Expectation Maximisation (EM)

CS 486/686: Introduction to Artificial Intelligence University of Waterloo

Incomplete Data

So far we have seen problems where

- Values of all attributes are known
- Learning is relatively easy

Many real-world problems have hidden variables

- Incomplete data
- Missing attribute values

Maximum Likelihood Learning

Learning of Bayes nets parameters

• $\Theta_{V=true, Par(V)=x} = P(V=true|Par(V)=x)$

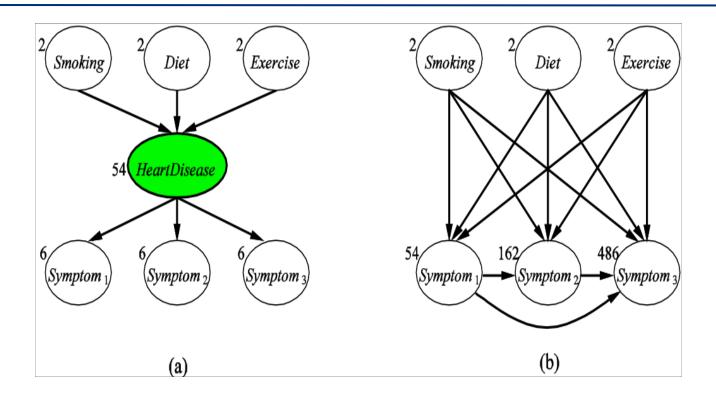
Assumes all attributes have values

- What if some values are missing?

Naïve Solutions

- Ignore examples with missing attribute values
 - What if all examples have missing attribute values?
- Ignore hidden variables
 - Model might become much more complex

Hidden Variables Heart disease example



a) Uses a Hidden Variable, simpler (fewer CPT parameters)b) No Hidden Variable, complex (many CPT parameters)

"Direct" ML

Maximize likelihood directly where E are the evidence variables and Z are the hidden variables

$$h_{ML} = \arg \max_{h} P(E|h)$$

= $\arg \max_{h} \sum_{Z} P(E, Z|h)$
= $\arg \max_{h} \sum_{Z} \prod_{i} CPT(V_i)$
= $\arg \max_{h} \log \sum_{z} \prod_{i} CPT(V_i)$

Expectation-Maximization (EM)

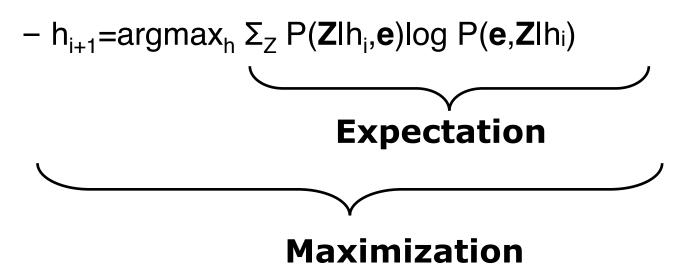
If we knew the missing values computing h_{ML} is trivial

- Guess h_{ML}
- Iterate
 - Expectation: based on h_{ML} compute expectation of (missing) values
 - Maximization: based on expected (missing) values compute new h_{ML}

Expectation-Maximization (EM)

Formally

- Approximate maximum likelihood
- Iteratively compute:



EM Derivation

$$\log P(\mathbf{e}|h) = \log \left[\frac{P(\mathbf{e}, \mathbf{Z}|h)}{P(\mathbf{Z}|\mathbf{e}, h)}\right]$$

= log P(\mathbf{e}, \mathbf{Z}|h) - log P(\mathbf{Z}|\mathbf{e}, h)
= $\sum_{Z} P(\mathbf{Z}|\mathbf{e}, h) \log P(\mathbf{e}, \mathbf{Z}|h) - \sum_{Z} P(\mathbf{Z}|\mathbf{e}, h) \log P(\mathbf{Z}|\mathbf{e}, h)$
\ge \sum_{Z} P(\mathbf{Z}|\mathbf{e}, h) \log P(\mathbf{e}, \mathbf{Z}|h)

EM finds a local maxima of $\sum_{Z} P(\mathbf{Z}|\mathbf{e}, h) \log P(\mathbf{e}, \mathbf{Z}|h)$

which is a lower bound of $\log P(\mathbf{e}|h)$

EM

Log inside can linearize the product $h_{i+1} = \arg \max_{h} \sum_{\mathbf{Z}} P(\mathbf{Z}|h, \mathbf{e}) \log P(\mathbf{e}, \mathbf{Z}|h)$ $= \arg \max_{h} \sum_{\mathbf{Z}} P(\mathbf{Z}|h, \mathbf{e}) \log \prod_{j} \mathrm{CPT}_{j}$ $= \arg \max_{h} \sum_{Z} P(\mathbf{Z}|h, \mathbf{e}) \sum_{j} \log \mathrm{CPT}_{j}$ Monotonic improvement of likelihood $P(\mathbf{e}|h_{i+1}) \geq P(\mathbf{e}|h_i)$

- Assume we have two coins, A and B
- The probability of getting heads with A is θ_A
- The probability of getting heads with B is θ_B
- We want to find θ_A and θ_B by performing a number of trials

Example from S. Zafeiriohu, Advanced Statistical Machine Learning, Imperial College

Coin A and Coin B

- нтттннтнтн
- ннннтнннн
- нтннннтнн
- нтнтттннтт
- тнннтнннтн

Coin A	Coin B
	5 H, 5 T
9 H, 1 T	
8 H, 2 T	
	4 H, 6 T
7 H, 3 T	
24 H, 6 T	9 H, 11 T

Coin A	Coin B
	5 H, 5 T
9 H, 1 T	
8 H, 2 T	
	4 H, 6 T
7 H, 3 T	
24 H, 6 T	9 H, 11 T

$$\theta_A = \frac{24}{24+6} = 0.8$$
$$\theta_B = \frac{9}{9+11} = 0.45$$

Now assume we do not know which coin was used in which trial (hidden variable)

- нтттннтнтн
- ннннтннннн
- нтннннтнн
- нтнтттннтт
- тнннтнннтн

Initialization: $\theta_A^0 = 0.60$ $\theta_B^0 = 0.50$

E Step: Compute the Expected counts of Heads and Tails

Trial 1: H T T T H H T H T H

$$P(A|\text{Trial 1}) = \frac{P(\text{Trial 1}|A)P(A)}{\sum_{i \in \{A,B\}} P(\text{Trial 1}|i)P(i)} = 0.45$$

$$P(B|\text{Trial 1}) = \frac{P(\text{Trial 1}|B)P(B)}{\sum_{i \in \{A,B\}} P(\text{Trial 1}|i)P(i)} = 0.55$$

$$Coin A \quad Coin B \quad 2.2 \text{ H}, \quad 2.8 \text{ H},$$

- HTTTHHTHTH
 (0.55 A, 0.45 B)
- **HHHHTHHHH** (0.80 A, 0.20 B)
- **HTHHHHHHHH** (0.73 A, 0.27 A)
- **HTHTTTHHTT** (0.35 A, 0.65 B)
- **THHHTHHHTH** (0.65 A, 0.35 B)

Coin A	Coin B
2.2H, 2.2T	2.8H, 2.8T
7.2H, 0.8T	1.8H, 0.2T
5.9H, 1.5T	2.1H, 0.5T
1.4H, 2.1T	2.6H, 3.9T
4.5H, 1.9T	2.5H, 1.1T
21.3H, 8.6T	11.7H, 8.4T

M Step: Compute parameters based on expected counts

Coin A	Coin B
2.2H, 2.2T	2.8H, 2.8T
7.2H, 0.8T	1.8H, 0.2T
5.9H, 1.5T	2.1H, 0.5T
1.4H, 2.1T	2.6H, 3.9T
4.5H, 1.9T	2.5H, 1.1T
21.3H, 8.6T	11.7H, 8.4T

$$\begin{split} \theta^1_A &= \frac{21.3}{21.3+8.6} = 0.71 \\ \theta^1_B &= \frac{11.7}{11.7+8.4} = 0.58 \end{split}$$

Repeat

$$heta_A^{10} = 0.80 \ heta_B^{10} = 0.52$$

EM: k-means Algorithm

Input

- Set of examples, E
- Input features X₁,...,X_n
- val(e,X)=value of feature j for example e
- k classes

Output

- Function *class*:E-> {1,...,k} where *class(e)=i* means example e belongs to class i
- Function *pval* where *pval(i,X_j)* is the predicted value of feature X_j for each example in class i

k-means Algorithm

Sum-of-squares error for class i and pval is

$$\sum_{e \in E} \sum_{j=1}^{n} (\operatorname{pval}(\operatorname{class}(e), X_j) - \operatorname{val}(e, X_j))^2$$

 Goal: Final *class* and *pval* that minimizes sum-of-squares error.

Minimizing the error

$$\sum_{e \in E} \sum_{j=1}^{n} (\operatorname{pval}(\operatorname{class}(e), X_j) - \operatorname{val}(e, X_j))^2$$

- Given *class*, the *pval* that minimizes sum-of-square error is the mean value for that class
- Given *pval*, each example can be assigned to the *class* that minimizes the error for that example

k-means Algorithm

- Randomly assign the examples to classes
- Repeat the following two steps until E step does not change the assignment of any example
 - **M**: For each class i and feature Xj

 \boldsymbol{n}

$$\operatorname{pval}(i, X_j) = \frac{\sum_{e:\operatorname{class}(e)=i} \operatorname{val}(e, X_j)}{|\{e: \operatorname{class}(e)=i\}|}$$

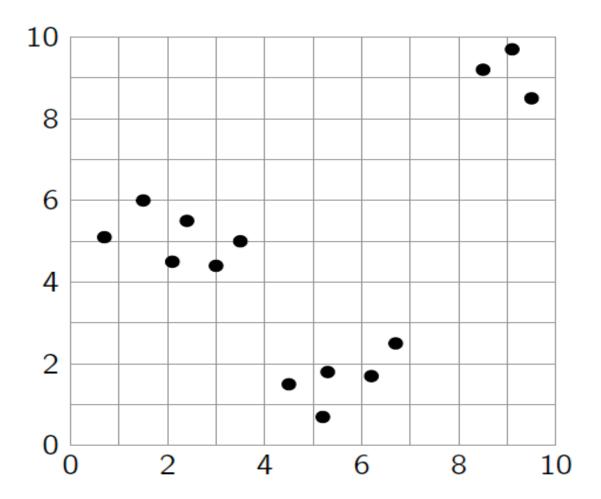
- E: For each example e, assign e to the class that minimizes

$$\sum_{j=1}^{n} (\operatorname{pval}(\operatorname{class}(e), X_j) - \operatorname{val}(e, X_j))^2$$

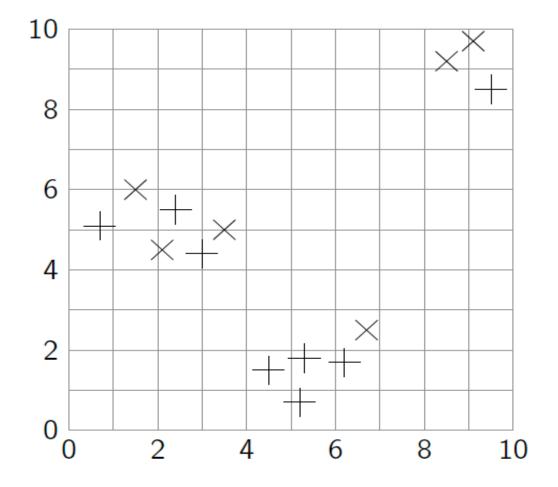
k-means Example

- Data set: (X,Y) pairs
 - $\begin{array}{l} & (0.7,5.1) \ (1.5,6), \ (2.1, \ 4.5), \ (2.4, \ 5.5), \ (3, \ 4.4), \\ & (3.5, \ 5), \ (4.5, \ 1.5), \ (5.2, \ 0.7), \ (5.3, \ 1.8), \ (6.2, \\ & 1.7), \ (6.7, \ 2.5), \\ & (8.5, \ 9.2), \ (9.1, \ 9.7), \ (9.5, \\ & 8.5) \end{array}$

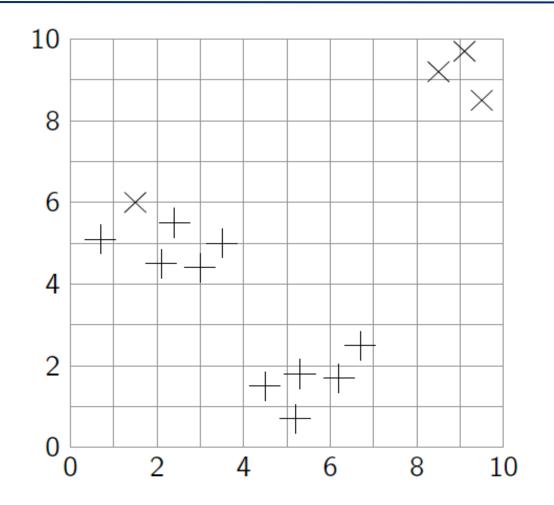
Example Data



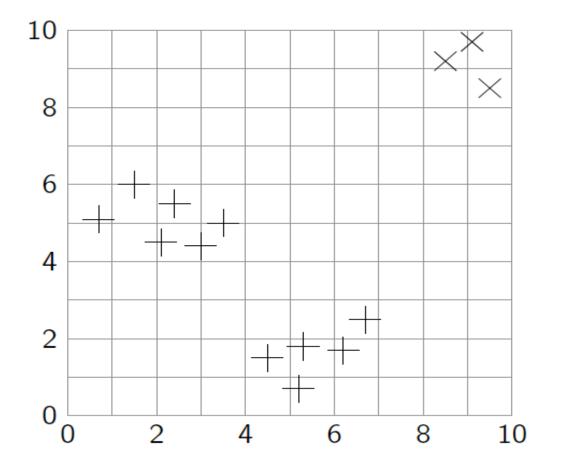
Random Assignment to Classes



Assign Each Example to Closest Mean



Reassign each example



Properties of k-means

- An assignment is stable if both M step and E step do not change the assignment
 - Algorithm will eventually converge to a stable local minimum
 - No guarantee that it will converge to a global minimum
- Increasing k can always decrease error until k is the number of different examples