### 4.0 More about Hidden Markov Models

**Reference**: 1. 6.1-6.6, Rabiner and Juang

2. 4.4.1 of Huang

### Markov Model

• An example : a 3-state Markov Chain  $\lambda$ 

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

$$\mathbf{A} = \begin{bmatrix} 0.6 & 0.3 & 0.1 \\ 0.1 & 0.7 & 0.2 \\ 0.3 & 0.2 & 0.5 \end{bmatrix}$$
$$\pi = \begin{bmatrix} 0.4 & 0.5 & 0.1 \end{bmatrix}$$

ly  $0.3 \qquad 0.3 \qquad 0.3$   $0.7 \qquad \begin{array}{c} 0.3 \\ 0.2 \\ \hline 0.2 \\ \hline \end{array} \qquad \begin{array}{c} 0.5 \\ \hline \end{array}$   $0.5 \qquad \begin{array}{c} 0.5 \\ \hline \end{array}$ 

0.6

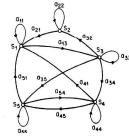
- Given a sequence of observed symbols  $O=\{CABBCABC\}$ , the **only one** corresponding state sequence is  $\{S_3S_1S_2S_3S_1S_2S_3\}$ , and the corresponding probability is  $P(O|\lambda)=P(q_0=S_3)$ 

$$P(S_1|S_3)P(S_2|S_1)P(S_2|S_2)P(S_3|S_2)P(S_1|S_3)P(S_2|S_1)P(S_3|S_2) \\ = 0.1 \times 0.3 \times 0.3 \times 0.7 \times 0.2 \times 0.3 \times 0.3 \times 0.2 = 0.00002268$$

### 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  $P(q_i = j | q_{i-1} = i, q_{i-2} = k, \dots) = P(q_i = j | q_{i-1} = i) \equiv \mathbf{a}_{ii}$
  - $\pi$  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.

### **Hidden Markov Model**

#### • HMM, an extended version of Markov Model

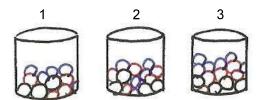
- The observation is a probabilistic function (discrete or continuous)
   of a state instead of an one-to-one correspondence of a state
- The model is a doubly embedded stochastic process with an underlying stochastic process that is not directly observable (hidden)
  - What is hidden? The State Sequence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

### • Elements of an HMM {S,A,B,π}

- 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

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# Simplified HMM



RGBGGBBGRRR....

**Hidden Markov Model** 

• An example : a 3-state discrete HMM λ

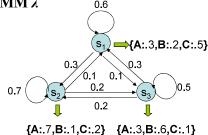
$$\mathbf{A} = \begin{bmatrix} 0.6 & 0.3 & 0.1 \\ 0.1 & 0.7 & 0.2 \\ 0.3 & 0.2 & 0.5 \end{bmatrix}$$

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

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

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

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



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

$$\begin{split} P(\overline{\mathbf{O}}|\lambda) &= \sum_{i=1}^{27} P(\overline{\mathbf{O}}, \mathbf{q}_i | \lambda) = \sum_{i=1}^{27} P(\overline{\mathbf{O}}|\mathbf{q}_i, \lambda) P(\mathbf{q}_i | \lambda) \qquad \mathbf{q}_i : \text{state sequence} \\ e.g. \text{ when } \mathbf{q}_i &= \left\{ S_2 S_2 S_3 \right\}, P(\overline{\mathbf{O}}|\mathbf{q}_i, \lambda) = P\left(\mathbf{A}|S_2\right) P\left(\mathbf{B}|S_2\right) P\left(\mathbf{C}|S_3\right) = 0.7*0.1*0.1 = 0.007 \\ P\left(\mathbf{q}_i | \lambda\right) &= P\left(\mathbf{q}_0 = S_2\right) P\left(S_2 | S_2\right) P\left(S_3 | S_2\right) = 0.5*0.7*0.2 = 0.07 \end{split}$$

## **Hidden Markov Model**

• 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  $V = \{v_1, v_2, v_3, \dots, v_M\}, v_k \in \mathbb{R}^D$
- the set of observation probability distributions  $B=\{b_j(\mathbf{v}_k)\}$  is defined as  $b_j(\mathbf{v}_k)=P(\mathbf{o}_i=\mathbf{v}_k|\mathbf{q}_i=j), \ 1 \le k \le M, \ 1 \le j \le N$ **o**,: observation at time t,  $\mathbf{q}_i$ : state at time t
- $\Rightarrow$  for state j,  $b_i(\mathbf{v}_k)$  consists of only M probability values

#### Continuous and infinite observations:

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

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

## **Hidden Markov Model**

- Three Basic Problems for HMMs Given an observation sequence  $\overline{O}$ =(o<sub>1</sub>,o<sub>2</sub>,....,o<sub>T</sub>), and an HMM  $\lambda$  =(A,B, $\pi$ )
  - Problem 1 :

How to *efficiently* compute  $P(\overline{\mathbf{O}}|\lambda)$ ?

- *⇒* Evaluation problem
- Problem 2:

How to choose an optimal state sequence  $\mathbf{q} = (q_1, q_2, \dots, q_T)$ ?

- *⇒ 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* 

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#### **Basic Problem 1 for HMM**

 $\overline{\mathbf{O}} = o_1 o_2 o_3.....o_t....o_T$  observation sequence

 $\overline{q} = q_1q_2q_3....q_t....q_T$  state sequence

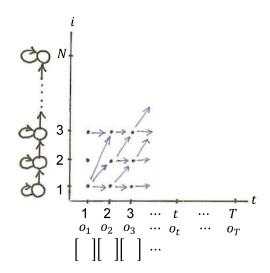
• **Problem 1:** Given  $\lambda$  and  $\overline{O}$ , find  $P(\overline{O}|\lambda) = \text{Prob[observing } \overline{O} \text{ given } \lambda]$ 

• Direct Evaluation: considering all possible state sequence  $\overline{\mathbf{q}}$ 

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

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

# Basic Problem 1



#### **Basic Problem 1 for HMM**

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

$$\alpha_{t}(i) = P(o_1 o_2 \dots o_t, q_t = i | \lambda)$$

=Prob[observing  $o_1o_2...o_t$ , state i at time  $t|\lambda$ ]

- Initialization

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

- 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

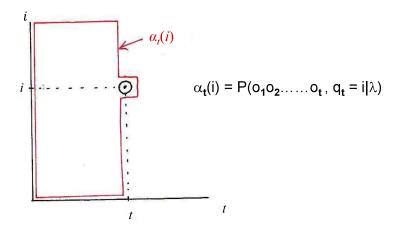
$$P(\overline{O}|\lambda) = \sum_{i=1}^{N} \alpha_{T}(i)$$

See Fig. 6.5 of Rabiner and Juang

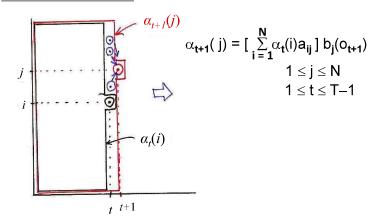
- All state sequences, regardless of how long previously, merge to the N state at each time instant t

# Basic Problem 1

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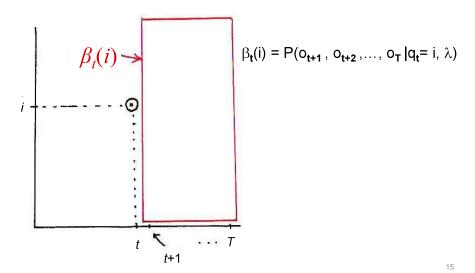


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Forward Algorithm

# Basic Problem 2



#### **Basic Problem 2 for HMM**

- Problem 2: Given  $\lambda$  and  $\overline{O} = o_1 o_2 ... o_T$ , find a best state sequence  $\overline{q} = q_1 q_2 ... q_T$
- Backward Algorithm : defining a backward variable β<sub>t</sub>(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 
$$\beta_T(i)=1,\ 1\leq i\leq N \qquad (\ \beta_{T\text{-}1}(i)=\sum\limits_{i=1}^N a_{ij}\ b_j(o_T)\ )$$

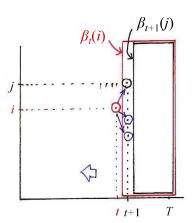
- Induction 
$$\beta_{t}(i) = \sum_{j=1}^{N} a_{ij} \ b_{j}(o_{t+1})\beta_{t+1}(j)$$
 
$$t = T-1, T-2, \dots, 2, 1, \qquad 1 \leq i \leq N$$
 See Fig. 6.6 of Rabiner and Juang

• Combining Forward/Backward Variables

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

## Basic Problem 2

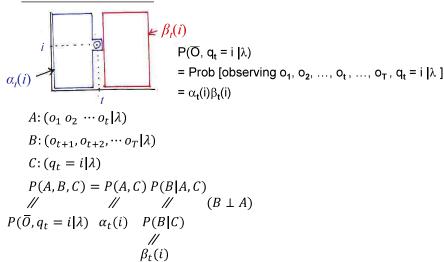
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$$\beta_{t+1}(j) \qquad \beta_{t}(i) = \sum_{j=1}^{N} a_{ij} b_{j}(o_{t+1}) \beta_{t+1}(j)$$

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

## Backward Algorithm



#### **Basic Problem 2 for HMM**

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

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

- Solution

$$q_t^* = \arg \max_{1 \le i \le N} [\gamma_t(i)], 1 \le t \le T$$

in fact

$$\begin{aligned} & \underset{q_t^*}{\text{milter}} = \arg \max_{1 \le i \le N} \left[ P(\overline{O}, q_t = i | \lambda) \right] \\ & = \arg \max_{1 \le i \le N} \left[ \alpha_t(i) \beta_t(i) \right] \end{aligned}$$

- Problem

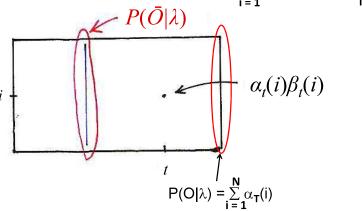
maximizing the probability at each time t individually

$$\overline{q}{}^*{=}\ q_1{}^*q_2{}^*{\dots}q_T{}^*$$
 may not be a valid sequence

(e.g. 
$$a_{q_t*q_{t+1}*} = 0$$
)

## Basic Problem 2

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



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### **Basic Problem 2 for HMM**

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

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

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

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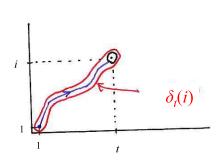
$$\delta_{t+1}(j) = \max \left[ \delta_t(i) a_{ij} \right] \bullet b_j(o_{t+1})$$

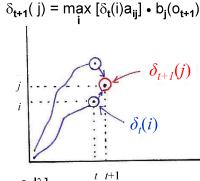
- Backtracking

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

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





$$\delta_t(i) = \max_{\textbf{q}_1,\textbf{q}_2,\dots\textbf{q}} \Pr_{\textbf{t}\text{-}1}[\textbf{q}_1,\textbf{q}_2,\dots\textbf{q}_{\textbf{t}\text{-}1},\ \textbf{q}_t = i,\ \textbf{o}_1,\textbf{o}_2,\dots,\textbf{o}_t\ \big| \boldsymbol{\lambda}]$$

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#### **Basic Problem 2 for HMM**

- Complete Procedure for Viterbi Algorithm
- Initialization

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

- Recursion

$$\delta_{t+1}(j) = \max_{1 \le i \le N} \left[ \delta_t(i) a_{ij} \right] \bullet b_j(o_{t+1})$$

$$1 \le t \le T-1$$
,  $1 \le j \le N$ 

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

$$1 \le t \le T-1$$
,  $1 \le j \le N$ 

- Termination

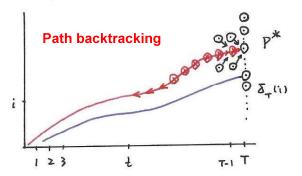
$$P^* = \max_{1 \le i \le N} [\delta_T(i)]$$

$$q_T^* = \arg \max_{1 \le i \le N} [\delta_T(i)]$$

- Path backtracking

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

# Viterbi Algorithm



#### **Basic Problem 2 for HMM**

- Application Example of Viterbi Algorithm
- Isolated word recognition

$$\lambda_0 = (A_0, B_0, \pi_0)$$

$$\boldsymbol{\lambda}_1 = (\boldsymbol{A}_1, \boldsymbol{B}_1, \boldsymbol{\pi}_1)$$

$$\lambda_{n} = (A_{n}, B_{n}, \boldsymbol{\pi}_{n})$$

observation

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

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

Basic Problem 1 Forward Algorithm Basic Problem 2 Viterbi Algorithm

(for all paths) (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.

#### **Basic Problem 3 for HMM**

- **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

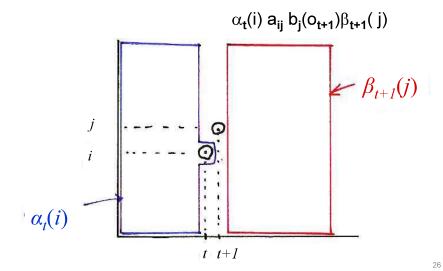
$$\begin{split} \boldsymbol{\epsilon}_{t}(\;i,j\;) &=\; P(q_{t}=i,\,q_{t+1}=j\;|\;\overline{O},\,\lambda) \\ &=\; \frac{\alpha_{t}(i)\;a_{ij}\;b_{j}(o_{t+1})\beta_{t+1}(j)}{\sum\limits_{i=1}^{N}\;\sum\limits_{j=1}^{N}\left[\alpha_{t}(i)a_{ij}\;b_{j}(o_{t+1})\beta_{t+1}(j)\right]} \\ &=\; \frac{Prob[\overline{O},\;q_{t}=i,\,q_{t+1}=j|\lambda]}{P(\overline{O}|\lam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

See Fig. 6.7 of Rabiner and Juang

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

 $\sum_{t=1}^{T-1} \epsilon_t(i,j)$  = expected number of transitions from state i to state j in  $\overline{O}$ 

# Basic Problem 3



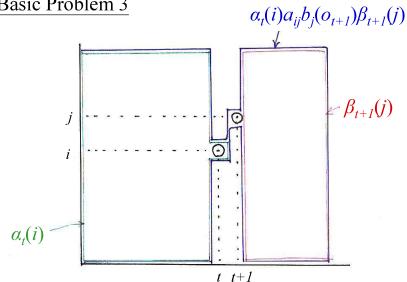
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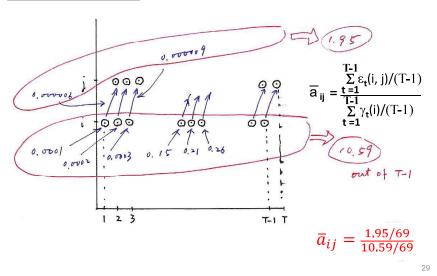
# Basic Problem 3

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

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

Basic Problem 3





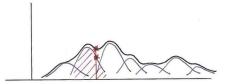
**Basic Problem 3 for HMM** 

• 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

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



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- Results

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

See Fig. 6.9 of Rabiner and Juang

**Basic Problem 3 for HMM** 

- Results

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

• Continuous Density HMM

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

N(): Multi-variate Gaussian

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

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

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

**Basic Problem 3 for HMM** 

• Continuous Density HMM

$$\begin{split} \overline{\mu}_{jk} &= \frac{\sum\limits_{t=1}^{T} \left[ \gamma_{t}(j,k) \bullet o_{t} \right]}{\sum\limits_{t=1}^{T} \gamma_{t}(j,k)} \\ \overline{U}_{jk} &= \frac{\sum\limits_{t=1}^{T} \left[ \gamma_{t}(j,k)(o_{t} - \mu_{jk}) \left( o_{t} - \mu_{jk} \right)' \right]}{\sum\limits_{t=1}^{T} \gamma_{t}(j,k)} \end{split}$$

• Iterative Procedure

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

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

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

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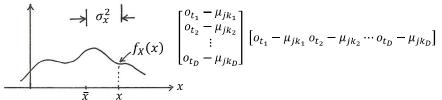
$$\frac{-}{\mu_{jk}} = \frac{\sum\limits_{t=1}^{T} \left[ \begin{array}{c} \gamma_t(j,k) \\ \sum\limits_{t=1}^{T} \gamma_t(j,k) \end{array} \right] \bullet o_t]}{\sum\limits_{t=1}^{T} \gamma_t(j,k)}$$

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

$$\overline{U}_{jk} = \frac{\sum_{t=1}^{T} [\gamma_{t}(j,k)] (o_{t} - \mu_{jk}) (o_{t} - \mu_{jk})^{t}}{\sum_{t=1}^{T} \gamma_{t}(j,k)}$$

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

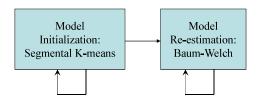
### $f_X(x)$ : prob. density function



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### **Basic Problem 3 for HMM**

- 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

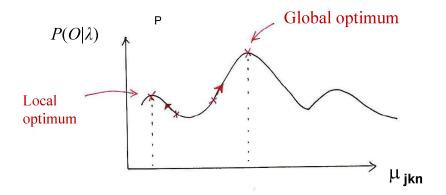


# Basic Problem 3

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

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

Basic Problem 3



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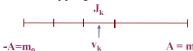
# **Vector Quantization (VQ)**

### • 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



$$S = \bigcup_{k=1}^{L} J_{k}, V = \{ v_{1}, v_{2}, ..., v_{L} \}$$

$$Q : S \to V$$

$$Q(x[n]) = v_{k} \text{ if } x[n] \in J_{k}$$

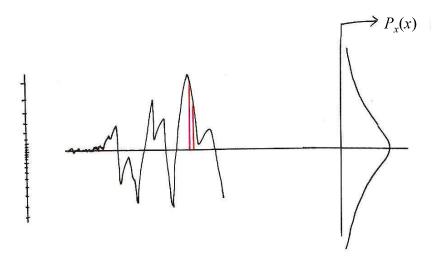
$$L = 2^{R}$$

Each v<sub>k</sub> represented by an R-bit pattern

- -Quantization characteristics (codebook)  $\{J_1, J_2, ..., J_L\}$  and  $\{v_1, v_2, ..., v_L\}$ 
  - designed considering at least
  - 1. error sensitivity
  - 2. probability distribution of x[n]

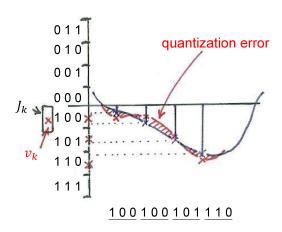
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# **Vector Quantization**



# **Vector Quantization**

Scalar Quantization: Pulse Coded Modulation (PCM)



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# **Vector Quantization (VQ)**

### 2-dim Vector Quantization (VQ)

Example:

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

$$S = {\overline{X}_n = (x[n], x[n+1]); |x[n]| < A, |x[n+1]| < A}$$

•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

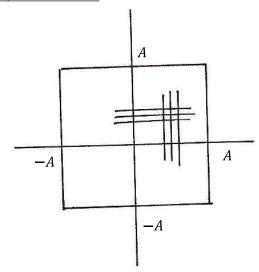
 $Q(\overline{x}_n) = \overline{v}_k \text{ if } \overline{x}_n \in J_k$  $J_k = 2^R$ 

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

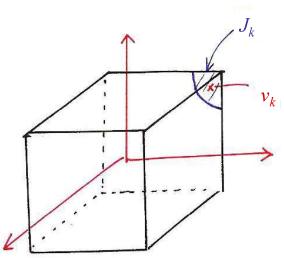
- Considerations
  - 1.error sensitivity may depend on x[n], x[n+1] jointly
  - 2.distribution of x[n], x[n+1] may be correlated statistically
  - 3.more flexible choice of J<sub>k</sub>
- Quantization Characteristics (codebook)

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

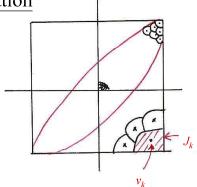
# **Vector Quantization**



# **Vector Quantization**



# **Vector Quantization**



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

$$1024=2^{10}$$

# **Vector Quantization (VQ)**

### N-dim Vector Quantization

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

# Codebook Trained by a Large

### Training Set

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

$$\begin{array}{c} d(\;\overline{x},\overline{y}\;):S{\times}S \to R^{+}\,(\text{non-negative}\\ \text{real numbers}) \end{array}$$

-desired properties

$$d(\overline{x}, \overline{y}) \ge 0$$

$$d(\overline{x}, \overline{x}) = 0$$

$$d(\overline{x}, \overline{y}) = d(\overline{y}, \overline{x})$$

$$d(\overline{x}, \overline{y}) + d(\overline{y}, \overline{z}) \ge d(\overline{x}, \overline{z})$$

examples:

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

$$d(\overline{x}, \overline{y}) = \sum_{i=1}^{n} |x_i - y_i|$$

$$d(\overline{x}, \overline{y}) = (\overline{x} - \overline{y})^t \sum_{i=1}^{t} (\overline{x} - \overline{y})$$

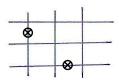
Mahalanobis Distance

 $\Sigma$ : Co-variance Matrix

## **Distance Measures**

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

city block distance

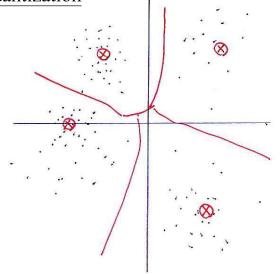


$$d(\bar{x}, \bar{y}) = (\bar{x} - \bar{y})^t \Sigma^{-1} (\bar{x} - \bar{y})$$
 Mahalanobis distance

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

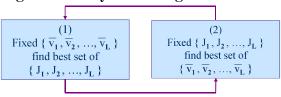
$$\sum = \begin{bmatrix} \sigma_1^2 & \cdots & 0 \\ \vdots & \sigma_2^2 & \vdots \\ 0 & \cdots & \sigma_n^2 \end{bmatrix}, d(\bar{x}, \bar{y}) = \sum_i \frac{(x_i - y_i)^2}{\sigma_i^2}$$

**Vector Quantization** 



# **Vector Quantization (VQ)**

• K-Means Algorithm/Lloyd-Max Algorithm



$$(1) J_{\mathbf{k}} = \{ \overline{x} \mid d(\overline{x}, \overline{v}_{\mathbf{k}}) < d(\overline{x}, \overline{v}_{\mathbf{j}}), j \neq k \}$$

$$\rightarrow D = \sum_{\text{all } x} d(\overline{x}, Q(\overline{x})) = \min$$

nearest neighbor condition

(2) For each k  $\rightarrow D_{\mathbf{k}} = \sum d(\overline{x}, \overline{v}_{\mathbf{k}}) = \min$ centroid condition

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(3) Convergence condition

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

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

• Iterative Procedure to Obtain Codebook from a Large Training Set

# **Vector Quantization (VQ)**

- 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

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

- step 2: Splitting.

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

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

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

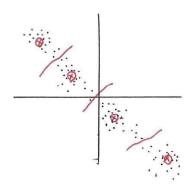
$$v_k - v_k(1+\varepsilon)$$
 $v_k = v_k(1-\varepsilon)$ 

$$\mathbf{v}_{k}^{(1)} = \mathbf{v}_{k}$$
 $\mathbf{v}_{k}^{(2)}$ : the vector  $\mathbf{v}_{k}$ 

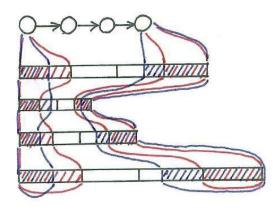
 $\frac{1}{V_k}$ (2): the vector most far apart

- 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



# Segmental K-Means



# **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

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

- •For continuous density HMM (M Gaussian mixtures per state)
  - ⇒ cluster the observation vectors within each state j into a set of M clusters (e.g. with vector quantiziation)

 $c_{jm}$  = number of vectors classified in cluster m of state j

divided by number of vectors in state j

 $\mu_{\text{jm}} = \text{sample}\,\text{mean}\,\text{of}\,\,\text{the}\,\,\text{vectors}\,\text{classified}\,\text{in}\,\,\text{cluster}\,\text{m}\,\,\text{of}\,\,\text{state}\,\,j$ 

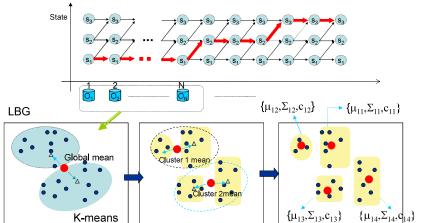
 $\sum_{im}$  = sample covariance matrix of the vectors classified in cluster m of state j

- 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/λ):
   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

# **Initialization in HMM Training**

### • An example for Continuous HMM

- 3 states and 4 Gaussian mixtures per state



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# **Initialization in HMM Training**

## • An example for discrete HMM

- 3 states and 2 codewords

