

CSC 4020 Fundamental of Machine Learning: Linear Regression

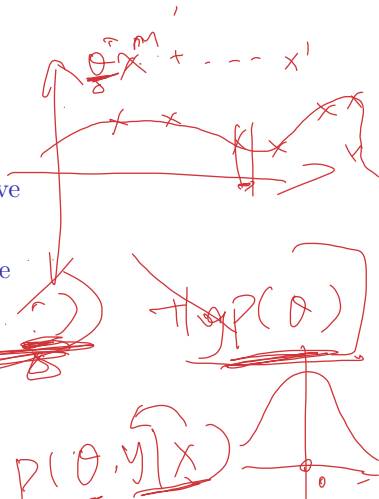
Baoyuan Wu
School of Data Science, CUHK-SZ

January 25/27, 2021

Outline

- 1 Some illustrations and review of last week
- 2 Linear Regression: A Deterministic Perspective
- 3 Linear Regression: A Probabilistic Perspective
 - Probabilistic modeling
 - Robust linear regression
 - Ridge regression
 - Lasso regression
- 4 Generalized Linear Regression

max a post (MAP)



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- Welcome to the office hour at **Wednesday 10:30–11:30am in DY 411.**

Review of last week

- **Probability theory:**

- Discrete probability distributions: Bernoulli, Binomial, Beta
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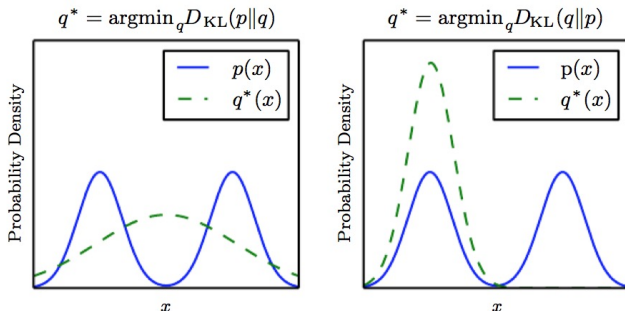
- Discrete probability distributions: Bernoulli, Binomial, Beta
- Continuous probability distributions: Gaussian, Student t , Laplace

- **Information theory:**

- Information
- Entropy, marginal/conditional/joint entropy, relative entropy (KL divergence, mutual information)

Properties of KL divergence

- $D(p_X(x)||q_X(x)) \geq 0$ with equality if and only if $p_X(x) = q_X(x)$.
- $D(p_X(x)||q_X(x)) \neq D(q_X(x)||p_X(x))$



One constraint with respect to q is missing at last time, *i.e.*, it is the **single mode distribution!** More detailed derivations could be found at <https://dibyaghosh.com/blog/probability/kldivergence.html>.

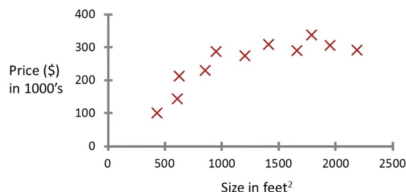
Linear regression

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- m indicates the number of training samples; x denotes the input variable/feature; y denotes the output variable.

Size in feet ² (x)	Price in 1000's (y)
2104	460
1416	232
1514	315
852	178
...	...



Linear hypothesis function

- Our goal is to find a linear hypothesis function to well fit the training data D , *i.e.*,

$$h_{\theta}(x) = \theta_0 + \theta_1\phi(x) = [\theta_0, \theta_1][1; \phi(x)] = \hat{\phi}(x)^{\top} \theta \quad (1)$$

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- Given x , $h_{\theta}(x)$ is a **linear function** of $\theta = [\theta_0; \theta_1]$. This is why it is called **linear regression**.

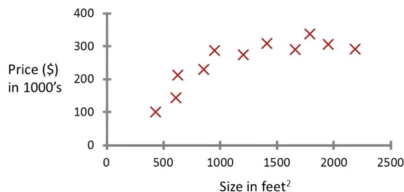
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- Then, given D , how to learn θ ?



Cost function

- We design the following **cost function** to minimize the difference between the prediction $h_{\boldsymbol{\theta}}(x_i)$ and the ground-truth value y_i , *i.e.*,

$$J(\boldsymbol{\theta}) = \frac{1}{2} \sum_{i=1}^m (h_{\boldsymbol{\theta}}(x_i) - y_i)^2 \quad (2)$$

$$= \frac{1}{2} \sum_{i=1}^m (\theta_0 + \theta_1 x_i - y_i)^2, \quad (3)$$

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- $J(\boldsymbol{\theta})$ is a **convex** or **non-convex** function? What is the shape of it?

- The linear regression is formulated to the following optimization problem

$$\boldsymbol{\theta}^* = \arg \min_{\boldsymbol{\theta}} J(\boldsymbol{\theta}) = \frac{1}{2} \sum_{i=1}^m (\bar{\mathbf{x}}_i^\top \boldsymbol{\theta} - y_i)^2. \quad (5)$$

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- $\boldsymbol{\theta}$ can be updated by **gradient descent algorithm**,

$$\boldsymbol{\theta} \leftarrow \boldsymbol{\theta} - \alpha \frac{\partial J(\boldsymbol{\theta})}{\partial \boldsymbol{\theta}}, \quad \frac{\partial J(\boldsymbol{\theta})}{\partial \boldsymbol{\theta}} = \sum_{i=1}^m (\bar{\mathbf{x}}_i^\top \boldsymbol{\theta} - y_i) \bar{\mathbf{x}}_i \quad (6)$$

where α is called step-size or learning rate.

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- Does gradient descent always converge to the optimal solution? (Plot the trajectory of gradient descent on curve or contours)

Analytical solution

- If we set the gradient to 0, then we can get the following solution

$$\frac{\partial J(\boldsymbol{\theta})}{\partial \boldsymbol{\theta}} = \sum_{i=1}^m (\bar{\mathbf{x}}_i^\top \boldsymbol{\theta} - y_i) \bar{\mathbf{x}}_i = \mathbf{X}^\top \mathbf{X} \boldsymbol{\theta} - \mathbf{X}^\top \mathbf{y} = 0 \quad (7)$$

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which are called **normal equation** and **ordinary least squares** (OLS) solution, respectively. $\mathbf{X} = [\bar{\mathbf{x}}_1^\top; \bar{\mathbf{x}}_2^\top; \dots; \bar{\mathbf{x}}_m^\top] \in \mathbb{R}^{m \times d}$.

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- Since there is a closed-form solution, why do we need gradient descent algorithm?

Geometric interpretation

- Since $\boldsymbol{\theta}^* = (\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top \mathbf{y}$, then the predictions of \mathbf{X} can be obtained by

$$\hat{\mathbf{y}} = \mathbf{X} \boldsymbol{\theta}^* = \mathbf{X} (\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top \mathbf{y}, \quad (9)$$

which corresponds to the **orthogonal projection** of \mathbf{y} onto the column space of \mathbf{X} .

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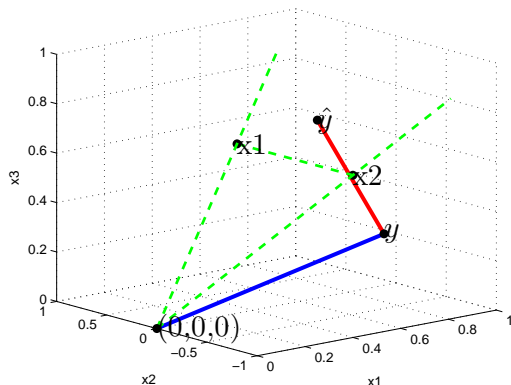
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$$\mathbf{X} = \begin{pmatrix} 1 & 2 \\ 1 & -2 \\ 1 & 2 \end{pmatrix},$$

$$\mathbf{y} = \begin{pmatrix} 8.89 \\ 0.61 \\ 1.77 \end{pmatrix}$$



Normal equation vs. gradient descent

	Size (feet ²)	Number of bedrooms	Number of floors	Age of home (years)	Price (\$1000)
x_0	x_1	x_2	x_3	x_4	y
1	2104	5	1	45	460
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1	1534	3	2	30	315
1	852	2	1	36	178

$X = \begin{bmatrix} 1 & 2104 & 5 & 1 & 45 \\ 1 & 1416 & 3 & 2 & 40 \\ 1 & 1534 & 3 & 2 & 30 \\ 1 & 852 & 2 & 1 & 36 \end{bmatrix}$
 $m \times (n+1)$

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 m -dimensional vector

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$m \times (n+1)$ m -dimensional vector

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Gradient Descent	Normal Equation
Need to choose alpha	No need to choose alpha
Needs many iterations	No need to iterate
$O(kn^2)$	$O(n^3)$, need to calculate inverse of $X^T X$
Works well when n is large	Slow if n is very large

- We assume that the relationship between the input variable/feature \mathbf{x} and the output variable y is

$$y = \boldsymbol{\theta}^\top \mathbf{x} + e, \text{ where } e \sim \mathcal{N}(0, \sigma^2), \quad (10)$$

where e is called **observation noise** or **residual error**, and it is independent with any specific input \mathbf{x} .

Probabilistic modeling

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- Thus, the output y can also be seen as a random variable, and its conditional probability is formulated as

$$p(y|\mathbf{x}, \boldsymbol{\theta}) = \mathcal{N}(\boldsymbol{\theta}^\top \mathbf{x}, \sigma^2) \quad (11)$$

Maximum log-likelihood estimation

- The parameter $\boldsymbol{\theta}$ can be learned by maximum log-likelihood estimation (MLE), given the training dataset $D = \{(\mathbf{x}_i, y_i)\}_{i=1}^m$, as follows

$$\boldsymbol{\theta}_{MLE} = \arg \max_{\boldsymbol{\theta}} \log \mathcal{L}(\boldsymbol{\theta} | D) \quad (12)$$

$$= \sum_i^m \log p(y | \mathbf{x}, \boldsymbol{\theta}) = \sum_i^m \log \mathcal{N}(\boldsymbol{\theta}^\top \mathbf{x}, \sigma^2) \quad (13)$$

$$= -\log(\sigma^m (2\pi)^{\frac{m}{2}}) - \frac{1}{2\sigma^2} \sum_i^m (y_i - \boldsymbol{\theta}^\top \mathbf{x}_i)^2 \quad (14)$$

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- Removing the constants w.r.t. $\boldsymbol{\theta}$,

$$\boldsymbol{\theta}_{MLE} = \arg \min_{\boldsymbol{\theta}} \frac{1}{2} \sum_i^m (y_i - \boldsymbol{\theta}^\top \mathbf{x}_i)^2, \quad (15)$$

which is exactly same with the cost function from the deterministic perspective.

Robust linear regression

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- Let's see the loss curve of the residual sum of squares (RSS),

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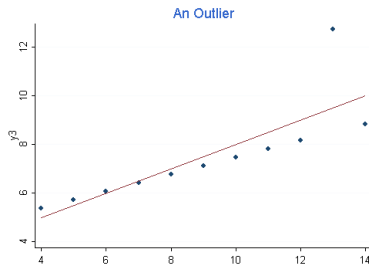
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- How to alleviate the significant influence of outliers?

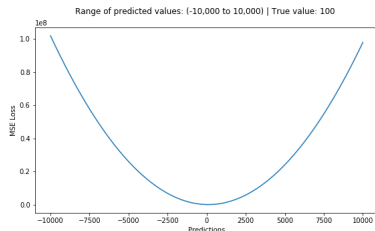
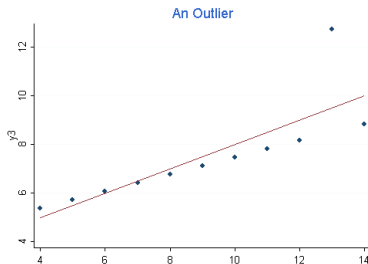


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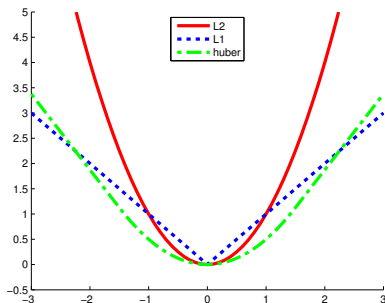
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- The curves of ℓ_1 and ℓ_2 losses are shown ad follows.
- When the residual is large, the ℓ_1 loss is much smaller than the ℓ_2 loss, such that the influence of outliers could be alleviated.



Robust linear regression

- Actually, the above ℓ_1 loss can also be derived from the probabilistic perspective, by assuming that

$$p(y|\mathbf{x}, \boldsymbol{\theta}, b) = \text{Lap}(y|\mathbf{x}, \boldsymbol{\theta}, b) \propto \exp\left(-\frac{1}{b}|y - \boldsymbol{\theta}^\top \mathbf{x}|\right) \quad (18)$$

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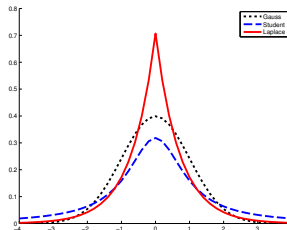
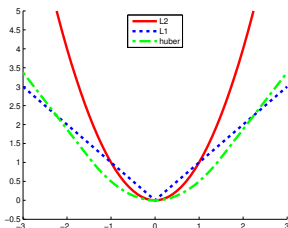
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- Applying the maximum log-likelihood estimation (MLE), we will obtain

$$\boldsymbol{\theta}_{MLE} = \arg \max_{\boldsymbol{\theta}} \log \mathcal{L}(\boldsymbol{\theta}|D) = \sum_i^m \log p(y_i|\mathbf{x}_i, \boldsymbol{\theta}) \quad (19)$$

$$\equiv \arg \min_{\boldsymbol{\theta}} \sum_{i=1}^m |\bar{\mathbf{x}}_i^\top \boldsymbol{\theta} - y_i| \quad (20)$$



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$$\boldsymbol{\theta}_{MLE} = \arg \min_{\boldsymbol{\theta}} \sum_{i=1}^m |\mathbf{x}_i^{\top} \boldsymbol{\theta} - y_i| \quad (21)$$

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- However, the ℓ_1 loss function is non-differentiable and non-linear. The gradient descent algorithm cannot be adopted.

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$$\min_{\boldsymbol{\theta}, \mathbf{t}} \sum_i^m t_i \quad (22)$$

$$s.t. \quad -t_i \leq \mathbf{x}_i^\top \boldsymbol{\theta} - y_i \leq t_i, 1 \leq i \leq m. \quad (23)$$

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- It is called **iteratively reweighted least squares** method.

Ridge regression

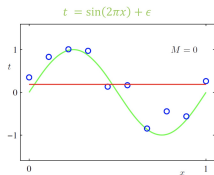
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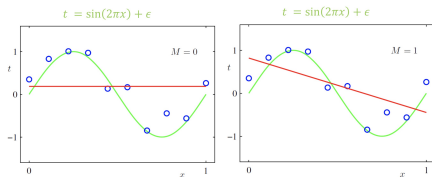
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- Is there other more theoretical approaches? SURE!



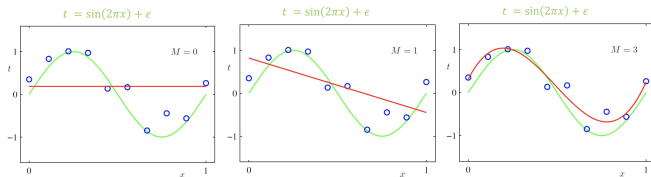
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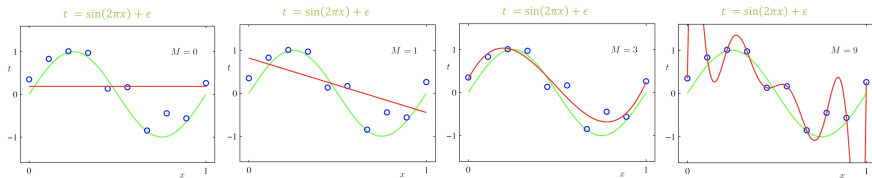
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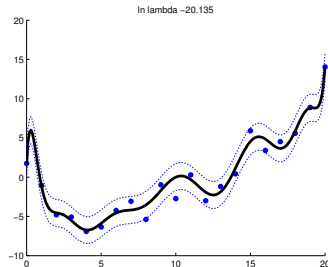


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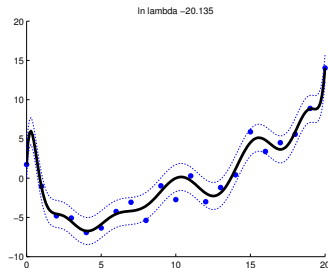


Ridge regression



- Let's see one simple example, we use a polynomial function with 14 degree to fit $m = 21$ data points. The learned curve is very “wiggly” (see above).

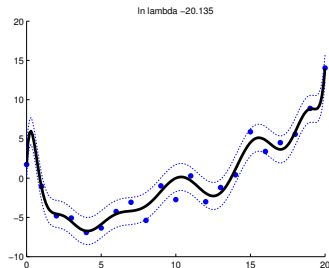
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- There are many large positive/negative values, such that a small change of features could lead to significant change of output.

Ridge regression

- How to get smaller parameter values?

Ridge regression

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- We can assume that the parameter follow a zero-mean Gaussian prior

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$$\boldsymbol{\theta}_{MAP} = \arg \max_{\boldsymbol{\theta}} \sum_i^m \log p(y|\mathbf{x}, \boldsymbol{\theta}) + \log p(\boldsymbol{\theta}) \quad (28)$$

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- The corresponding closed-form solution is given by

$$\boldsymbol{\theta}_{MAP} = (\lambda \mathbf{I} + \mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top \mathbf{y}. \quad (31)$$

Ridge regression

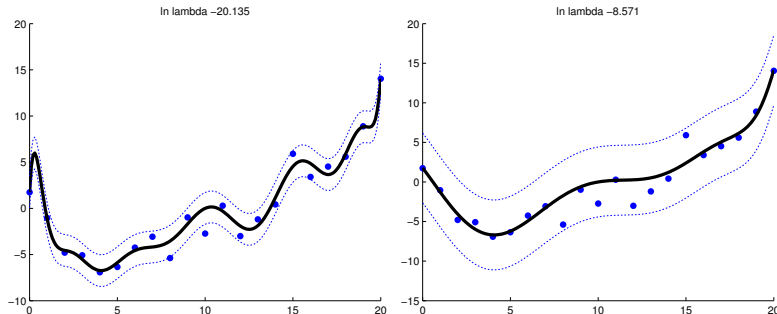
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Ridge regression

- The above method is also known as **ridge regression**, or **penalized least squares**.
- In general, adding a Gaussian prior to the parameters of a model to encourage them to be small is called ℓ_2 **regularization** or **weight decay**.

Ridge regression

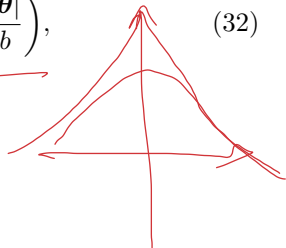
- The above method is also known as **ridge regression**, or **penalized least squares**.
- In general, adding a Gaussian prior to the parameters of a model to encourage them to be small is called ℓ_2 **regularization** or **weight decay**.
- As shown below, when we set a larger λ , *i.e.*, more weight on the prior, the resulting curve will be smoother.



Lasso regression

- We can replace the Gaussian prior by a Laplacian prior, *i.e.*,

$$p(\boldsymbol{\theta}) = \text{Lap}(\boldsymbol{\theta}|\mathbf{0}, b) = \frac{1}{2b} \exp\left(-\frac{|\boldsymbol{\theta}|}{b}\right), \quad (32)$$



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$$p(\boldsymbol{\theta}) = \text{Lap}(\boldsymbol{\theta}|\mathbf{0}, b) = \frac{1}{2b} \exp\left(-\frac{|\boldsymbol{\theta}|}{b}\right), \quad (32)$$

- The combination of the Gaussian distribution of $p(y|\mathbf{x}, \boldsymbol{\theta})$ and the Laplacian prior, leading to

assume $t = |\boldsymbol{\theta}|$

$$\boldsymbol{\theta}_{MAP} = \arg \max_{\boldsymbol{\theta}} \sum_{i=1}^m \log p(y|\mathbf{x}, \boldsymbol{\theta}) + \log p(\boldsymbol{\theta}) \quad (33)$$

① min t
s.t. $-\tau \leq \theta \leq \tau$

$$= \sum_{i=1}^m \log \mathcal{N}(\boldsymbol{\theta}^\top \mathbf{x}, \sigma^2) + \text{Lap}(\boldsymbol{\theta}|\mathbf{0}, b) \quad (34)$$

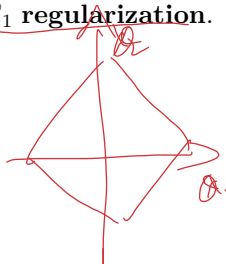
② $|a| = \min_u \left(\frac{a^2}{u} + u \right)$ $\Rightarrow u = |a|$ $\theta = a$

$$\equiv \arg \min_{\boldsymbol{\theta}} \sum_{i=1}^m (\bar{\mathbf{x}}_i^\top \boldsymbol{\theta} - y_i)^2 + \lambda |\boldsymbol{\theta}| \quad (35)$$

$\min_{\boldsymbol{\theta}} \min_u \left(\frac{a^2}{u} + u \right) \left[\min_{\boldsymbol{\theta}} f(\boldsymbol{\theta}) + R(a), \text{ s.t. } \boldsymbol{\theta} = a \right]$

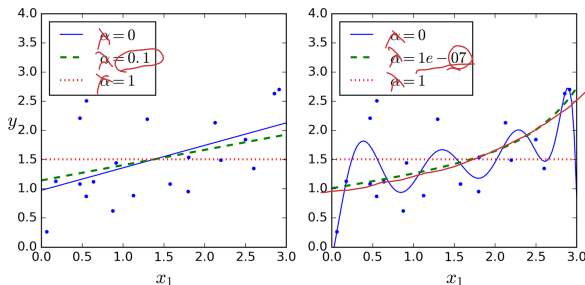
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Lasso regression

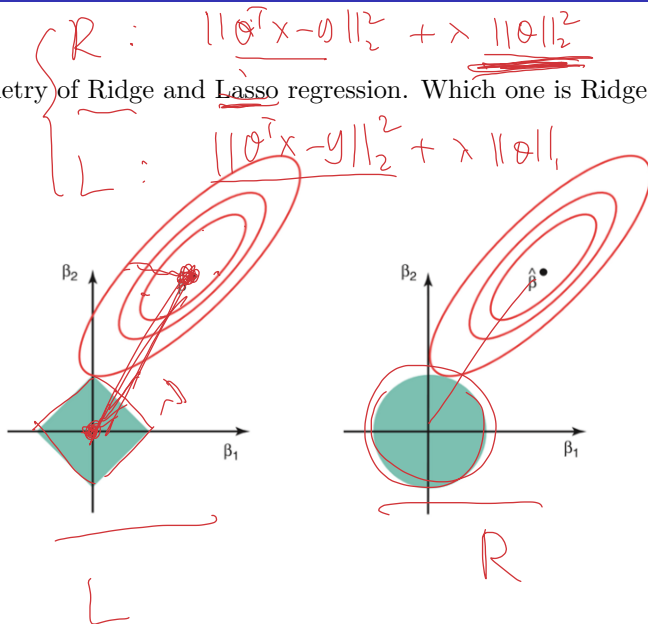
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$$y = \underline{Q} x_1 + \theta_0$$

Geometry of Ridge and Lasso regression

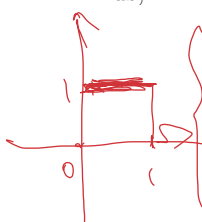
- Geometry of Ridge and Lasso regression. Which one is Ridge?



Summary of different linear regressions

Note that the uniform distribution will not change the mode of the likelihood.

Thus, MAP estimation with a uniform prior corresponds to MLE.

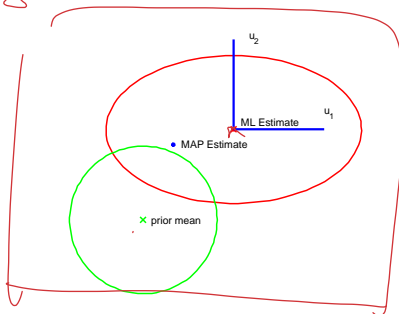


$p(y x, \theta)$	$p(\theta)$	regression method
Gaussian	<u>Uniform</u>	<u>Least squares</u>
Gaussian	<u>Gaussian</u>	Ridge regression
Gaussian	Laplace	<u>Lasso regression</u>
<u>Laplace</u>	<u>(Uniform)</u>	<u>Robust regression</u>
Student	Uniform	Robust regression

$$\min_{\theta} \frac{1}{2} \|\theta^T x - y\|_2^2$$

$$+ \lambda \|\theta\|_2^2$$

$$+ \lambda \|\theta\|_1$$



$$\theta^T \phi(x)$$

$$[x^m \dots x^1]$$

Generalized linear regression

- Linear model:

$$\left\{ \begin{array}{l} \mu(\mathbf{x}|\boldsymbol{\theta}) = \boldsymbol{\theta}^\top \phi(\mathbf{x}), \\ y(\mathbf{x}|\boldsymbol{\theta}) \sim f(\mu(\mathbf{x}|\boldsymbol{\theta})), \end{array} \right. \quad \begin{array}{l} (36) \\ (37) \end{array}$$

where f denotes a distribution function.

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- Generalized linear model (GLM):

$$\mu(\mathbf{x}|\boldsymbol{\theta}) = \underline{g}^{-1}(\boldsymbol{\theta}^\top \phi(\mathbf{x})), \quad (38)$$

$$\underline{y(\mathbf{x}|\boldsymbol{\theta})} \sim \underline{f(\mu(\mathbf{x}|\boldsymbol{\theta}))}, \quad (39)$$

where g is called **link function**, which is required to be monotonically increasing differentiable.

②

$g(x) = x$

①

Generalized linear regression

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where g is called **link function**, which is required to be monotonically increasing differentiable.

- The standard linear model is a special case of GLM with $g(a) = a$.

Why we need generalized linear regression

- Why we need generalized linear model? Let's see one example.

In the early stages of a disease epidemic, the rate at which new cases occur can often increase exponentially through time. Hence, if μ_i is the expected number of new cases on day t_i , a model of the form

$$\mu_i = \gamma \exp(\delta t_i)$$

seems appropriate.

- Such a model can be turned into GLM form, by using a **log link** so that

$$\log(\mu_i) = \log(\gamma) + \delta t_i = \beta_0 + \beta_1 t_i.$$

- Since this is a count, the Poisson distribution (with expected value μ_i) is probably a reasonable distribution to try.

$$u = \theta^T x$$

$$\beta^T t$$

Log linear regression

- **Poisson distribution** The Poisson distribution is popular for modeling the number of times an event occurs in (an interval of time) or space.

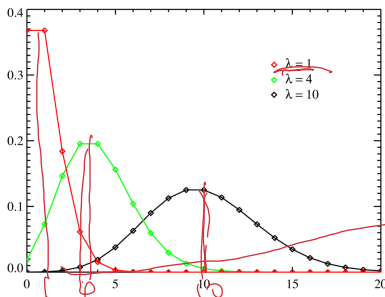


Log linear regression

- **Poisson distribution** The Poisson distribution is popular for modeling the number of times an event occurs in an interval of time or space.
- A discrete random variable X is said to have a Poisson distribution with parameter $\lambda > 0$ if for $k = 0, 1, 2, \dots$, the probability mass function of X is given by

$$f(k; \lambda) = P(X = k | \lambda) = \frac{\lambda^k e^{-\lambda}}{k!}, \quad (40)$$

where e is Euler's number ($e = 2.71828\dots$), we k is the number of occurrences, $k!$ is the factorial of k .



$$E(k) = \lambda$$
$$\text{Var}(k) = \lambda$$

Log linear regression

- We assume that the conditional probability follows

$$P(y_i|\mathbf{x}_i, \boldsymbol{\theta}) = \textit{Poisson}(\lambda_i) = \frac{\lambda^{y_i} e^{-\lambda}}{y_i!}, \quad \ln \lambda_i = \boldsymbol{\theta}^\top \mathbf{x}_i \quad (41)$$

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- The log-likelihood function is formulated as follows

$$\text{max} \quad \mathcal{L}(\boldsymbol{\theta}) = \sum_{i=1}^m \ln P(y_i|\mathbf{x}_i, \boldsymbol{\theta}) = \sum_{i=1}^m \underbrace{y_i \ln \lambda_i - \lambda_i - \log y_i!}_{\text{log-likelihood for one sample}} \quad (42)$$

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- We have

$$\frac{\partial \mathcal{L}(\boldsymbol{\theta})}{\partial \boldsymbol{\theta}} = \sum_{i=1}^m (y_i - e^{\boldsymbol{\theta}^\top \mathbf{x}_i}) \mathbf{x}_i = 0 \Rightarrow \ln y_i = (\boldsymbol{\theta}^*)^\top \mathbf{x}_i \quad (43)$$

$$\frac{\partial \mathcal{L}(\boldsymbol{\theta})}{\partial \lambda_i} = \left(\frac{y_i}{\lambda_i} - 1 \right); \quad \frac{\partial \lambda_i}{\partial \boldsymbol{\theta}} = \frac{\exp(\boldsymbol{\theta}^\top \mathbf{x}_i)}{\lambda_i} \cdot \mathbf{x}_i$$

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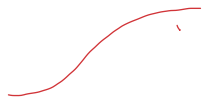
- Plot the log-linear regression as below.

Logistic regression

- We assume that the conditional probability follows

$$\mu_i = g(\theta^T x_i)$$

$$P(y_i | x_i, \theta, N) = \text{Bin}(y_i | N, \mu_i) = \binom{N}{y_i} \mu_i^{y_i} (1 - \mu_i)^{N - y_i}, \quad \mu_i = \frac{1}{1 + e^{-\theta^T x_i}}. \quad (44)$$



$$\mu(1 + e^{-a}) = 1$$

$$g(a) = \frac{1}{1 + e^{-a}}$$

$$\Rightarrow e^{-a} = \frac{1}{\mu} - 1$$

$$a = \ln\left(\frac{1}{\mu} - 1\right)$$

$$g(a) = \dots$$

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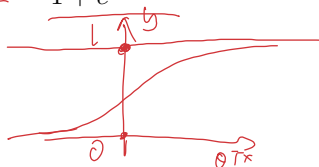
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$$p(y_i = k | \boldsymbol{\theta}) =$$



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$$\mathcal{L}(\boldsymbol{\theta}) = \sum_{i=1}^m \log P(y_i|\mathbf{x}_i, \boldsymbol{\theta}) = y_i \log \mu_i + (N - y_i) \log(1 - \mu_i) \quad (45)$$

- We have

$$\frac{\partial \mathcal{L}(\boldsymbol{\theta})}{\partial \boldsymbol{\theta}} = \sum_{i=1}^m (y_i - N\mu_i) \mathbf{x}_i = 0 \quad \Rightarrow \quad \frac{y_i}{N} = \mu_i = \frac{1}{1 + e^{-\boldsymbol{\theta}^\top \mathbf{x}_i}}. \quad (46)$$

- Since the $\sigma(a) = \frac{1}{1+e^{-a}}$ is called **sigmoid function** or **logit function**, the above model is called **logit regression** or **logistic regression**.

Logistic regression

- We assume that the conditional probability follows

$$P(y_i|\mathbf{x}_i, \boldsymbol{\theta}, N) = \text{Bin}(y_i|N, \mu_i) = \binom{N}{y_i} \mu_i^{y_i} (1 - \mu_i)^{N-y_i}, \quad \mu_i = \frac{1}{1 + e^{-\boldsymbol{\theta}^\top \mathbf{x}_i}}. \quad (44)$$

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- Since the $\sigma(a) = \frac{1}{1+e^{-a}}$ is called **sigmoid function** or **logit function**, the above model is called **logit regression** or **logistic regression**.
- Since $\frac{y_i}{N} \in [0, 1]$, it can be seen as the posterior probability. Thus, logistic regression is a **classification model**, rather than regression.

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$$g(x) = x$$

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- Linear model is the linear function of the parameter θ , rather than the input feature
- Linear model is a special case of generalized linear model, while generalized linear model is not always linear
- Choosing different linear models is equivalent to choosing different distributions of $p(y|\mathbf{x}, \theta)$ and $p(\theta)$, according to the task and the data

Reading material

- <https://www.stat.cmu.edu/~ryantibs/advmethods/notes/glm.pdf>