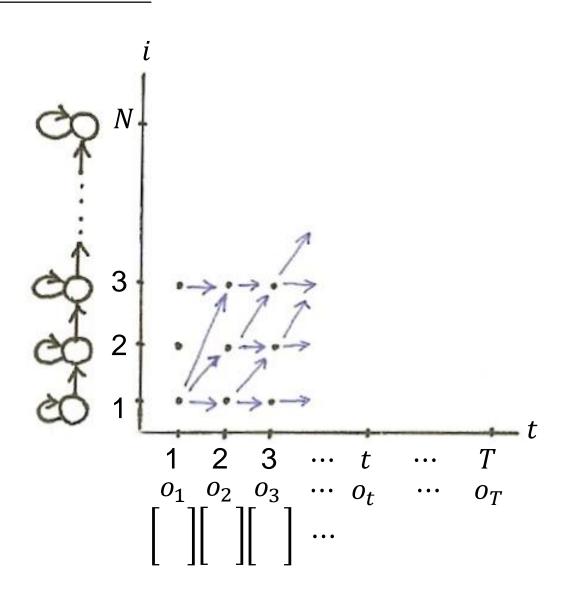
$$1 \le t \le T-1$$

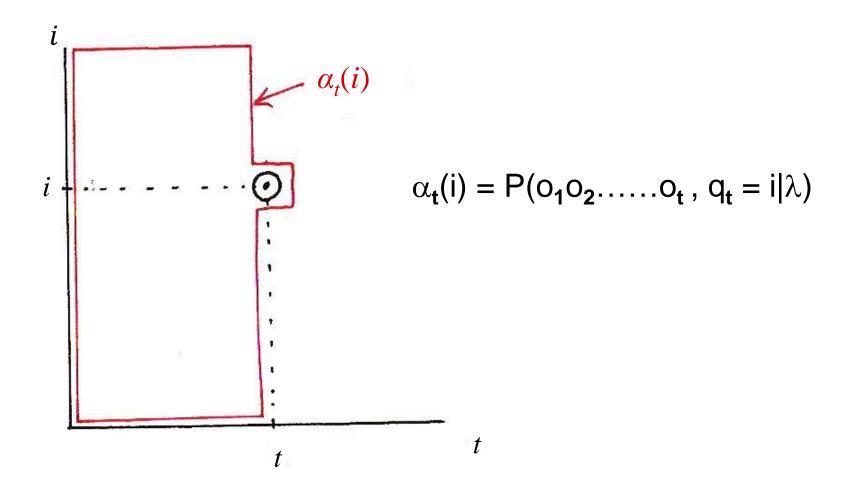
- Termination

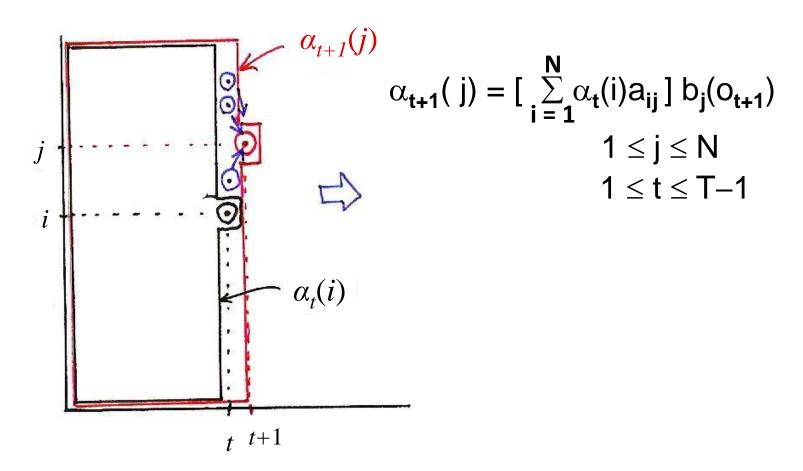
$$P(\overline{O}|\lambda) = \sum_{i=1}^{N} \alpha_{T}(i)$$

See Fig. 6.5 of Rabiner and Juang

- All state sequences, regardless of how long previously, merge to the N state at each time instant t







# Forward Algorithm

- **Problem 2:** Given  $\lambda$  and  $\overline{O} = o_1 o_2 \dots o_T$ , find a best state sequence  $\overline{q} = q_1 q_2 \dots q_T$
- Backward Algorithm : defining a backward variable  $\beta_t(i)$

$$\beta_{t}(i) = P(o_{t+1}, o_{t+2}, ..., o_{T} | q_{t}=i, \lambda)$$

$$= Prob[observing o_{t+1}, o_{t+2}, ..., o_{T} | state i at time t, \lambda]$$

- Initialization

$$\beta_{\mathbf{T}}(i) = 1, \quad 1 \le i \le N$$
  $(\beta_{\mathbf{T-1}}(i) = \sum_{i=1}^{N} a_{ij} b_{j}(o_{\mathbf{T}}))$ 

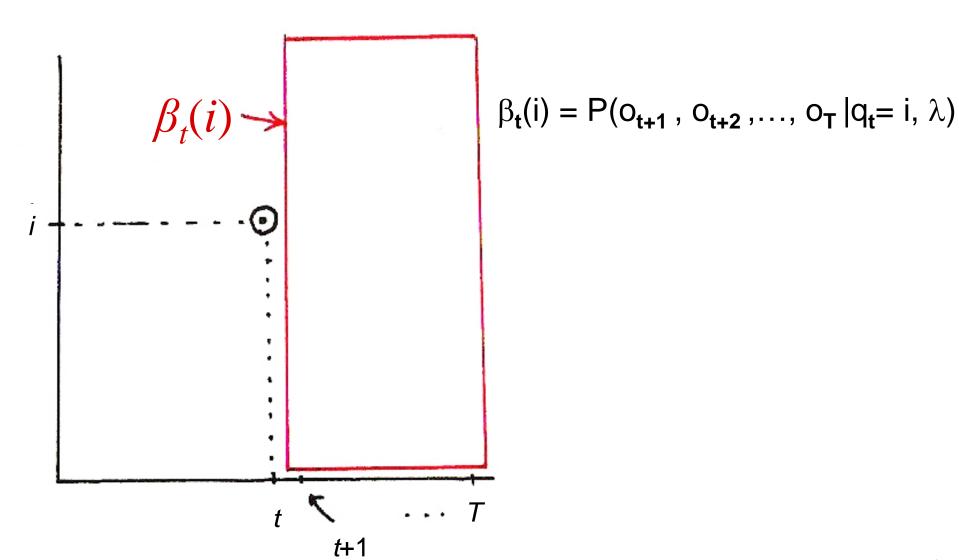
- Induction

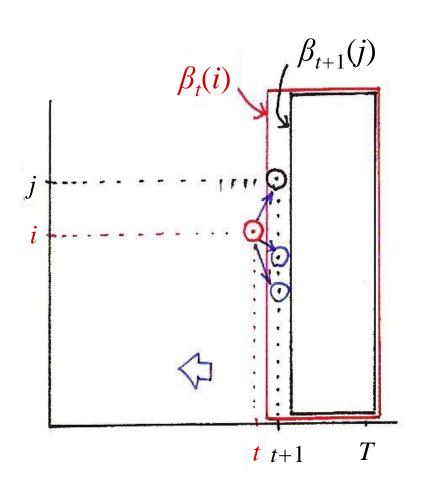
$$\begin{split} \beta_{t}(i) = & \sum_{j=1}^{N} a_{ij} \ b_{j}(o_{t+1}) \beta_{t+1}(j) \\ t = T-1, \ T-2, \dots, \ 2, \ 1, \qquad 1 \leq i \leq N \end{split}$$

See Fig. 6.6 of Rabiner and Juang

#### Combining Forward/Backward Variables

$$\begin{split} &P(\overline{O}, q_t = i \mid \lambda) \\ &= \text{Prob [observing o_1, o_2, ..., o_t, ..., o_T, q_t = i \mid \lambda]} \\ &= \alpha_t(i)\beta_t(i) \\ &P(\overline{O}\mid \lambda) = \sum_{i=1}^N P(\overline{O}, q_t = i \mid \lambda) = \sum_{i=1}^N \left[\alpha_t(i)\beta_t(i)\right] \end{split}$$

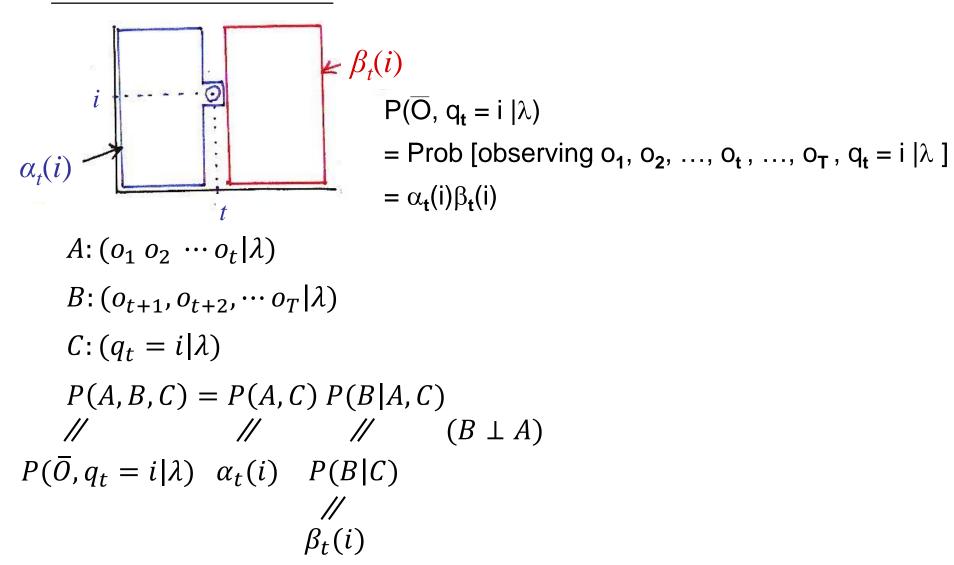




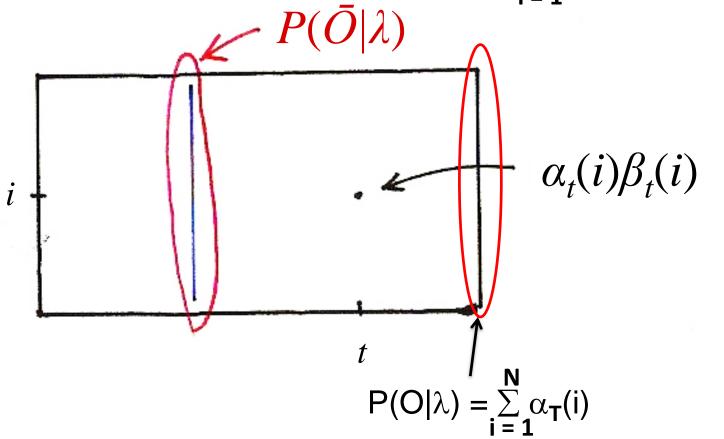
$$\beta_{t}(i) = \sum_{j=1}^{N} a_{ij} b_{j}(o_{t+1}) \beta_{t+1}(j)$$

$$t = T-1, T-2,..., 2, 1, 1 \le i \le N$$

Backward Algorithm



$$P(\overline{O}|\lambda) = \sum_{i=1}^{N} P(\overline{O}, q_t = i | \lambda) = \sum_{i=1}^{N} [\alpha_t(i)\beta_t(i)]$$



- Approach 1 Choosing state q<sub>t</sub>\* individually as the most likely state at time t
  - Define a new variable  $\gamma_t(i) = P(q_t = i \mid \overline{O}, \lambda)$

$$\gamma_{t}(i) = \frac{\alpha_{t}(i)\beta_{t}(i)}{\sum_{i=1}^{N} \alpha_{t}(i)\beta_{t}(i)} = \frac{P(O, q_{t}=i|\lambda)}{P(O|\lambda)}$$

- Solution

$$\begin{aligned} q_t^* &= \arg\max_{1 \leq i \leq N} \left[ \!\! \left[ \!\! \left[ t(i) \right] \!\! \right], \, 1 \leq t \leq T \right] \\ &\inf \text{fact} \\ q_t^* &= \arg\max_{1 \leq i \leq N} \left[ P(\overline{O}, \, q_t = i | \lambda) \right] \\ &= \arg\max_{1 \leq i \leq N} \left[ \alpha_t(i) \beta_t(i) \right] \end{aligned}$$

- Problem maximizing the probability at each time t individually  $\overline{q}^*=q_1^*q_2^*...q_T^* \text{ may not be a valid sequence}$  (e.g.  $a_{q_t^*q_{t+1}^*}=0$ )

- Approach 2 —Viterbi Algorithm finding the single best sequence  $\overline{q}^* = q_1^* q_2^* \dots q_T^*$ 
  - Define a new variable  $\delta_t(i)$

$$\delta_{t}(i) = \max_{q_{1}, q_{2}, \dots, q_{t-1}} P[q_{1}, q_{2}, \dots, q_{t-1}, q_{t} = i, o_{1}, o_{2}, \dots, o_{t} | \lambda]$$

- = the highest probability along a certain single path ending at state i at time t for the first t observations, given  $\lambda$
- Induction

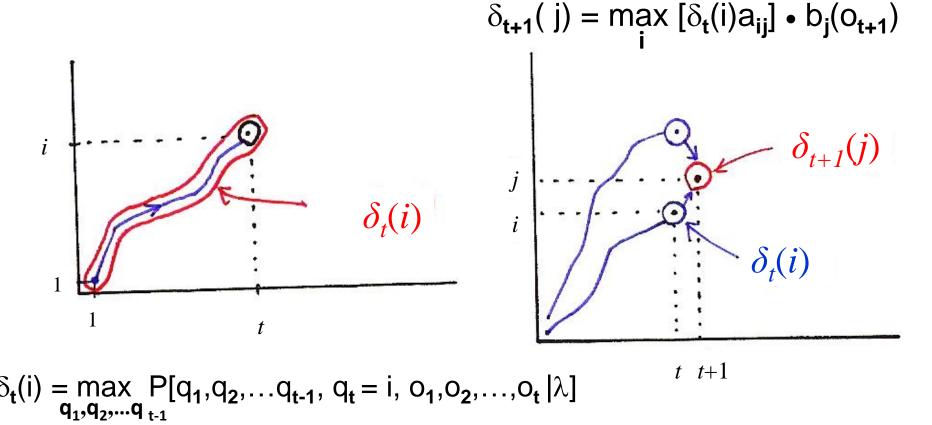
$$\delta_{t+1}(j) = \max_{i} [\delta_{t}(i)a_{ij}] \cdot b_{j}(o_{t+1})$$

- Backtracking

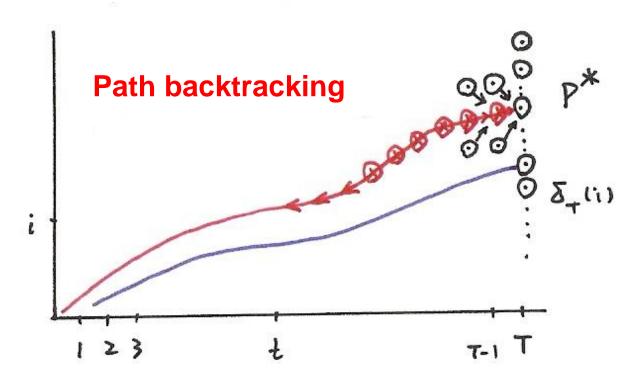
$$\psi_{t+1}(j) = \arg \max_{1 \le i \le N} [\delta_t(i)a_{ij}]$$

the best previous state at t-1 given at state j at time t keeping track of the best previous state for each j and t

# Viterbi Algorithm



# Viterbi Algorithm



- Complete Procedure for Viterbi Algorithm
  - Initialization

$$\delta_1(i) = \pi_i b_i(o_1) , \quad 1 \le i \le N$$

- Recursion

$$\begin{split} \delta_{t+1}(j) &= \max_{1 \leq i \leq N} \left[ \delta_t(i) a_{ij} \right] \bullet b_j(o_{t+1}) \\ &1 \leq t \leq T\text{-}1, \quad 1 \leq j \leq N \\ \psi_{t+1}(j) &= \underset{1 \leq i \leq N}{arg \max} \left[ \delta_t(i) a_{ij} \right] \\ &1 \leq t \leq T\text{-}1, \quad 1 \leq j \leq N \end{split}$$

- Termination

$$\begin{aligned} P^* &= \max_{1 \leq i \leq N} \left[ \delta_T(i) \right] \\ q_T^* &= arg \max_{1 \leq i \leq N} \left[ \delta_T(i) \right] \end{aligned}$$

- Path backtracking

$$q_t^* = \psi_{t+1}(q^*_{t+1}), \quad t = T-1, t-2, \dots, 2, 1$$

#### • Application Example of Viterbi Algorithm

- Isolated word recognition

$$\lambda_0 = (A_0, B_0, \boldsymbol{\pi}_0)$$

$$\lambda_1 = (A_1, B_1, \boldsymbol{\pi}_1)$$

$$\vdots$$

$$\lambda_n = (A_n, B_n, \boldsymbol{\pi}_n)$$

observation

$$\overline{\mathbf{O}} = (o_1, o_2, ... o_T)$$

$$k^* = \arg \max_{\mathbf{1} \le \mathbf{i} \le \mathbf{n}} \mathbf{P}[\overline{\mathbf{O}} \mid \lambda_i] \approx \arg \max_{\mathbf{1} \le \mathbf{i} \le \mathbf{n}} [\mathbf{P}^* \mid \lambda_i]$$

$$\widehat{\Box}$$

Basic Problem 1 Basic Problem 2
Forward Algorithm (for all paths) Basic Problem 2
Viterbi Algorithm (for a single best path)

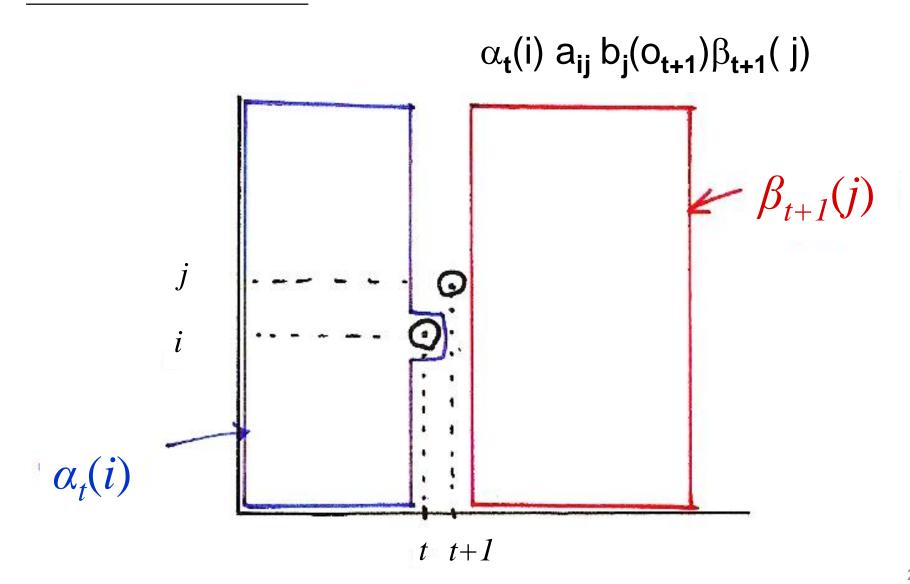
-The model with the highest probability for the most probable path usually also has the highest probability for all possible paths.

- **Problem 3:** Give  $\overline{O}$  and an initial model  $\lambda = (A, B, \pi)$ , adjust  $\lambda$  to maximize  $P(\overline{O}|\lambda)$ 
  - Baum-Welch Algorithm (Forward-backward Algorithm)
  - Define a new variable

$$\begin{split} & \mathcal{E}_{t}(\ i,j\ ) = \ P(q_{t}=i,\,q_{t+1}=j\ |\ \overline{O},\,\lambda) \\ & = \frac{\alpha_{t}(i)\ a_{ij}\ b_{j}(o_{t+1})\beta_{t+1}(j)}{\sum\limits_{i=1}^{N}\sum\limits_{j=1}^{N}\left[\alpha_{t}(i)a_{ij}\ b_{j}(o_{t+1})\beta_{t+1}(j)\right]} \\ & = \frac{Prob[\overline{O},\,q_{t}=i,\,q_{t+1}=j|\lambda]}{P(\overline{O}|\lambda)} \end{split}$$

See Fig. 6.7 of Rabiner and Juang

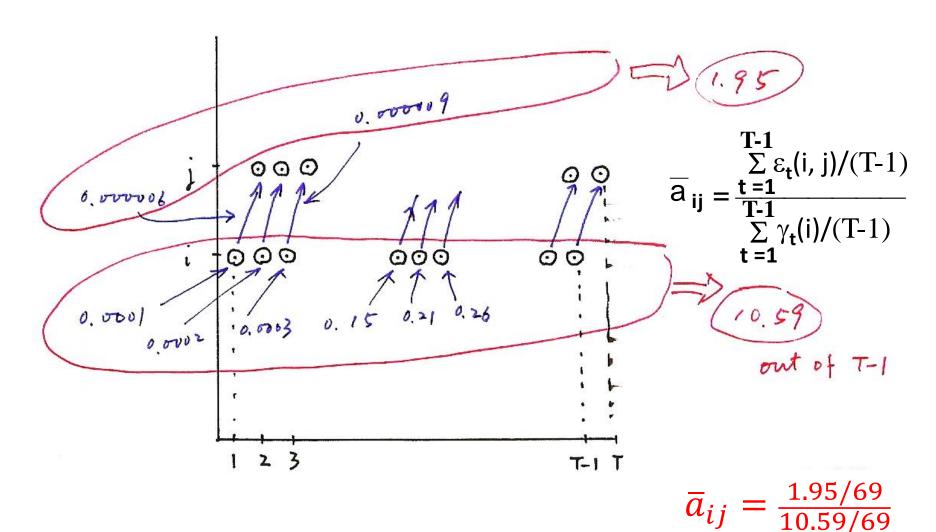
- Recall  $\gamma_t(i) = P(q_t = i \mid \overline{O}, \lambda)$   $\sum_{t=1}^{T-1} \gamma_t(i) = \text{expected number of times that state } i \text{ is visited in } \overline{O} \text{ from } t = 1 \text{ to } t = T-1$   $= \text{expected number of transitions from state } i \text{ in } \overline{O}$   $\sum_{t=1}^{T-1} \epsilon_t(i,j) = \text{expected number of transitions from state } i \text{ to state } j \text{ in } \overline{O}$ 



# Basic Problem 3 $\alpha_t(i)a_{ij}b_j(o_{t+1})\beta_{t+1}(j)$ $\beta_{t+1}(j)$ $\alpha_t(i)$ t t+1

$$\gamma_t(i) = \frac{\alpha_t(i) \beta_t(i)}{\sum_{i=1}^{N} [\alpha_t(i) \beta_t(i)]} = \frac{P(\overline{O}, q_t = i | \lambda)}{P(\overline{O} | \lambda)} = P(q_t = i | \overline{O}, \lambda)$$

$$\begin{split} \varepsilon_{t}(i,j) &= \frac{\alpha_{t}(i) \; a_{ij} \; b_{j}(o_{t+1}) \; \beta_{t+1}(j)}{\sum_{j=1}^{N} \sum_{i=1}^{N} \alpha_{t}(i) \; a_{ij} \; b_{j}(o_{t+1}) \; \beta_{t+1}(j)} \\ &= \frac{P(\bar{O}, q_{t} = i, q_{t+1} = j | \lambda)}{P(\bar{O} | \lambda)} = P(q_{t} = i, q_{t+1} = j | \bar{O}, \lambda) \end{split}$$



- Results

$$\begin{split} \overline{\pi}_i &= \gamma_1(i) \\ \overline{a}_{ij} &= \frac{\sum\limits_{t=1}^{T-1} \epsilon_t(i,j)}{\sum\limits_{t=1}^{T-1} \gamma_t(i)} \\ \overline{b}_j(k) &= \text{Prob}[o_t = v_k \mid q_t = j \ ] = \frac{\sum\limits_{t=1}^{T} \gamma_t(j)}{\sum\limits_{t=1}^{T} \gamma_t(j)} \end{split}$$
 (for discrete HMM)

#### Continuous Density HMM

$$b_{j}(o) = \sum_{k=1}^{M} c_{jk} N(o; \mu_{jk}, U_{jk})$$

N(): Multi-variate Gaussian

 $\mu_{jk}$ : mean vector for the k-th mixture component

U<sub>jk</sub>: covariance matrix for the k-th mixture component

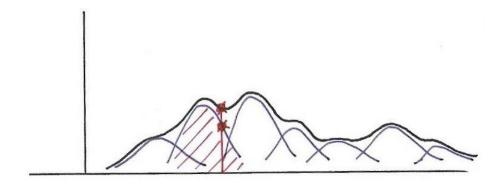
$$\sum_{k=1}^{M} c_{jk} = 1 \text{ for normalization}$$

#### Continuous Density HMM

- Define a new variable

 $\gamma_t(j, k) = \gamma_t(j)$  but including the probability of  $o_t$  evaluated in the k-th mixture component out of all the mixture components

$$= \left(\frac{\alpha_t(j)\beta_t(j)}{\sum\limits_{j=1}^{N}\alpha_t(j)\beta_t(j)}\right) \left(\frac{c_{jk}N(o_t;\,\mu_{jk},\,U_{jk})}{\sum\limits_{m=1}^{M}c_{jm}N(o_t;\,\mu_{jm},\,U_{jm})]}\right)$$



- Results

$$\overline{c}_{jk} = \frac{\sum_{t=1}^{T} \gamma_t(j, k)}{\sum_{t=1}^{T} \sum_{k=1}^{M} \gamma_t(j, k)}$$

See Fig. 6.9 of Rabiner and Juang

## Continuous Density HMM

$$\overline{\mu}_{jk} = \frac{\sum_{t=1}^{T} [\gamma_t(j, k) \cdot o_t]}{\sum_{t=1}^{T} \gamma_t(j, k)}$$

$$\overline{U}_{jk} = \frac{\sum_{t=1}^{T} [\gamma_t(j, k)(o_t - \mu_{jk}) (o_t - \mu_{jk})']}{\sum_{t=1}^{T} \gamma_t(j, k)}$$

#### • Iterative Procedure

$$\lambda = (A, B, \pi) \xrightarrow{\overline{\lambda}} \overline{\lambda} = (\overline{A}, \overline{B}, \overline{\pi})$$

$$\overline{O} = o_1 o_2 \dots o_T$$

- It can be shown (by EM Theory (or EM Algorithm))  $P(\overline{O}|\overline{\lambda}) \ge P(\overline{O}|\lambda) \text{ after each iteration}$ 

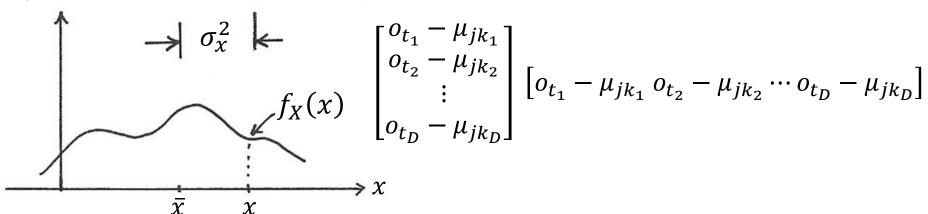
$$\overline{\mu}_{jk} = \frac{\sum_{t=1}^{T} \left[ \gamma_t(j,k) \bullet o_t \right]}{\sum_{t=1}^{T} \gamma_t(j,k)}$$

$$\int_{-\infty}^{\infty} x \left[ f_X(x) \right] dx = \bar{x}$$

$$\overline{U}_{jk} = \frac{\sum_{t=1}^{T} [\gamma_t(j,k)] (o_t - \mu_{jk}) (o_t - \mu_{jk})'}{\sum_{t=1}^{T} \gamma_t(j,k)} \qquad \int_{-\infty}^{\infty} [(x - \overline{x})^2] f_X(x) dx = \sigma_X^2$$

$$\int_{-\infty}^{\infty} \left[ (x - \bar{x})^2 \right] f_X(x) \ dx = \sigma_X^2$$

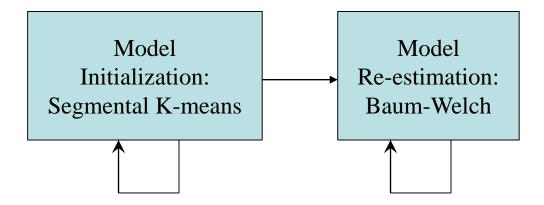
## $f_X(x)$ : prob. density function

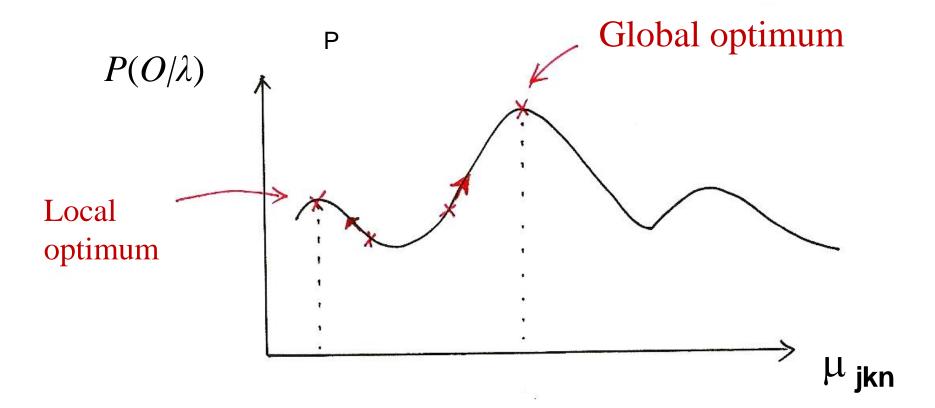


$$\overline{U} = \begin{bmatrix}
\vdots \\
\vdots \\
\vdots \\
\vdots \\
\vdots \\
\vdots
\end{bmatrix} = E(\begin{bmatrix} x_1 - \overline{x}_1 \\ x_2 - \overline{x}_2 \\ \vdots \\ \vdots \end{bmatrix} [x_1 - \overline{x}_1, x_2 - \overline{x}_2, \cdots])$$

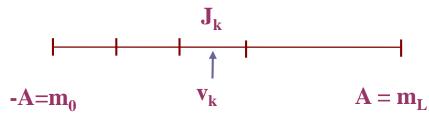
$$\bar{u}_{lm} = E[(x_l - \bar{x}_l)(x_m - \bar{x}_m)]$$

- No closed-form solution, but approximated iteratively
- An initial model is needed-model initialization
- May converge to local optimal points rather than global optimal point
  - heavily depending on the initialization
- Model training





- An Efficient Approach for Data Compression
  - replacing a set of real numbers by a finite number of bits
- An Efficient Approach for Clustering Large Number of Sample Vectors
  - grouping sample vectors into clusters, each represented by a single vector (codeword)
- Scalar Quantization
  - replacing a single real number by an R-bit pattern
  - a mapping relation



$$S = \bigcup_{k=1}^{L} J_k, V = \{ v_1, v_2, ..., v_L \}$$

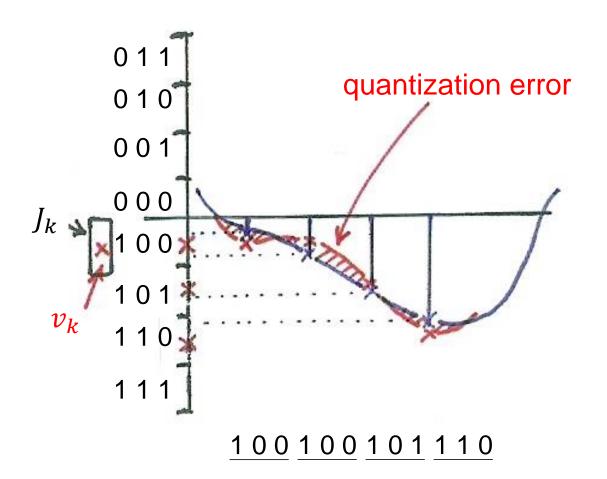
$$Q: S \to V$$

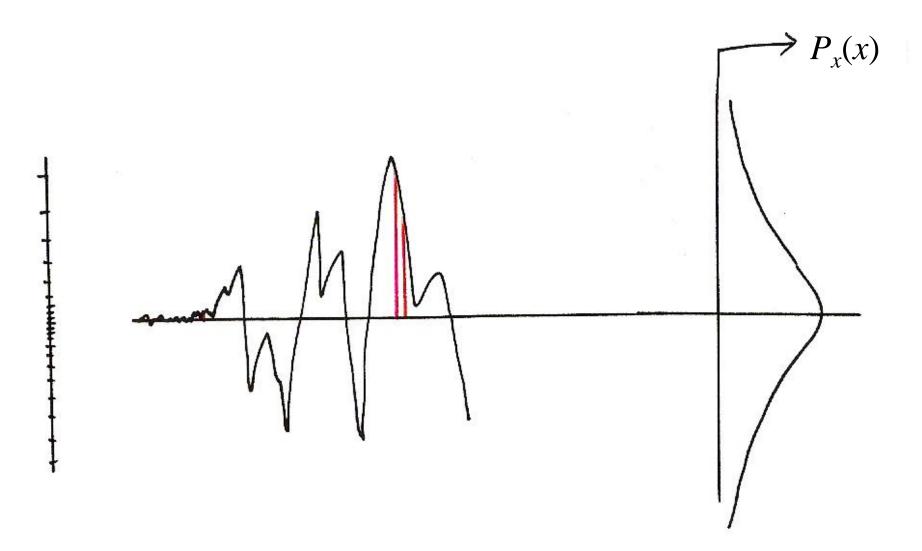
$$Q(x[n]) = v_k \text{ if } x[n] \in J_k$$

$$L = 2^R$$
Each  $v_k$  represented by an R-bit pattern

- Quantization characteristics (codebook)
   { J<sub>1</sub>, J<sub>2</sub>, ..., J<sub>L</sub> } and { v<sub>1</sub>, v<sub>2</sub>, ..., v<sub>L</sub> }
   designed considering at least
   1 error sensitivity
  - 1. error sensitivity
  - 2. probability distribution of x[n]

Scalar Quantization: Pulse Coded Modulation (PCM)





### 2-dim Vector Quantization (VQ)

#### Example:

```
\overline{x}_{n} = (x[n], x[n+1])

S = {\overline{x}_{n} = (x[n], x[n+1]); |x[n]| < A, |x[n+1]| < A}
```

### •VQ

-S divided into L 2-dim regions  $\{J_1, J_2, ..., J_k, ..., J_L\}$ 

$$S = \bigcup_{k=1}^{L} J_k$$

each with a representative

vector 
$$\overline{v}_k \in J_k$$
,  $V = \{\overline{v}_1, \overline{v}_2, ..., \overline{v}_L\}$ 

$$-Q: S \rightarrow V$$

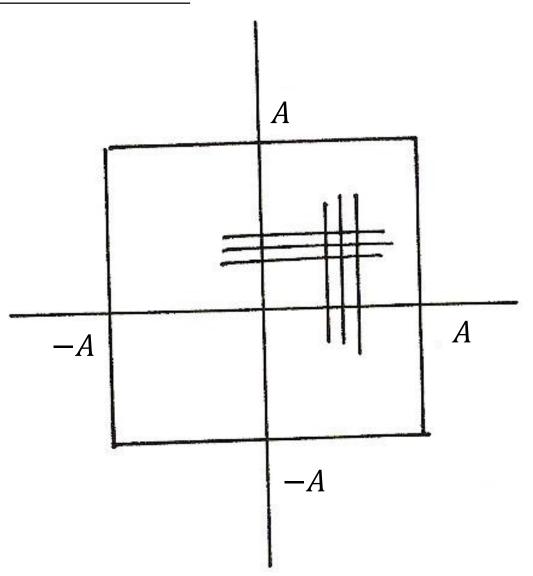
$$Q(\overline{x}_n) = \overline{v}_k$$
 if  $\overline{x}_n \in J_k$ 

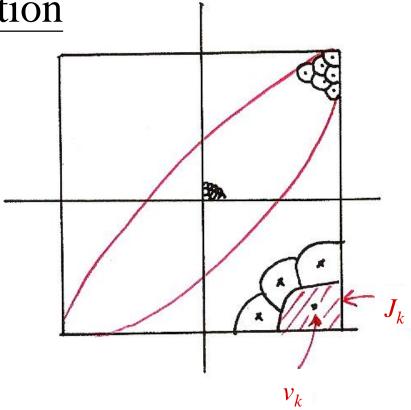
$$L=2^{\mathbf{R}}$$

each  $\overline{v}_k$  represented by an R-bit pattern

- Considerations
  - 1.error sensitivity may depend on x[n], x[n+1] jointly
  - 2.distribution of x[n], x[n+1] may be correlated statistically
  - 3.more flexible choice of  $J_k$
- Quantization Characteristics (codebook)

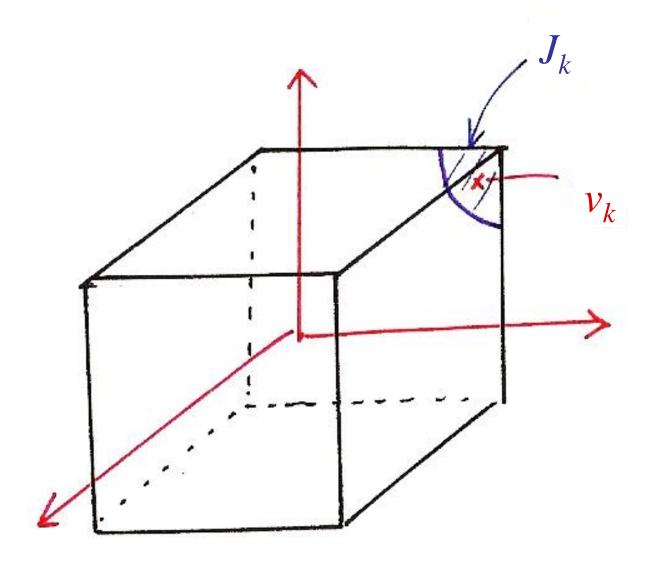
$$\{ J_1, J_2, ..., J_L \}$$
 and  $\{ \overline{v}_1, \overline{v}_2, ..., \overline{v}_L \}$ 





$$(256)^2 = (2^8)^2 = 2^{16}$$

$$1024=2^{10}$$



### N-dim Vector Quantization

$$\begin{split} \overline{x} &= (x_1\,,\,x_2\,,\,\ldots,\,x_N\,)\\ S &= \{\overline{x} = (x_1\,,\,x_2\,,\,\ldots,\,x_N)\,,\\ &\quad |\,x_k\,| < A\,,\,k = 1,2,\ldots N\}\\ S &= \mathop{\cup}\limits_{k=1}^L \,J_{_k}\\ V &= \{\overline{v}_1\,,\,\overline{v}_2\,,\,\ldots,\,\overline{v}_L\,\}\\ Q &: S \to V\\ Q(\overline{x}) &= \overline{v}_k \ \ \text{if} \ \ \overline{x} \in J_k\\ L &= 2^R\,,\, \text{each}\,\overline{v}_k \,\, \text{represented}\\ \text{by an R-bit pattern} \end{split}$$

# Codebook Trained by a Large Training Set

### Training Set

# Define distance measure between two vectors $\overline{\mathbf{x}}$ , $\overline{\mathbf{y}}$

$$d(\overline{x}, \overline{y}): S \times S \rightarrow R^+ \text{ (non-negative real numbers)}$$

-desired properties

$$\begin{aligned} &d(\overline{x}, \overline{y}\ ) \geq 0 \\ &d(\overline{x}, \overline{x}\ ) = 0 \\ &d(\overline{x}, \overline{y}\ ) = d(\overline{y}, \overline{x}\ ) \\ &d(\overline{x}, \overline{y}\ ) + d(\overline{y}, \overline{z}\ ) \geq d(\overline{x}, \overline{z}\ ) \end{aligned}$$

examples:

$$d(\overline{x}, \overline{y}) = \sum_{i} (x_{i} - y_{i})^{2}$$

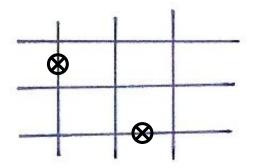
$$d(\overline{x}, \overline{y}) = \sum_{i} |x_{i} - y_{i}|$$

$$d(\overline{x}, \overline{y}) = (\overline{x} - \overline{y})^{t} \sum_{i} (\overline{x} - \overline{y})$$
Mahalanobis Distance

 $\Sigma$ : Co-variance Matrix

### **Distance Measures**

$$d(\bar{x}, \bar{y}) = \sum_{i} |x_i - y_i|$$
 city block distance

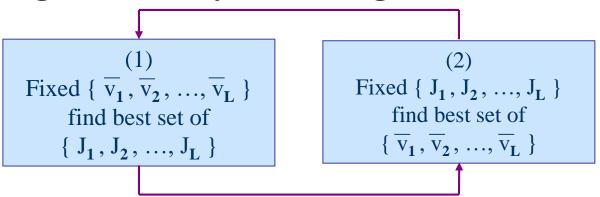


$$d(\bar{x}, \bar{y}) = (\bar{x} - \bar{y})^t \Sigma^{-1} (\bar{x} - \bar{y})$$
 Mahalanobis distance

$$\sum = \begin{bmatrix} 1 & \cdots & 0 \\ \vdots & \ddots & \vdots \\ 0 & \cdots & 1 \end{bmatrix}, d(\bar{x}, \bar{y}) = \sum_{i} (x_i - y_i)^2$$

$$\sum = \begin{vmatrix} \sigma_1^2 & \cdots & 0 \\ \vdots & \sigma_2^2 & \vdots \\ 0 & \cdots & \sigma_n^2 \end{vmatrix}, d(\bar{x}, \bar{y}) = \sum_i \frac{(x_i - y_i)^2}{\sigma_i^2}$$

### K-Means Algorithm/Lloyd-Max Algorithm



$$\begin{array}{l} (1) \; J_{k} = \{ \; \overline{x} \; | \; d(\overline{x} \; , \; \overline{v}_{k} \; ) < d(\overline{x} \; , \; \overline{v}_{j}) \; , \; j \neq k \; \} \\ \rightarrow D = & \sum_{\text{all } \overline{x}} d(\overline{x} \; , \; Q(\overline{x}) \; ) = min \\ \text{nearest neighbor condition} \\ \end{array}$$

(2) For each k
$$\overline{v}_{k} = \frac{1}{M} \sum_{x \in J_{k}} \overline{x}$$

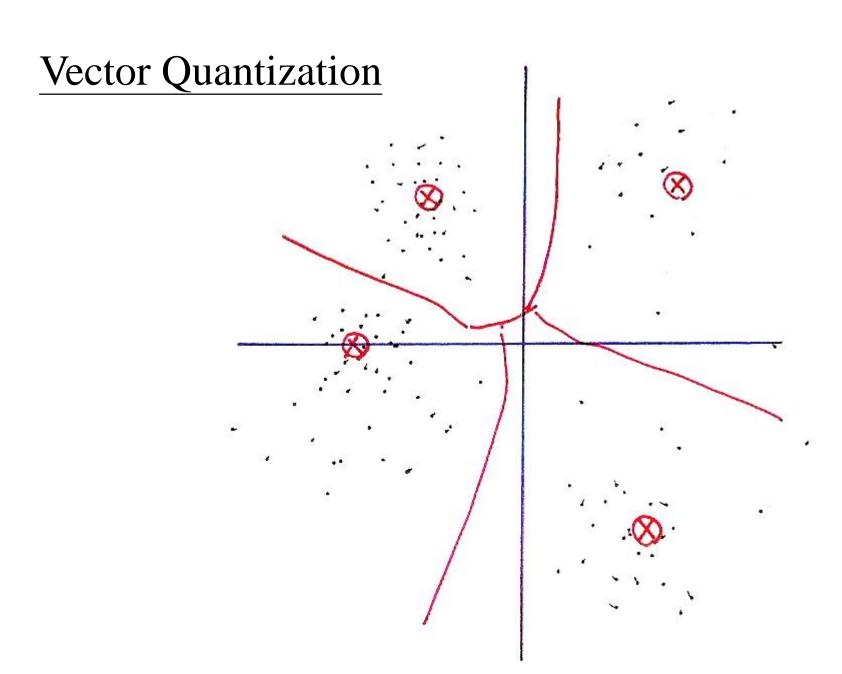
$$\rightarrow D_{k} = \sum_{\overline{x} \in J_{k}} d(\overline{x}, \overline{v}_{k}) = \min$$
centroid condition

(3) Convergence condition

$$D = \sum_{k=1}^{L} D_k$$

after each iteration D is reduced, but  $D \ge 0$   $\mid D^{(m+1)} - D^{(m)} \mid < \in$ , m : iteration

• Iterative Procedure to Obtain Codebook from a Large Training Set



- K-means Algorithm may Converge to Local Optimal Solutions
  - depending on initial conditions, not unique in general
- Training VQ Codebook in Stages— LBG Algorithm
  - step 1: Initialization. L = 1, train a 1-vector VQ codebook

$$\overline{v} = \frac{1}{N} \sum_{j} \overline{x}_{j}$$

- step 2: Splitting.

Splitting the L codewords into 2L codewords, L = 2L

• example 1

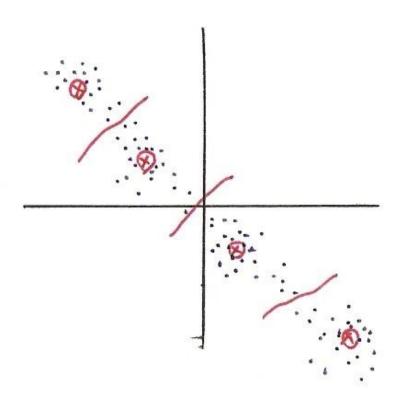
$$\overline{v}_{k}^{(1)} = \overline{v}_{k}(1+\varepsilon)$$

$$\overline{v}_{k}^{(2)} = \overline{v}_{k}(1-\varepsilon)$$

• example 2

- step 3: K-means Algorithm: to obtain L-vector codebook
- step 4: Termination. Otherwise go to step 2
- Usually Converges to Better Codebook

# **LBG Algorithm**



## **Initialization in HMM Training**

### • An Often Used Approach—Segmental K-Means

- Assume an initial estimate of all model parameters (e.g. estimated by segmentation of training utterances into states with equal length)
  - •For discrete density HMM

```
b_{j}(k) = \frac{\text{number of vectors in state } j \text{ associated with codeword } k}{\text{total number of vectors in state } j}
```

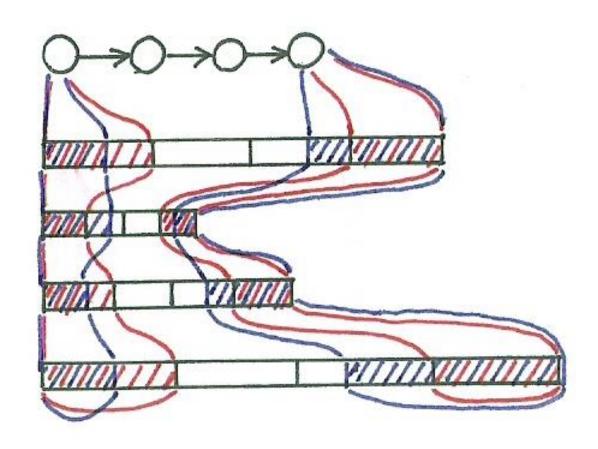
•For continuous density HMM (M Gaussian mixtures per state)

```
\Rightarrow \text{cluster the observation vectors within each state } j \text{ into a set of } M \text{ clusters}  (e.g. with vector quantization) c_{jm} = \text{number of vectors classified in cluster m of state } j  divided by number of vectors in state j \mu_{jm} = \text{sample mean of the vectors classified in cluster m of state } j  \sum_{jm} = \text{sample covariance matrix of the vectors classified in cluster m of state } j
```

- Step 1 : re-segment the training observation sequences into states based on the initial model by Viterbi Algorithm
- Step 2 : Reestimate the model parameters (same as initial estimation)
- Step 3: Evaluate the model score  $P(O|\lambda)$ :

  If the difference between the previous and current model scores exceeds a threshold, go back to Step 1, otherwise stop and the initial model is obtained

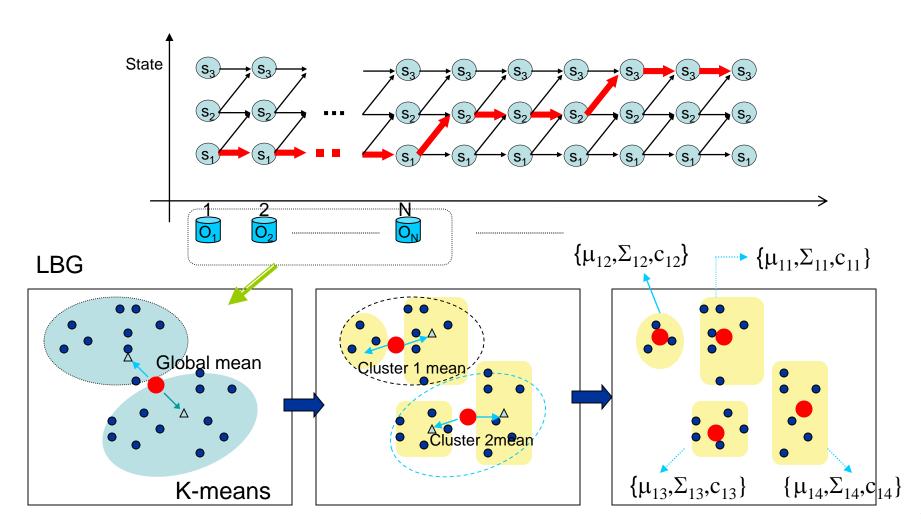
## Segmental K-Means



### **Initialization in HMM Training**

### An example for Continuous HMM

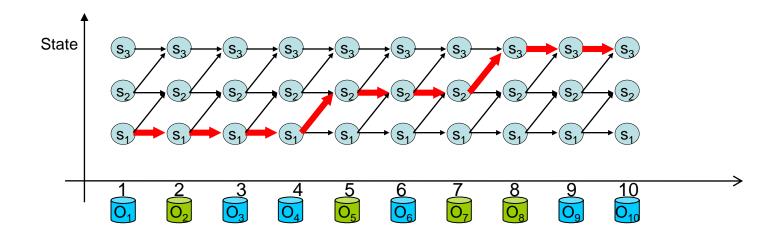
- 3 states and 4 Gaussian mixtures per state



## **Initialization in HMM Training**

### • An example for discrete HMM

- 3 states and 2 codewords



$$b_1(\mathbf{v}_1) = 3/4, b_1(\mathbf{v}_2) = 1/4$$
  
 $b_2(\mathbf{v}_1) = 1/3, b_2(\mathbf{v}_2) = 2/3$   
 $b_3(\mathbf{v}_1) = 2/3, b_3(\mathbf{v}_2) = 1/3$ 

