<u>Week : 9</u>

Part one

Anomaly detection:

Suppose we have some training sets we want to select which of them are faulty and which are not. In that case we need to make such graph as below using their features and predict new test sets.

Anomaly detection example

Aircraft engine features:

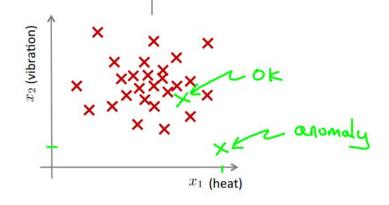
 $\rightarrow x_1$ = heat generated

 \rightarrow x_2 = vibration intensity

...

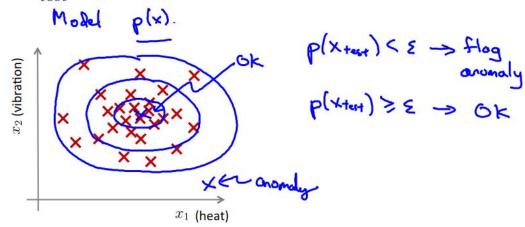
Dataset: $\{x^{(1)}, x^{(2)}, \dots, x^{(m)}\}$

New engine: x_{test}



Density estimation

- → Dataset: $\{x^{(1)}, x^{(2)}, \dots, x^{(m)}\}$
- \rightarrow Is x_{test} anomalous?



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Anomaly detection example

- → Fraud detection:
 - $\rightarrow x^{(i)}$ = features of user i 's activities
 - \rightarrow Model p(x) from data.
 - → Identify unusual users by checking which have $p(x) < \varepsilon$
- → Manufacturing
- Monitoring computers in a data center.
 - $\rightarrow x^{(i)}$ = features of machine i

 x_1 = memory use, x_2 = number of disk accesses/sec,

 x_3 = CPU load, x_4 = CPU load/network traffic.

... p(x) < &

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Gaussian Distribution:

Gaussian distribution is a probability distribution by which we can simulate the density of our training sets. There we used normal gaussian distribution where two perimeter are used: Mewe and sigma. There mewe is the center of the density and sigma is the width of the curve.

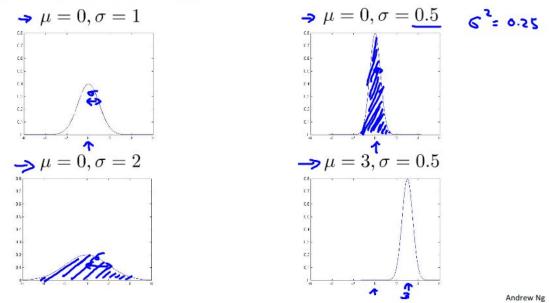
Gaussian (Normal) distribution

Say $x \in \mathbb{R}$. If x is a distributed Gaussian with mean μ , variance $\underline{\sigma}^2$.

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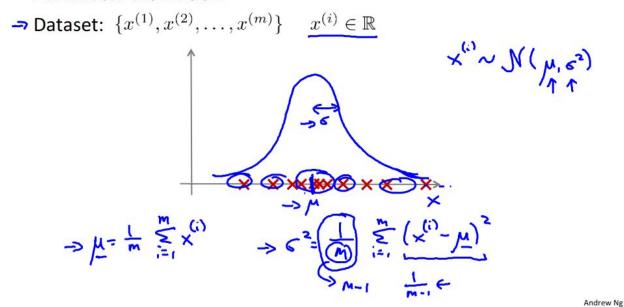
From the image below we can see that as sigma grows bigger, the width of the curve becomes thicker.

Gaussian distribution example



If we have a graph with some random training sets and want to find mewe and sigma, can apply the formula of mewe and sigma.

Parameter estimation



Anomaly detection algorithm:

In this part we will learn an algorithm which will help us to separate test set anomaly and non-anomaly using gaussian distribution.

Density estimation

Training set:
$$\{x^{(1)}, \dots, x^{(m)}\}$$

Each example is $\underline{x} \in \mathbb{R}^n$

$$\Rightarrow P(x)$$

$$= P(x_1; \mu_1, e_1^2) P(x_2; \mu_2, e_2^2) P(x_3; \mu_3, e_2^2) \cdots P(x_n; \mu_n, e_n^2)$$

$$= P(x_1; \mu_1, e_1^2) P(x_2; \mu_2, e_2^2) P(x_3; \mu_3, e_2^2) \cdots P(x_n; \mu_n, e_n^2)$$

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$$= P(x_1; \mu_1, e_$$

Anomaly detection algorithm

- \rightarrow 1. Choose features $\underline{x_i}$ that you think might be indicative of anomalous examples.
- $\Rightarrow \text{ 2. Fit parameters } \mu_1, \dots, \mu_n, \sigma_1^2, \dots, \sigma_n^2$ $\Rightarrow \underbrace{\mu_j = \frac{1}{m} \sum_{i=1}^m x_j^{(i)}}_{\sigma_j^2} \qquad \text{p(x_j; \mu_j, c_j^2)}$ $\Rightarrow \sigma_j^2 = \frac{1}{m} \sum_{i=1}^m (x_j^{(i)} \mu_j)^2 \qquad \text{for } \mu_j = \frac{1}{m} \underbrace{\sum_{i=1}^m x_j^{(i)}}_{\sigma_j^2} = \frac{1}{m$
- $\Rightarrow \text{ 3. Given new example } x \text{ , compute } \underline{p(x)} \text{ : } \\ \underline{p(x)} = \prod_{j=1}^n \underline{p(x_j; \mu_j, \sigma_j^2)} = \prod_{j=1}^n \frac{1}{\sqrt{2\pi}\sigma_j} \exp{(-\frac{(x_j \mu_j)^2}{2\sigma_j^2})} \\ \text{ Anomaly if } \underline{p(x)} < \varepsilon$

Anomaly detection example p(x) $\frac{\varepsilon = 0.02}{p(x_{test}^{(1)})} = 0.0426 \geqslant \varepsilon$ $p(x_{test}^{(2)}) = 0.0021 < \varepsilon$

evaluating an anomaly detection **Developing** and system:

In order to develop an anomaly evaluation system we need to follow the same process of supervised learning algorithms. First we need some training sets with labels to train our algorithm and then we will have a graph from which we can plot a normal gaussian distribution graph. And this gaussian distribution graph will help us to say if a test set is anomaly or not!

The importance of real-number evaluation

When developing a learning algorithm (choosing features, etc.), making decisions is much easier if we have a way of evaluating our learning algorithm.

- → Assume we have some labeled data, of anomalous and nonanomalous examples. (y = 0 if normal, y = 1 if anomalous).
- \rightarrow Training set: $\underline{x^{(1)}, x^{(2)}, \dots, x^{(m)}}$ (assume normal examples/not

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If we have 10k training sets with labels then we can divide them into 60-20-20 for training sets, cross-validation sets and test sets consecutively. Thus we can develop our algorithm.

Aircraft engines motivating example

- → 10000 good (normal) engines
- Training set: 6000 good engines (y = 0), 10 anomalous (y = 1)Test: 2000 good engines (y = 0), 10 anomalous (y = 1)

Alternative:

Training set: 6000 good engines

- CV: 4000 good engines (y=0), 10 anomalous (y=1) Test: 4000 good engines (y=0) 10 anomalous (y=1)

For evaluating, we need to fit the unlabeled training sets into gaussian distribution and then evaluate our cross validation or test sets.

Algorithm evaluation

- \Rightarrow Fit model $\underline{p(x)}$ on training set $\{\underline{x^{(1)},\ldots,x^{(m)}}\}$ $(x_{\text{test}}^{(i)},y_{\text{test}}^{(i)})$ \Rightarrow On a cross validation/test example \underline{x} , predict

$$y = \begin{cases} \frac{1}{0} & \text{if } p(x) < \bigcirc \text{(anomaly)} \\ 0 & \text{if } p(x) \ge \bigcirc \text{(normal)} \end{cases}$$

Possible evaluation metrics:

- -> True positive, false positive, false negative, true negative
- → Precision/Recall

→ - F₁-score <</p>

Test set

CV

Can also use cross validation set to choose parameter ε

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Anomaly vs supervised learning:

First few steps of anomaly and supervised learning are the same. So it may be confusing that when should we use anomaly detection and when supervised learning? It that case we need to evaluate some fact:

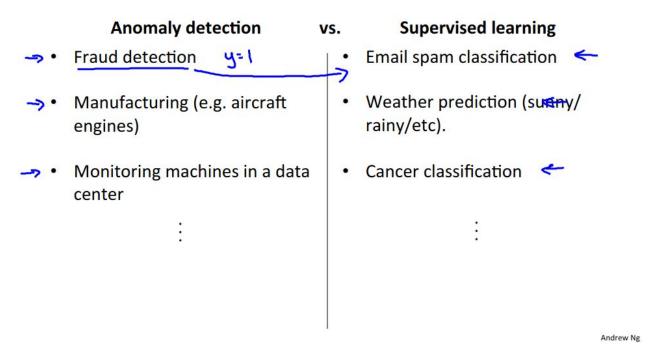
Anomaly detection

- Very small number of positive examples (y = 1). (0-20 is common).
- \Rightarrow Large number of negative $(\underline{y} = 0)$ examples.
- > Many different "types" of anomalies. Hard for any algorithm to learn from positive examples what the anomalies look like;
- future anomalies may look nothing like any of the anomalous examples we've seen so far.

Supervised learning VS.

Large number of positive and < negative examples.

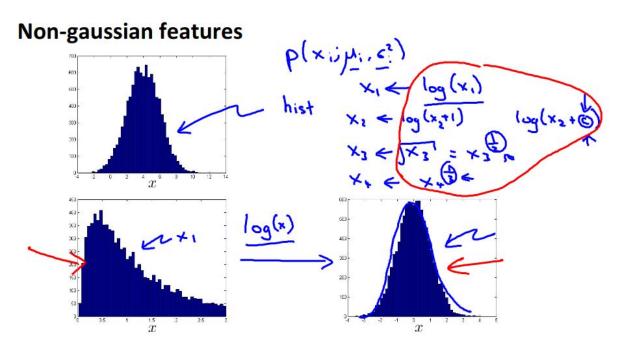
Enough positive examples for < algorithm to get a sense of what positive examples are like, future <positive examples likely to be similar to ones in training set.



Choosing what feature to use:

Sometimes it may be necessary that we need to choose the right features or debug some feature. In that case we need to follow some technique:

1. If our datasets don't give a bowl shape graph then we need to customize our features.



2. Sometimes we need to choose a new feature to evaluate new test case:

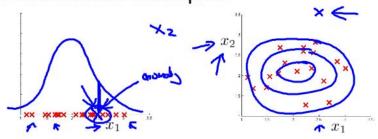
Error analysis for anomaly detection

Want p(x) large for normal examples x .

p(x) small for anomalous examples x.

Most common problem:

 $p(x)\$ is comparable (say, both large) for normal and anomalous examples



3. If you think that one or more features become faulty then you need to customize them with new features. In the picture below, we get x3,x4 linearly high. It that case, to detect which feature is faulty we customize them with x5 or x6 feature.

Monitoring computers in a data center

Choose features that might take on unusually large or small values in the event of an anomaly.

 \rightarrow x_1 = memory use of computer

 $\rightarrow x_2$ = number of disk accesses/sec

 $\rightarrow x_3 = CPU load <$

 $\rightarrow x_4$ = network traffic \leftarrow

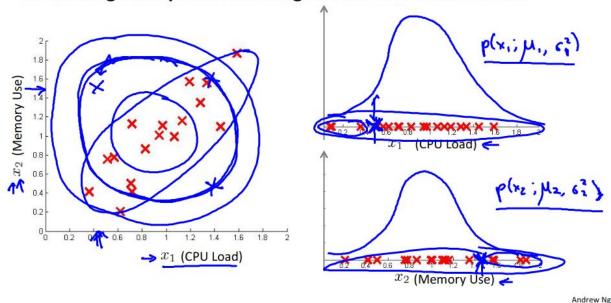
Xs = CPU load network tobic

Multivariate Gaussian Distribution:

Sometimes normal gaussian distribution may mark some test cases as non anomaly cases though they are not. It happens because of doing multiplication. In that case, we

can use multivariate gaussian distribution to get more perfect result.

Motivating example: Monitoring machines in a data center



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Multivariate Gaussian (Normal) distribution

 $x \in \mathbb{R}^n$. Don't model $p(x_1), p(x_2), \ldots$, etc. separately. Model p(x) all in one go.

Parameters: $\mu \in \mathbb{R}^n$, $\Sigma \in \mathbb{R}^{n \times n}$ (covariance matrix)

$$\frac{1}{(2\pi)^{n/2}} = \exp(-\frac{1}{2}(x-\mu)^{T} \mathcal{E}^{-1}(x-\mu))$$

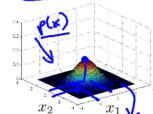
$$|\mathcal{E}| = \det(\sin n \alpha t) = \int_{0}^{\infty} \det(\sin n \alpha t)$$

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Here are some graphs of multivariate gaussian distribution. As we notice, the graphs change with the values of mewe and sigma.

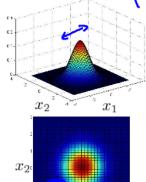
Multivariate Gaussian (Normal) examples

$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}$$



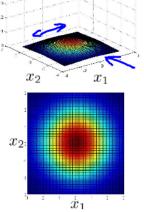
$$x_2$$
 x_1
 x_2
 x_1
 x_2

$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 0.6 & 0 \\ 0 & 0.6 \end{bmatrix}$$



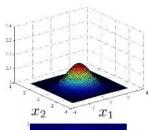
$$x_2$$
 x_1 x_2

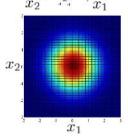
$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 2 & 0 \\ 0 & 2 \end{bmatrix}$$



Multivariate Gaussian (Normal) examples $\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \qquad \mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 0 & 0 \\ 0 & 1 \end{bmatrix} \qquad \mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 0 & 0 \\ 0 & 1 \end{bmatrix}$

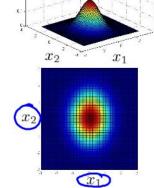
$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}$$



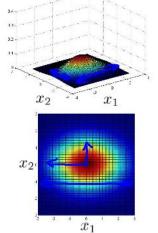


$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 0 & 0 \\ 0 & 1 \end{bmatrix}$$

 \dot{x}_1

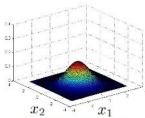


$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 2 & 0 \\ 0 & 1 \end{bmatrix}$$



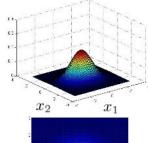
Multivariate Gaussian (Normal) examples

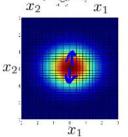
$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}$$



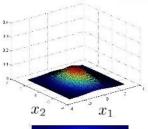
$$x_2$$
 x_1 x_2 x_1 x_2 x_1

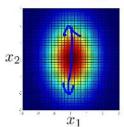
$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \qquad \qquad \mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 0.6 \end{bmatrix}$$





$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 2 \end{bmatrix}$$

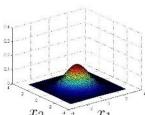


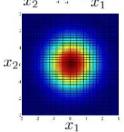


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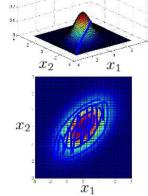
Multivariate Gaussian (Normal) examples

$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}$$

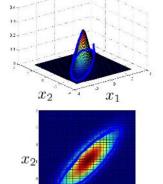




$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0.5 \\ 0.5 & 1 \end{bmatrix}$$



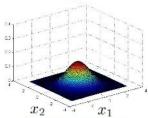
$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \qquad \qquad \mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0.5 \\ 0.5 & 1 \end{bmatrix} \qquad \qquad \mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0.8 \\ 0 & 1 \end{bmatrix}$$



 \mathring{x}_1

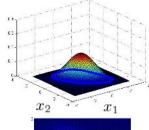
Multivariate Gaussian (Normal) examples

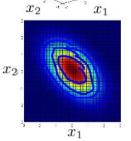
$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \underbrace{\Sigma}_{ } = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}$$



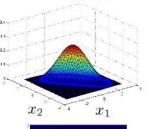
$$x_2$$
 x_1 x_2 x_1 x_2 x_1

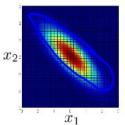
$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0.5 \\ 0.5 & 1 \end{bmatrix}$$





$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \qquad \qquad \mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & -0.5 \\ 0 & 1 \end{bmatrix} \qquad \qquad \mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & -0.8 \\ -0.8 & 1 \end{bmatrix}$$

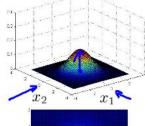


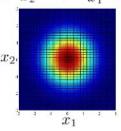


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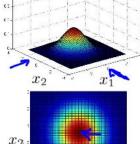
Multivariate Gaussian (Normal) examples

$$\mu = \begin{bmatrix} 0 \\ 0 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}$$



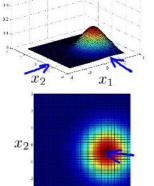


$$\mu = \begin{bmatrix} 0 \\ 0.5 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}$$



$$x_2$$

$$\mu = \begin{bmatrix} 1.5 \\ -0.5 \end{bmatrix} \Sigma = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}$$



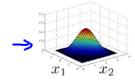
 $\dot{x_1}$

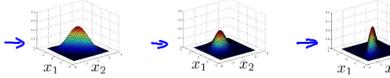
Anomaly direction using multivariate gaussian distribution:

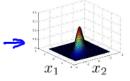
Multivariate Gaussian (Normal) distribution

Parameters μ, Σ

$$p(x;\mu,\Sigma) = \frac{1}{(2\pi)^{\frac{n}{2}}|\Sigma|^{\frac{1}{2}}} \exp\left(-\frac{1}{2}(x-\mu)^T \Sigma^{-1}(x-\mu)\right)$$







Parameter fitting:

Given training set $\{x^{(1)}, x^{(2)}, \dots, x^{(m)}\}$ \leftarrow

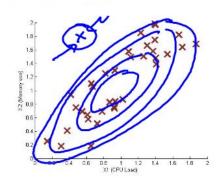
$$\boxed{\mu} = \frac{1}{m} \sum_{i=1}^{m} x^{(i)} \quad \boxed{\Sigma} = \frac{1}{m} \sum_{i=1}^{m} (x^{(i)} - \mu)(x^{(i)} - \mu)^{T}$$

Anomaly detection with the multivariate Gaussian

1. Fit model $\underline{p(x)}$ by setting

$$\mu = \frac{1}{m} \sum_{i=1}^{m} x^{(i)}$$

$$\Sigma = \frac{1}{m} \sum_{i=1}^{m} (x^{(i)} - \mu)(x^{(i)} - \mu)^{T}$$



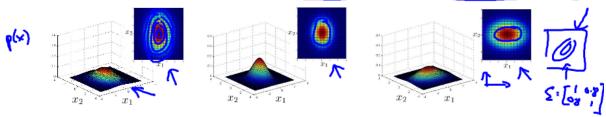
2. Given a new example x, compute

$$p(x) = \frac{1}{(2\pi)^{\frac{n}{2}} |\Sigma|^{\frac{1}{2}}} \exp\left(-\frac{1}{2}(x-\mu)^T \Sigma^{-1}(x-\mu)\right)$$

Flag an anomaly if $p(x) < \varepsilon$

Relationship to original model

Original model: $p(x) = p(x_1; \mu_1(\sigma_1^2) \times p(x_2; \mu_2, \sigma_2^2) \times \cdots \times p(x_n; \mu_n, \sigma_n^2)$



Corresponds to multivariate Gaussian

$$\Rightarrow p(x; \mu, \Sigma) = \frac{1}{(2\pi)^{\frac{n}{2}} |\Sigma|^{\frac{1}{2}}} \exp\left(-\frac{1}{2}(x-\mu)^T \Sigma^{-1}(x-\mu)\right)$$

where



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Original model

$$p(x_1; \mu_1, \sigma_1^2) \times \cdots \times p(x_n; \mu_n, \sigma_n^2)$$

Manually create features to capture anomalies where x_1, x_2 take unusual combinations of values.

(alternatively, scales better to large n=10,000, n=10,000

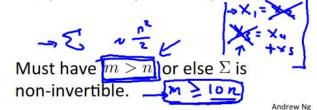
OK even if m (training set size) is small

vs. 🤝 Multivariate Gaussian

$$p(x;\mu,\Sigma) = \frac{1}{(2\pi)^{\frac{n}{2}}|\Sigma|^{\frac{1}{2}}} \exp\left(-\frac{1}{2}(x-\mu)^T \sum_{i=1}^{n} x - \mu)\right)$$

 Automatically captures correlations between features

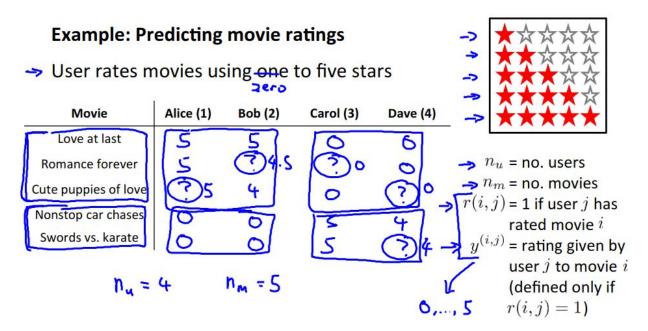
Computationally more expensive



Part two

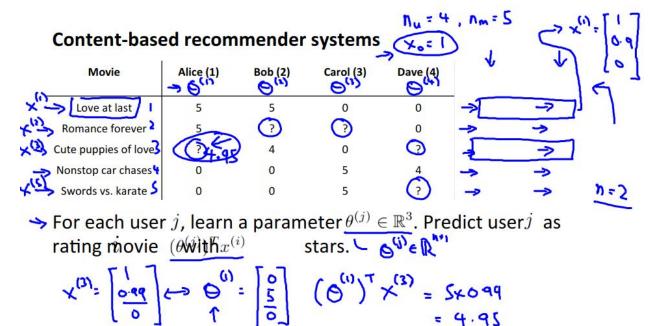
Recommender system:

Online tvs (e.g : netflix, amazon, HVO) use this system to suggest movies or series to their users. In this system they let their users rate movies or series from zero to five stars and then predict their want and choice.



Content based recommendation:

Previously we saw that users rate movies from zero to five stars. But they don't rate all of the movies. In this section we want to make a prediction by which we can say how much a user would rate an unrated movie based on their other movie rating.



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The prediction is likely to linear regression.

Problem formulation

- $\rightarrow r(i,j) = 1$ if user j has rated movie i (0 otherwise)
- \rightarrow $y^{(i,j)} = \text{rating by user } j \text{ on movie } i \text{ (if defined)}$
- \rightarrow $\theta^{(j)}$ = parameter vector for user j
- $\Rightarrow x^{(i)}$ = feature vector for movie i
- \rightarrow For user j, movie i, predicted rating: $\underline{(\theta^{(j)})^T(x^{(i)})}$
- $\rightarrow \underline{m^{(j)}}$ = no. of movies rated by user jTo learn $\underline{\theta^{(j)}}$:

$$\sum_{(i,j)}^{(i)} \frac{1}{2^{j}} \sum_{(i,i,j)=1}^{(i)} \frac{((Q_{(i)})_{i}(x_{(i)}) - Q_{(i,j)})_{j}}{((Q_{(i)})_{i}(x_{(i)}) - Q_{(i,j)})_{j}} + \frac{1}{2^{j}} \sum_{(i,j)=1}^{(i)} \frac{1}{2^{j}} (Q_{(i)}^{(i)})_{j}$$

We want to predict for all users as we calculate for theta 1,2,3,...

Optimization objective:

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$$\min_{\theta^{(1)},...,\theta^{(n_u)}} \frac{1}{2} \sum_{j=1}^{n_u} \sum_{i:r(i,j)=1} \left((\theta^{(j)})^T x^{(i)} - y^{(i,j)} \right)^2 + \frac{\lambda}{2} \sum_{j=1}^{n_u} \sum_{k=1}^{n} (\theta^{(j)}_k)^2$$

$$\mathsf{Gradient \ descent \ update:}$$

$$\theta^{(j)}_k := \theta^{(j)}_k - \alpha \sum_{i:r(i,j)=1} ((\theta^{(j)})^T x^{(i)} - y^{(i,j)}) x^{(i)}_k \ (\text{for } k = 0)$$

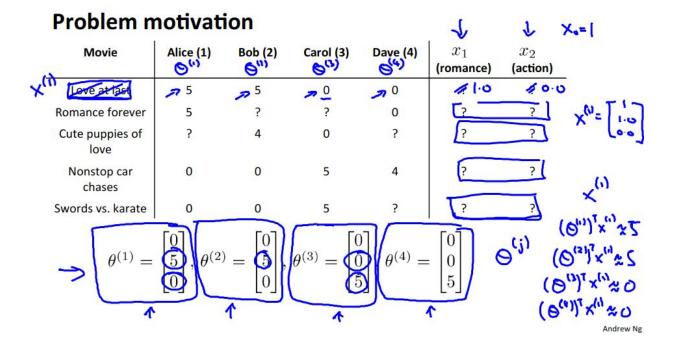
$$\theta^{(j)}_k := \theta^{(j)}_k - \alpha \left(\sum_{i:r(i,j)=1} ((\theta^{(j)})^T x^{(i)} - y^{(i,j)}) x^{(i)}_k + \lambda \theta^{(j)}_k \right) \ (\text{for } k \neq 0)$$

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Collaborative filtering:

Sometimes it may be difficult to set specific feature values for all movies. Instead of doing that we can let our users rate their taste on different features. Using those inputs we can calculate the specific values for different movies.

Problem motivation x_1 Movie Alice (1) Bob (2) Carol (3) **Dave (4)** x_2 (romance) (action) Love at last 5 5 0 0 0.9 0 Romance forever 5 ? ? 0 1.0 0.01 ? Cute puppies of ? 4 0 0.99 0 love Nonstop car 0 0 5 4 0.1 1.0 chases 5 ? 0 0 0.9 Swords vs. karate 0



Optimization algorithm

Given $\underline{\theta^{(1)},\ldots,\theta^{(n_u)}}$, to learn $\underline{x^{(i)}}$:

$$\Rightarrow \min_{x^{(i)}} \frac{1}{2} \sum_{j:r(i,j)=1} ((\underline{\theta^{(j)}})^T x^{(i)} - \underline{y^{(i,j)}})^2 + \frac{\lambda}{2} \sum_{k=1}^n (x_k^{(i)})^2$$

Given $\theta^{(1)}, \dots, \theta^{(n_u)}$, to learn $\underline{x^{(1)}, \dots, x^{(n_m)}}$:

$$\min_{x^{(1)},...,x^{(n_m)}} \frac{1}{2} \sum_{i=1}^{n_m} \sum_{j:r(i,j)=1} ((\theta^{(j)})^T x^{(i)} - y^{(i,j)})^2 + \frac{\lambda}{2} \sum_{i=1}^{n_m} \sum_{k=1}^n (x_k^{(i)})^2$$

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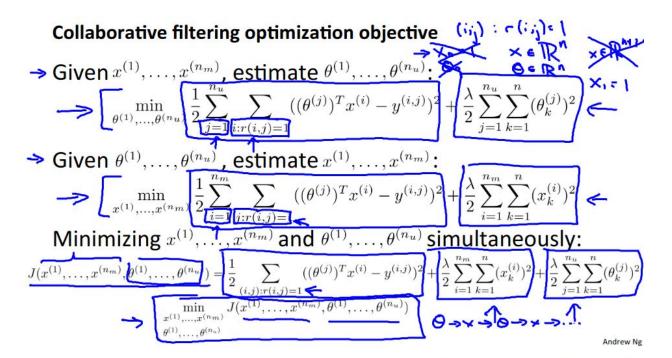
Collaborative filtering

Given $\underline{x^{(1)},\dots,x^{(n_m)}}$ (and movie ratings), can estimate $\underline{\theta^{(1)},\dots,\theta^{(n_u)}}$

Given
$$\frac{\theta^{(1)},\ldots,\theta^{(n_u)}}{\text{can estimate }x^{(1)},\ldots,x^{(n_m)}}$$

Collaborative filtering algorithm:

In the previous section we learned that by collaborative filtering we can train theta and X one after another. That means we need to use X to get min theta and to get min X need to use theta. But in collaborative filtering algorithm we can minimize our perimeter theta and X simultaneously.



Collaborative filtering algorithm

- \rightarrow 1. Initialize $x^{(1)}, \dots, x^{(n_m)}, \theta^{(1)}, \dots, \theta^{(n_u)}$ to small random values.
- ⇒ 2. Minimize $J(x^{(1)}, ..., x^{(n_m)}, \theta^{(1)}, ..., \theta^{(n_u)})$ using gradient descent (or an advanced optimization algorithm). E.g. for every $j = 1, ..., n_u, i = 1, ..., n_m$:

$$x_{k}^{(i)} := x_{k}^{(i)} - \alpha \left(\sum_{j:r(i,j)=1} ((\theta^{(j)})^{T} x^{(i)} - y^{(i,j)}) \theta_{k}^{(j)} + \lambda x_{k}^{(i)} \right)$$

$$\theta_{k}^{(j)} := \theta_{\underline{k}}^{(j)} - \alpha \left(\sum_{i:r(i,j)=1} ((\theta^{(j)})^{T} x^{(i)} - y^{(i,j)}) x_{k}^{(i)} + \lambda \theta_{k}^{(j)} \right)$$

$$\overline{\mathbf{d}}_{\underline{k}}^{(i)} \mathbf{d}_{\underline{k}}^{(i)} \mathbf{d}_{\underline{k}}^{(i)}$$

3. For a user with parameters $\underline{\theta}$ and a movie with (learned) features \underline{x} , predict a star rating of $\underline{\theta}^T \underline{x}$.

0,

Vectorize low rank matrix factorization:

Instead of using loops to calculate collaborative filtering we can use vectorize implementation. It will make computation easier.

nm: 5 **Collaborative filtering** Movie Alice (1) Bob (2) Carol (3) Dave (4) Love at last 5 5 0 0 Romance forever 5 Cute puppies of love Nonstop car chases Swords vs. karate

To improve user experience online tvs need to suggest related movies or videos. In that case they use existing user data to do that. Suppose, movie x1 is watched by a user. Now if you want to test if x2 movie is related to x1 then you need to following steps.

Finding related movies

For each product i, we learn a feature vector $x^{(i)} \in \mathbb{R}^n$.

How to find
$$\frac{\text{movies } j}{\|x^{(i)} - x^{(j)}\|} \rightarrow \text{movie } j$$
 and i are "similar"

5 most similar movies to movie i:

Find the 5 movies j with the smallest $\|x^{(i)} - x^{(j)}\|$.

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Implementation details : mean normalization :

Suppose a new user just signed up. He hasn't rated any movie yet. How do you suggest movies or other stuff? In that case, mean normalization will help us.

Users who have not rated any movies

Movie
 Alice (1)
 Bob (2)
 Carol (3)
 Dave (4)
 Eve (5)

 Love at last

$$\underline{5}$$
 $\underline{5}$
 0
 0
 ?
 0

 Romance forever
 5
 ?
 ?
 0
 ?
 0

 Cute puppies of love
 ?
 4
 0
 ?
 ?
 0
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 0
 ?
 0
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To apply this algorithm you need to compute the average rating for every single movie. Then subtract the avg value from original value. Then use this formula for user j, on movie i predict:

