

Machine Learning

1. Linear Regression

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Machine Learning



1. The Regression Problem

- 2. Simple Linear Regression
- 3. Multiple Regression
- 4. Variable Interactions
- 5. Model Selection
- 6. Case Weights

Example



Example: how does gas consumption depend on external temperature? (Whiteside, 1960s).

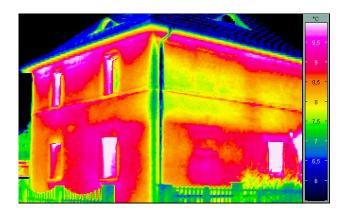
weekly measurements of

- average external temperature
- total gas consumption (in 1000 cubic feets)

A third variable encodes two heating seasons, before and after wall insulation.

How does gas consumption depend on external temperature?

How much gas is needed for a given temperature ?

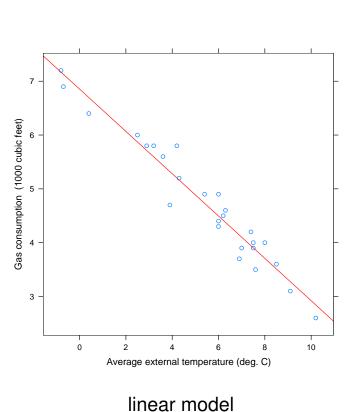


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Example

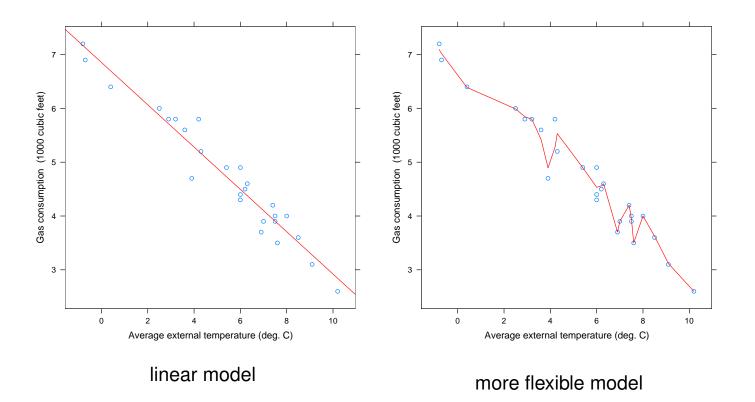
Machine Learning / 1. The Regression Problem





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Machine Learning / 1. The Regression Problem

Variable Types and Coding



The most common variable types:

numerical / interval-scaled / quantitative

where differences and quotients etc. are meaningful, usually with domain $\mathcal{X} := \mathbb{R}$, e.g., temperature, size, weight.

nominal / discrete / categorical / qualitative / factor

where differences and quotients are not defined, usually with a finite, enumerated domain, e.g., $\mathcal{X} := \{\text{red}, \text{green}, \text{blue}\}$ or $\mathcal{X} := \{a, b, c, \dots, y, z\}.$

ordinal / ordered categorical

where levels are ordered, but differences and quotients are not defined,

usually with a finite, enumerated domain, e.g., $\mathcal{X} := \{$ small, medium, large $\}$

Variable Types and Coding

Nominals are usually encoded as binary dummy variables:

$$\delta_{x_0}(X) := \left\{ \begin{array}{ll} 1, \mbox{ if } X = x_0, \\ 0, \mbox{ else} \end{array} \right.$$

one for each $x_0 \in \mathcal{X}$ (but one).

Example: $\mathcal{X} := \{ red, green, blue \}$

Replace

one variable X with 3 levels: red, green, blue

by

```
two variables \delta_{red}(X) and \delta_{green}(X) with 2 levels each: 0, 1
```

X	$\delta_{\mathrm{red}}(X)$	$\delta_{\text{green}}(X)$
red	1	0
green blue	0	1
blue	0	0
	1	1

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Machine Learning / 1. The Regression Problem

The Regression Problem Formally

Let

 X_1, X_2, \ldots, X_p be random variables called **predictors** (or **inputs**, **covariates**, **features**).

Let $\mathcal{X}_1, \mathcal{X}_2, \ldots, \mathcal{X}_p$ be their domains.

We write shortly

$$X := (X_1, X_2, \dots, X_p)$$

for the vector of random predictor variables and

 $\mathcal{X} := \mathcal{X}_1 \times \mathcal{X}_2 \times \cdots \times \mathcal{X}_p$

for its domain.

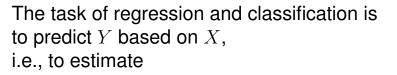
- Y be a random variable called **target** (or **output**, **response**). Let \mathcal{Y} be its domain.
- $\mathcal{D} \subseteq \mathcal{X} \times \mathcal{Y}$ be a (multi)set of instances of the unknown joint distribution p(X, Y) of predictors and target called **data**. \mathcal{D} is often written as enumeration

 $\mathcal{D} = \{(x_1, y_1), (x_2, y_2), \dots, (x_n, y_n)\}$





The Regression Problem Formally (v0)



$$\hat{y}(x) \coloneqq r(x) \coloneqq E(Y \,|\, X = x) = \int y \, p(y|x) dy$$

based on data (called regression function).

If *Y* is numerical, the task is called **regression**.

If Y is nominal, the task is called **classification**.

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Machine Learning / 1. The Regression Problem

The Regression Problem Formally (v1)



Let \mathcal{X} be any set (called **predictor space**). Given

- a set $\mathcal{D}^{\text{train}} \subseteq \mathcal{X} \times \mathbb{R}$ of data (called **training set**),

compute a regression function

$$\hat{y}: \mathcal{X} \to \mathbb{R}$$

s.t. for a set $\mathcal{D}^{\text{test}} \subseteq \mathcal{X} \times \mathbb{R}$ of data (called **test set**) not available during training, the test error

$$\operatorname{err}(\hat{y}; \mathcal{D}^{\mathsf{test}}) := \frac{1}{|D^{\mathsf{test}}|} \sum_{(x,y)\in\mathcal{D}^{\mathsf{test}}} (y - \hat{y}(x))^2$$

is minimal.

The Regression Problem Formally (v2)



Let \mathcal{X} be any set (called **predictor space**). Given

- a set $\mathcal{D}^{\text{train}} \subseteq \mathcal{X} \times \mathbb{R}$ of data (called **training set**),

– a loss function $\ell : \mathbb{R} \times \mathbb{R} \to \mathbb{R}$ that measures how bad it is to predict value \hat{y} if the true value is y,

compute a **regression function**

$$\hat{y}: \mathcal{X} \to \mathbb{R}$$

s.t. for a set $\mathcal{D}^{\text{test}} \subseteq \mathcal{X} \times \mathbb{R}$ of data (called **test set**) not available during training, the test error

$$\operatorname{err}(\hat{y}; \mathcal{D}^{\mathsf{test}}) := \frac{1}{|D^{\mathsf{test}}|} \sum_{(x,y)\in\mathcal{D}^{\mathsf{test}}} \ell(y, \hat{y}(x))$$

is minimal.

Examples:

$$\ell(y,\hat{y}) := (y - \hat{y})^2$$

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Machine Learning / 1. The Regression Problem

The Regression Problem Formally (v3)



Let \mathcal{X} be any set (called **predictor space**), and \mathcal{Y} be any set (called **target space**).

Given

- a set $\mathcal{D}^{\text{train}} \subseteq \mathcal{X} \times \mathcal{Y}$ of data (called **training set**),
- a loss function $\ell : \mathcal{Y} \times \mathcal{Y} \to \mathbb{R}$ that measures how bad it is to predict value \hat{y} if the true value is y,

compute a prediction function

$$\hat{y}: \mathcal{X} \to \mathcal{Y}$$

s.t. for a set $\mathcal{D}^{\text{test}} \subseteq \mathcal{X} \times \mathbb{R}$ of data (called **test set**) not available during training, the test error

$$\operatorname{err}(\hat{y}; \mathcal{D}^{\mathsf{test}}) := \frac{1}{|D^{\mathsf{test}}|} \sum_{(x,y)\in\mathcal{D}^{\mathsf{test}}} \ell(y, \hat{y}(x))$$

is minimal.

Examples:

$$\mathcal{Y} := \mathbb{R}, \quad \ell(y, \hat{y}) := (y - \hat{y})^2$$

The Regression Problem Formally (v4)

Let \mathcal{X} be any set (called **predictor space**),

- ${\mathcal Y}$ be any set (called **target space**), e.g., and
- $p:\mathcal{X}\times\mathcal{Y}
 ightarrow\mathbb{R}^+_0$ be a joint distribution / density.

Given

- a sample $\mathcal{D}^{\text{train}} \subseteq \mathcal{X} \times \mathcal{Y}$ (called **training set**), drawn from *p*,
- a loss function $\ell : \mathcal{Y} \times \mathcal{Y} \to \mathbb{R}$ that measures how bad it is to predict value \hat{y} if the true value is y,

compute a prediction function

$$\hat{y}: \mathcal{X} \to \mathcal{Y}$$

s.t. for another sample $\mathcal{D}^{\text{test}} \subseteq \mathcal{X} \times \mathbb{R}$ (called **test set**) drawn from the same distribution *p*, not available during training, the test error

$$\operatorname{err}(\hat{y}; \mathcal{D}^{\mathsf{test}}) := \frac{1}{|D^{\mathsf{test}}|} \sum_{(x,y)\in\mathcal{D}^{\mathsf{test}}} \ell(y, \hat{y}(x))$$

is minimal.

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Machine Learning / 1. The Regression Problem

The Regression Problem Formally (v5)

Let \mathcal{X} be any set (called **predictor space**), \mathcal{Y} be any set (called **target space**), and $p: \mathcal{X} \times \mathcal{Y} \to \mathbb{R}^+_0$ be a joint distribution / density.

Given

- a sample $\mathcal{D}^{\text{train}} \subseteq \mathcal{X} \times \mathcal{Y}$ (called **training set**), drawn from *p*,
- a loss function $\ell : \mathcal{Y} \times \mathcal{Y} \to \mathbb{R}$ that measures how bad it is to predict value \hat{y} if the true value is y,

compute a prediction function

$$\hat{y}: \mathcal{X} \to \mathcal{Y}$$

with minimal risk

$$\operatorname{risk}(\hat{y};p) := \int_{\mathcal{X} \times \mathcal{Y}} \ell(y,\hat{y}) \, p(x,y) \, d(x,y)$$

Explanation: $\operatorname{risk}(\hat{y}; p)$ can be estimated by the **empirical risk**

$$\operatorname{risk}(\hat{y}; \mathcal{D}^{\mathsf{test}}) := \frac{1}{|D^{\mathsf{test}}|} \sum_{(x,y) \in \mathcal{D}^{\mathsf{test}}} \ell(y, \hat{y}(x))$$

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Simple Examples: Single Predictor vs. Multiple Predictors 0 8.0 °0 0 7.5 7.0 6.5 6.0 5.5 x2 x1 5.0 0.2 0.4 0.6 0.8 0.0 1.0 х

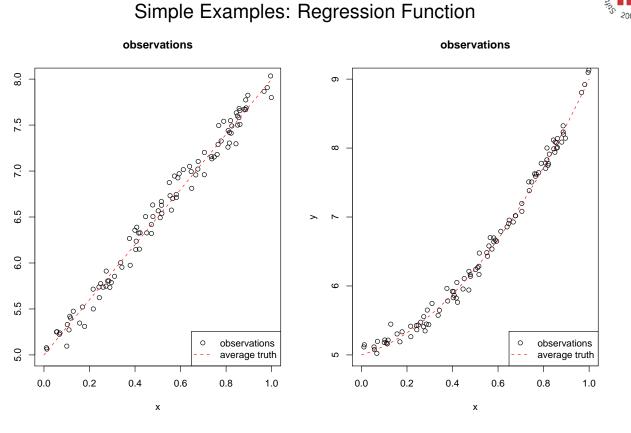
single predictor:

y = 3x + 5

multiple predictors:

 $y = x_1 + 2x_2 + 5$

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linear regression function:

y = 3x + 5

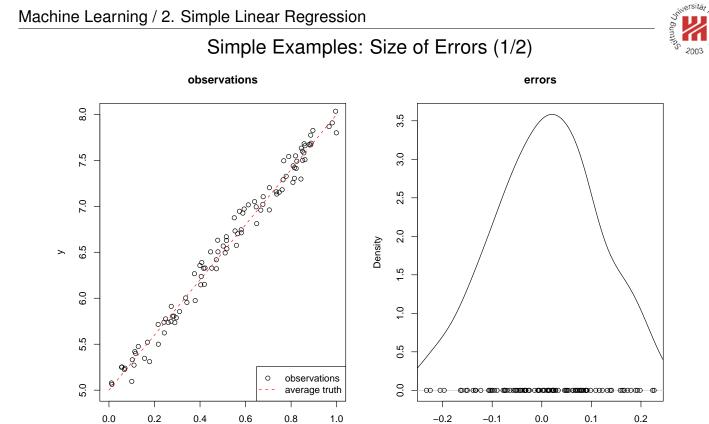
х

non-linear regression function:

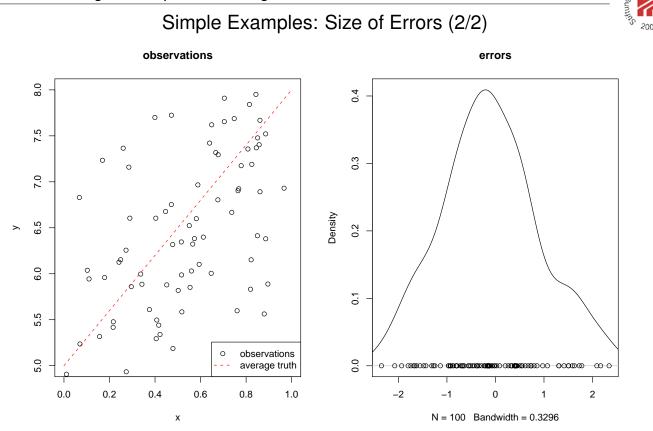
 $y = 3x^2 + x + 5$

N = 100 Bandwidth = 0.037

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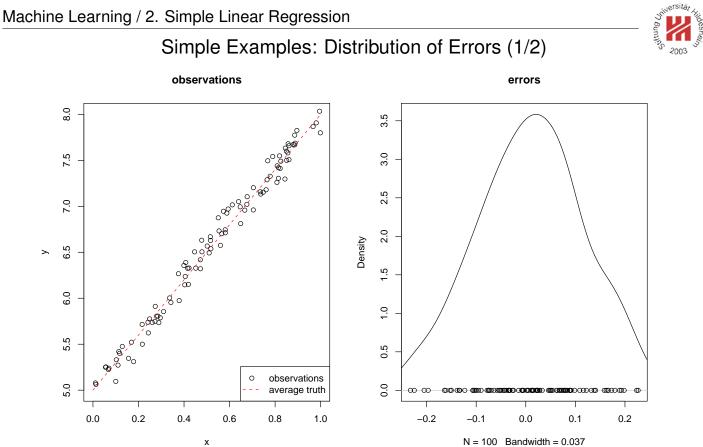


Small errors vs. ...

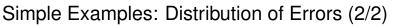


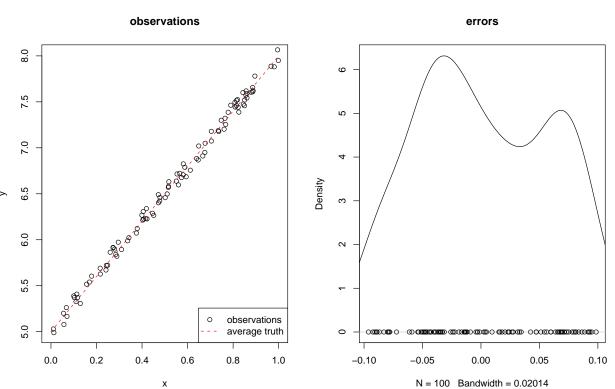
... large errors.

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Normally distributed errors vs. ...





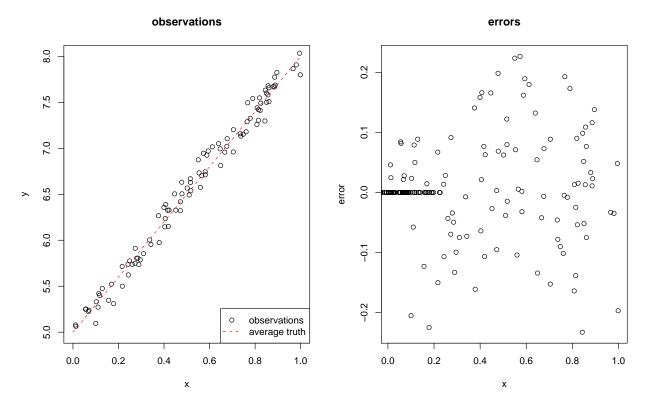
... uniformly distributed errors.

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Machine Learning / 2. Simple Linear Regression

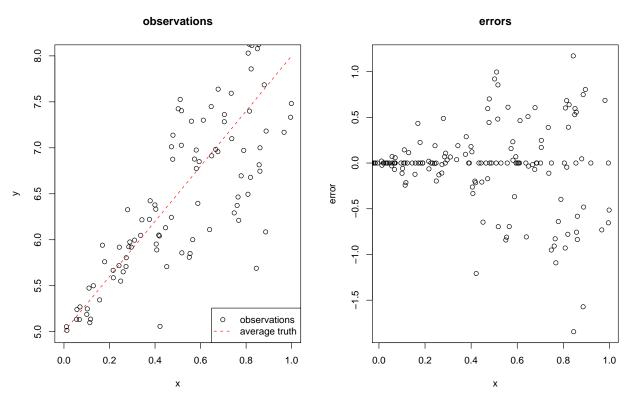


Simple Examples: Homoscedastic vs. Heteroscedastic Errors (1/2)



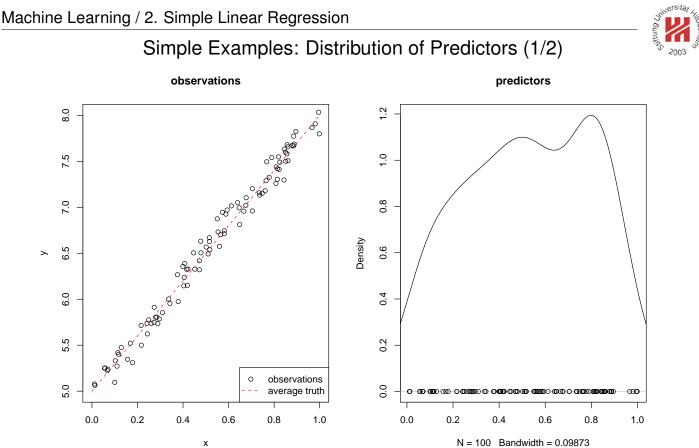
Errors do not depend on predictors (homoscedastic) vs. ...

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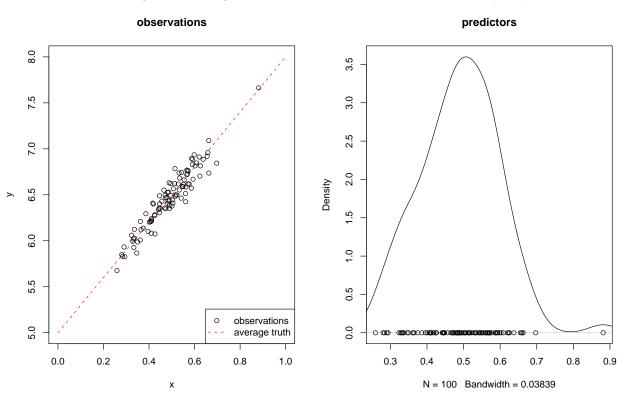
... errors do depend on predictors (heteroscedastic).

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Predictors are uniformly distributed vs. ...

Simple Examples: Distribution of Predictors (2/2)



... predictors are normally distributed.

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Machine Learning / 2. Simple Linear Regression

Simple Linear Regression Model

Make it simple:

- the predictor X is simple, i.e., one-dimensional $(X = X_1)$.
- r(x) is assumed to be linear:

$$r(x) = \beta_0 + \beta_1 x$$

• assume that the variance does not depend on *X*:

$$Y = \beta_0 + \beta_1 X + \epsilon, \quad E(\epsilon | X) = 0, V(\epsilon | X) = \sigma^2$$

- 3 parameters:
 - β_0 intercept (sometimes also called bias)

 β_1 slope

 σ^2 variance

Simple Linear Regression Model

parameter estimates

$$\hat{\beta}_0, \hat{\beta}_1, \hat{\sigma}^2$$

fitted line

$$\hat{r}(x) := \hat{\beta}_0 + \hat{\beta}_1 x$$

predicted / fitted values

 $\hat{y}_i := \hat{r}(x_i)$

residuals

$$\hat{\epsilon}_i := y_i - \hat{y}_i = y_i - (\hat{\beta}_0 + \hat{\beta}_1 x_i)$$

residual sums of squares (RSS) / square loss / L2 loss

$$RSS = \sum_{i=1}^{n} \hat{\epsilon}_i^2$$

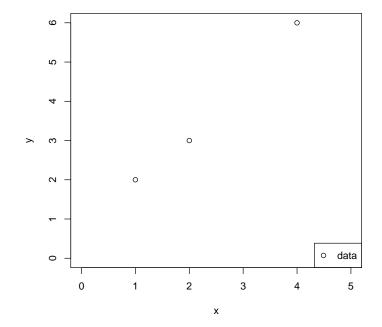
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Machine Learning / 2. Simple Linear Regression

How to estimate the parameters?

Example:

Given the data $\mathcal{D} := \{(1, 2), (2, 3), (4, 6)\}$, predict a value for x = 3.



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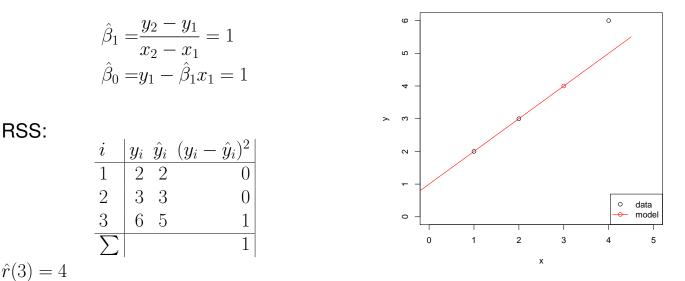


How to estimate the parameters?

Example:

Given the data $\mathcal{D} := \{(1, 2), (2, 3), (4, 6)\}$, predict a value for x = 3.

Line through first two points:



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Machine Learning / 2. Simple Linear Regression

How to estimate the parameters?

Example:

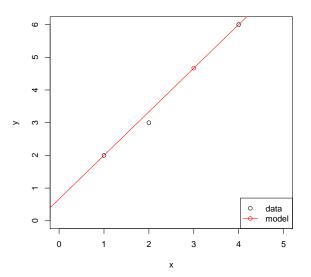
Given the data $\mathcal{D} := \{(1, 2), (2, 3), (4, 6)\}$, predict a value for x = 3.

Line through first and last point:

$$\hat{\beta}_1 = \frac{y_3 - y_1}{x_3 - x_1} = 4/3 = 1.333$$
$$\hat{\beta}_0 = y_1 - \hat{\beta}_1 x_1 = 2/3 = 0.667$$

RSS:

$$\frac{i}{1} \quad \frac{y_i}{2} \quad \frac{\hat{y}_i}{2} \quad \frac{(y_i - \hat{y}_i)^2}{0} \\
\frac{\hat{y}_i}{2} \quad \hat{y}_i \quad \frac{(y_i - \hat{y}_i)^2}{0} \\
\frac{\hat{y}_i}{2} \quad \hat{y}_i \quad \hat{y}_i \quad \frac{(y_i - \hat{y}_i)^2}{0} \\
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\hat{y}_i \quad \hat{y}_i \quad \hat{y}_i \quad \frac{(y_i - \hat{y}_i)^2}{0} \\
\hat{y}_i \quad \hat{y}_i$$







Least Squares Estimates / Definition

Journal South

In principle, there are many different methods to estimate the parameters $\hat{\beta}_0$, $\hat{\beta}_1$ and $\hat{\sigma}^2$ from data — depending on the properties the solution should have.

The **least squares estimates** are those parameters that minimize

RSS =
$$\sum_{i=1}^{n} \hat{\epsilon}_{i}^{2} = \sum_{i=1}^{n} (y_{i} - \hat{y}_{i})^{2} = \sum_{i=1}^{n} (y_{i} - (\hat{\beta}_{0} + \hat{\beta}_{1}x_{i}))^{2}$$

They can be written in closed form as follows:

$$\hat{\beta}_{1} = \frac{\sum_{i=1}^{n} (x_{i} - \bar{x})(y_{i} - \bar{y})}{\sum_{i=1}^{n} (x_{i} - \bar{x})^{2}}$$
$$\hat{\beta}_{0} = \bar{y} - \hat{\beta}_{1}\bar{x}$$
$$\hat{\sigma}^{2} = \frac{1}{n-2} \sum_{i=1}^{n} \hat{\epsilon}_{i}^{2}$$

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Machine Learning / 2. Simple Linear Regression

Least Squares Estimates / Proof

Proof (1/2):

$$RSS = \sum_{i=1}^{n} (y_i - (\hat{\beta}_0 + \hat{\beta}_1 x_i))^2$$
$$\frac{\partial RSS}{\partial \hat{\beta}_0} = \sum_{i=1}^{n} 2(y_i - (\hat{\beta}_0 + \hat{\beta}_1 x_i))(-1) \stackrel{!}{=} 0$$
$$\implies \quad n\hat{\beta}_0 = \sum_{i=1}^{n} (y_i - \hat{\beta}_1 x_i)$$



Least Squares Estimates / Proof



Proof (2/2):

$$RSS = \sum_{i=1}^{n} (y_i - (\hat{\beta}_0 + \hat{\beta}_1 x_i))^2$$

$$= \sum_{i=1}^{n} (y_i - (\bar{y} - \hat{\beta}_1 \bar{x}) - \hat{\beta}_1 x_i)^2$$

$$= \sum_{i=1}^{n} (y_i - \bar{y} - \hat{\beta}_1 (x_i - \bar{x}))^2$$

$$\frac{\partial RSS}{\partial \hat{\beta}_1} = \sum_{i=1}^{n} 2(y_i - \bar{y} - \hat{\beta}_1 (x_i - \bar{x}))(-1)(x_i - \bar{x}) \stackrel{!}{=} 0$$

$$\implies \hat{\beta}_1 = \frac{\sum_{i=1}^{n} (y_i - \bar{y})(x_i - \bar{x})}{\sum_{i=1}^{n} (x_i - \bar{x})^2}$$

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Machine Learning / 2. Simple Linear Regression

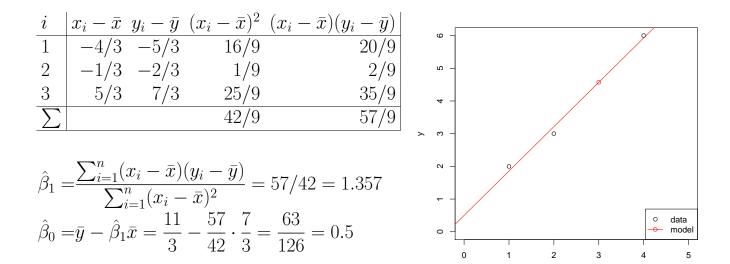
Least Squares Estimates / Example



Example:

Given the data $\mathcal{D} := \{(1, 2), (2, 3), (4, 6)\}$, predict a value for x = 3. Assume simple linear model.

 $\bar{x} = 7/3, \, \bar{y} = 11/3.$

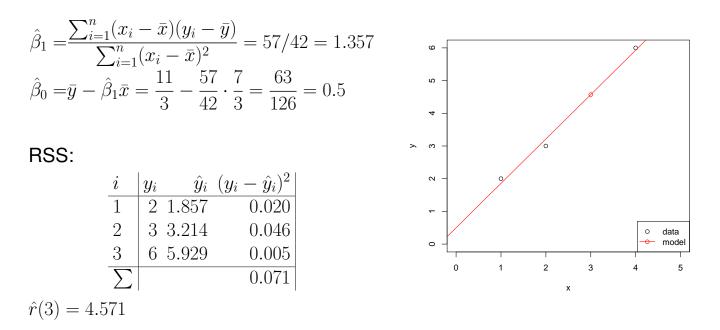


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Least Squares Estimates / Example

Example:

Given the data $\mathcal{D} := \{(1, 2), (2, 3), (4, 6)\}$, predict a value for x = 3. Assume simple linear model.



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A Generative Model



So far we assumed the model

$$Y = \beta_0 + \beta_1 X + \epsilon, \quad E(\epsilon | X) = 0, V(\epsilon | X) = \sigma^2$$

where we required some properties of the errors, but not its exact distribution.

If we make assumptions about its distribution, e.g.,

$$\epsilon | X \sim \mathcal{N}(0, \sigma^2)$$

and thus

$$Y|X = x \sim \mathcal{N}(\beta_0 + \beta_1 x, \sigma^2)$$

we can sample from this model.

Maximum Likelihood Estimates (MLE)



Let $p(X,Y\,|\,\theta)$ be a joint probability density function for X and Y with parameters $\theta.$

Likelihood:

$$L_{\mathcal{D}}(\theta) := \prod_{i=1}^{n} p(x_i, y_i \,|\, \theta)$$

The likelihood describes the probability of the data.

The **maximum likelihood estimates (MLE)** are those parameters that maximize the likelihood.

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Machine Learning / 2. Simple Linear Regression

Least Squares Estimates and Maximum Likelihood Estimates

Likelihood:

$$L_{\mathcal{D}}(\hat{\beta}_0, \hat{\beta}_1, \hat{\sigma}^2) := \prod_{i=1}^n \hat{p}(x_i, y_i) = \prod_{i=1}^n \hat{p}(y_i \mid x_i) p(x_i) = \prod_{i=1}^n \hat{p}(y_i \mid x_i) \prod_{i=1}^n p(x_i)$$

Conditional likelihood:

$$L_{\mathcal{D}}^{\mathsf{cond}}(\hat{\beta}_{0},\hat{\beta}_{1},\hat{\sigma}^{2}) := \prod_{i=1}^{n} \hat{p}(y_{i} \mid x_{i}) = \prod_{i=1}^{n} \frac{1}{\sqrt{2\pi}\hat{\sigma}} e^{-\frac{(y_{i}-\hat{y}_{i})^{2}}{2\hat{\sigma}^{2}}} = \frac{1}{\sqrt{2\pi}^{n}\hat{\sigma}^{n}} e^{\frac{1}{-2\hat{\sigma}^{2}}\sum_{i=1}^{n} (y_{i}-\hat{y}_{i})^{2}}$$

Conditional log-likelihood:

$$\log L_{\mathcal{D}}^{\mathsf{cond}}(\hat{\beta}_0, \hat{\beta}_1, \hat{\sigma}^2) \propto -n \log \hat{\sigma} - \frac{1}{2\hat{\sigma}^2} \sum_{i=1}^n (y_i - \hat{y}_i)^2$$

 \implies if we assume normality, the maximum likelihood estimates are just the least squares estimates.

Implementation Details

2003

(1) simple-regression(\mathcal{D}) : (2) sx := 0, sy := 0(3) for i = 1, ..., n do $\mathbf{s}\mathbf{x} := \mathbf{s}\mathbf{x} + x_i$ (4) $sy := sy + y_i$ (5) (6) Od (7) $\bar{x} := sx/n, \bar{y} := sy/n$ (8) a := 0, b := 0(9) for i = 1, ..., n do $a := a + (x_i - \bar{x})(y_i - \bar{y})$ (10) $b := b + (x_i - \bar{x})^2$ (11) (12) **Od** (13) $\beta_1 := a/b$ (14) $\beta_0 := \bar{y} - \beta_1 \bar{x}$ (15) return (β_0, β_1)

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Machine Learning / 2. Simple Linear Regression

Implementation Details



naive:

(1) simple-regression(\mathcal{D}) : (2) sx := 0, sy := 0(3) for i = 1, ..., n do $\mathbf{s}\mathbf{x} := \mathbf{s}\mathbf{x} + x_i$ (4)(5) $sy := sy + y_i$ (6) **Od** (7) $\bar{x} := sx/n, \bar{y} := sy/n$ (8) a := 0, b := 0(9) for i = 1, ..., n do $a := a + (x_i - \bar{x})(y_i - \bar{y})$ (10) $b := b + (x_i - \bar{x})^2$ (11) (12) **Od** (13) $\beta_1 := a/b$ (14) $\beta_0 := \bar{y} - \beta_1 \bar{x}$ (15) return (β_0, β_1)

single loop:

i simple-regression(\mathcal{D}) : sx := 0, sy := 0, sxx := 0, syy := 0, sxy := 03 for i = 1, ..., n do $\mathbf{sx} := \mathbf{sx} + x_i$ 4 $sy := sy + y_i$ 5 $\begin{aligned} \mathbf{sxx} &:= \mathbf{sxx} + x_i^2 \\ \mathbf{syy} &:= \mathbf{syy} + y_i^2 \end{aligned}$ 6 7 $sxy := sxy + x_i y_i$ 8 9 **od** 10 $\beta_1 := (n \cdot sxy - sx \cdot sy)/(n \cdot sxx - sx \cdot sx)$ $\prod \beta_0 := (\mathbf{s}\mathbf{y} - \beta_1 \cdot \mathbf{s}\mathbf{x})/n$ *12* **return** (β_0, β_1)



- 1. The Regression Problem
- 2. Simple Linear Regression
- **3. Multiple Regression**
- 4. Variable Interactions
- 5. Model Selection
- 6. Case Weights

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Machine Learning / 3. Multiple Regression

Several predictors



Several predictor variables X_1, X_2, \ldots, X_p :

$$Y = \beta_0 + \beta_1 X_1 + \beta_2 X_2 + \dots + \beta_P X_P + \epsilon$$
$$= \beta_0 + \sum_{i=1}^p \beta_i X_i + \epsilon$$

with p + 1 parameters $\beta_0, \beta_1, \ldots, \beta_p$.

Linear form



Several predictor variables X_1, X_2, \ldots, X_p :

$$Y = \beta_0 + \sum_{i=1}^p \beta_i X_i + \epsilon$$
$$= \langle \beta, X \rangle + \epsilon$$

where

$$\beta := \begin{pmatrix} \beta_0 \\ \beta_1 \\ \vdots \\ \beta_p \end{pmatrix}, \quad X := \begin{pmatrix} 1 \\ X_1 \\ \vdots \\ X_p \end{pmatrix},$$

Thus, the intercept is handled like any other parameter, for the artificial constant variable $X_0 \equiv 1$.

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Machine Learning / 3. Multiple Regression

Simultaneous equations for the whole dataset

For the whole dataset $(x_1, y_1), \ldots, (x_n, y_n)$:

$$\mathbf{Y} = \mathbf{X}\boldsymbol{\beta} + \boldsymbol{\epsilon}$$

where

$$\mathbf{Y} := \begin{pmatrix} y_1 \\ \vdots \\ y_n \end{pmatrix}, \quad \mathbf{X} := \begin{pmatrix} x_1 \\ \vdots \\ x_n \end{pmatrix} = \begin{pmatrix} x_{1,1} & x_{1,2} & \dots & x_{1,p} \\ \vdots & \vdots & \vdots & \vdots \\ x_{n,1} & x_{n,2} & \dots & x_{n,p} \end{pmatrix}, \quad \epsilon := \begin{pmatrix} \epsilon_1 \\ \vdots \\ \epsilon_n \end{pmatrix},$$

Least squares estimates



Least squares estimates
$$\hat{\beta}$$
 minimize
 $||\mathbf{Y} - \hat{\mathbf{Y}}||^2 = ||\mathbf{Y} - \mathbf{X}\hat{\beta}||^2$

The least squares estimates $\hat{\beta}$ are computed via

 $\mathbf{X}^T\mathbf{X}\hat{\boldsymbol{\beta}} = \mathbf{X}^T\mathbf{Y}$

Proof:

$$|\mathbf{Y} - \mathbf{X}\hat{eta}||^2 = \langle \mathbf{Y} - \mathbf{X}\hat{eta}, \mathbf{Y} - \mathbf{X}\hat{eta}
angle$$

$$\frac{\partial(\ldots)}{\partial\hat{\beta}} = 2\langle -\mathbf{X}, \mathbf{Y} - \mathbf{X}\hat{\beta} \rangle = -2(\mathbf{X}^T\mathbf{Y} - \mathbf{X}^T\mathbf{X}\hat{\beta}) \stackrel{!}{=} 0$$

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Machine Learning / 3. Multiple Regression

How to compute least squares estimates $\hat{\beta}$

Solve the $p \times p$ system of linear equations

$$\mathbf{X}^T \mathbf{X} \hat{\boldsymbol{\beta}} = \mathbf{X}^T \mathbf{Y}$$

i.e., Ax = b (with $A := \mathbf{X}^T \mathbf{X}, b = \mathbf{X}^T \mathbf{Y}, x = \hat{\beta}$).

There are several numerical methods available:

- 1. Gaussian elimination
- 2. Cholesky decomposition
- 3. QR decomposition

How to compute least squares estimates $\hat{\beta}$ / Example

Given is the following data:

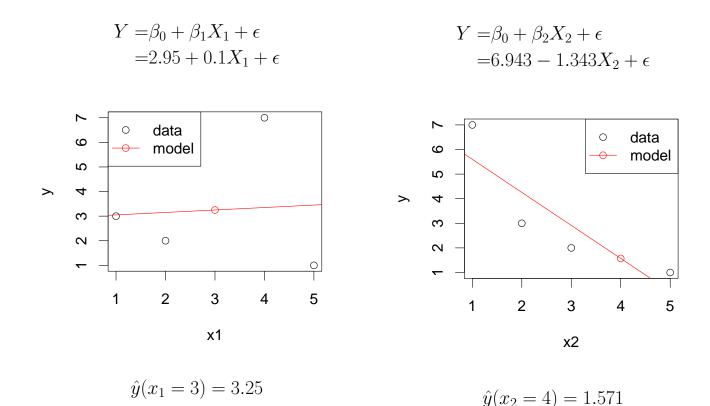
x_1	x_2	y
1	2	3
2	3	2
4	1	7
5	5	1

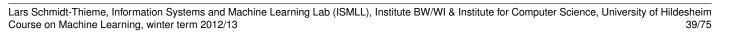
Predict a y value for $x_1 = 3, x_2 = 4$.

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Machine Learning / 3. Multiple Regression

How to compute least squares estimates $\hat{\beta}$ / Example









 $Y = \beta_0 + \beta_1 X_1 + \beta_2 X_2 + \epsilon$

How to compute least squares estimates $\hat{\beta}$ / Example



Now fit

to the data:

x_1	x_2	y
1	2	3
2	3	2
4	1	7
5	5	1

$$\mathbf{X} = \begin{pmatrix} 1 & 1 & 2 \\ 1 & 2 & 3 \\ 1 & 4 & 1 \\ 1 & 5 & 5 \end{pmatrix}, \quad \mathbf{Y} = \begin{pmatrix} 3 \\ 2 \\ 7 \\ 1 \end{pmatrix}$$
$$\mathbf{X}^{T}\mathbf{X} = \begin{pmatrix} 4 & 12 & 11 \\ 12 & 46 & 37 \\ 11 & 37 & 39 \end{pmatrix}, \quad \mathbf{X}^{T}\mathbf{Y} = \begin{pmatrix} 13 \\ 40 \\ 24 \end{pmatrix}$$

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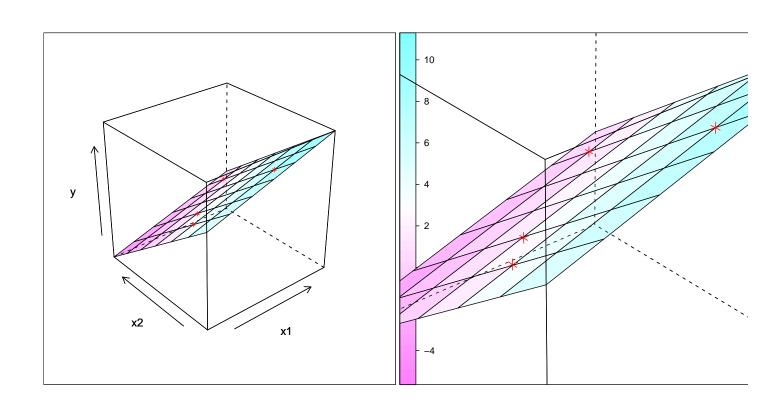
Machine Learning / 3. Multiple Regression



How to compute least squares estimates $\hat{\beta}$ / Example

$$\begin{pmatrix} 4 & 12 & 11 & | & 13 \\ 12 & 46 & 37 & | & 40 \\ 11 & 37 & 39 & | & 24 \end{pmatrix} \sim \begin{pmatrix} 4 & 12 & 11 & | & 13 \\ 0 & 10 & 4 & | & 1 \\ 0 & 16 & 35 & | & -47 \end{pmatrix} \sim \begin{pmatrix} 4 & 12 & 11 & | & 13 \\ 0 & 0 & 143 & | & -243 \end{pmatrix}$$
$$\sim \begin{pmatrix} 4 & 12 & 11 & | & 13 \\ 0 & 1430 & 0 & | & 1115 \\ 0 & 0 & 143 & | & -243 \end{pmatrix} \sim \begin{pmatrix} 286 & 0 & 0 & | & 1597 \\ 0 & 1430 & 0 & | & 1115 \\ 0 & 0 & 143 & | & -243 \end{pmatrix}$$
i.e.,
$$\hat{\beta} = \begin{pmatrix} 1597/286 \\ 1115/1430 \\ -243/143 \end{pmatrix} \approx \begin{pmatrix} 5.583 \\ 0.779 \\ -1.699 \end{pmatrix}$$

How to compute least squares estimates $\hat{\beta}$ / Example



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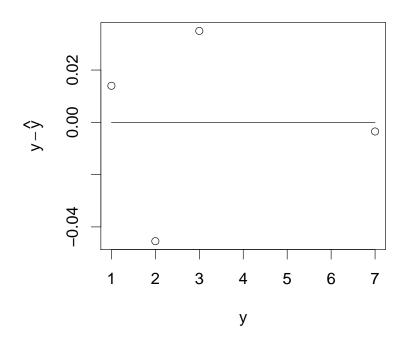
Machine Learning / 3. Multiple Regression

How to compute least squares estimates $\hat{\beta}$ / Example

To visually assess the model fit, a plot

residuals $\hat{\epsilon} = y - \hat{y}$ vs. true values y

can be plotted:



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The Normal Distribution (also Gaussian)

written as:

$$X \sim \mathcal{N}(\mu, \sigma^2)$$

with parameters:

 μ mean,

 σ standard deviance.

Ċ

probability density function (pdf):

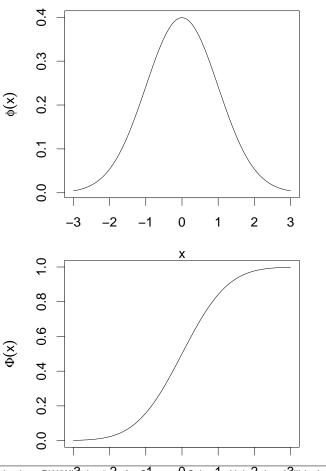
$$\phi(x) := \frac{1}{\sqrt{2\pi\sigma}} e^{-\frac{(x-\mu)^2}{2\sigma^2}}$$

cumulative distribution function (cdf):

$$\Phi(x) := \int_{-\infty}^x \phi(t) dt$$

 Φ^{-1} is called **quantile function**.

 Φ and Φ^{-1} have no analytical form, but have to be computed numerically.



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Machine Learning / 3. Multiple Regression

written as:

 $X \sim t_p$

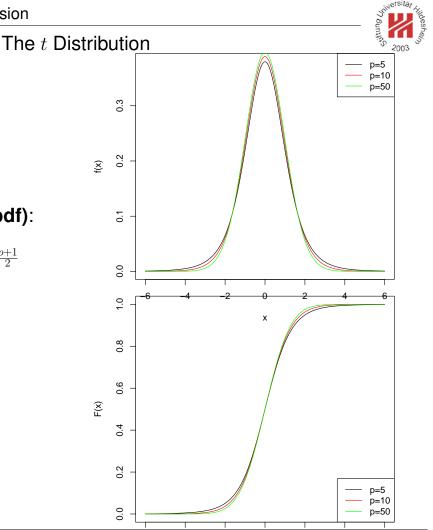
with parameter:

p degrees of freedom.

probability density function (pdf):

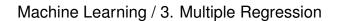
$$p(x) := \frac{\Gamma(\frac{p+1}{2})}{\sqrt{p \pi} \, \Gamma(\frac{p}{2})} (1 + \frac{x^2}{p})^{-\frac{p+1}{2}}$$

$$t_p \stackrel{p \to \infty}{\longrightarrow} \mathcal{N}(0, 1)$$



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written as:

$$X \sim \chi_p^2$$

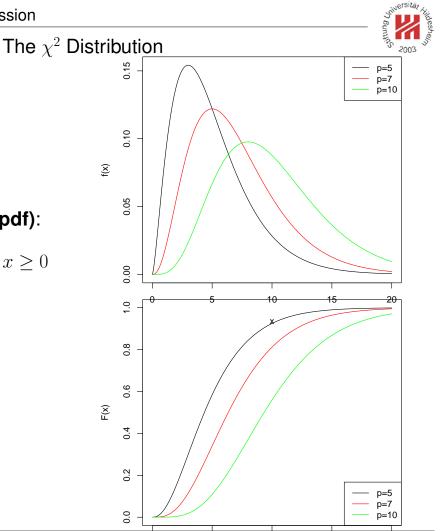
with parameter: *p* degrees of freedom.

probability density function (pdf):

$$p(x) := \frac{1}{\Gamma(p/2)2^{p/2}} x^{\frac{p}{2}-1} e^{-\frac{x}{2}}, \quad x \ge 0$$

If $X_1, \ldots, X_p \sim \mathcal{N}(0, 1)$, then

$$Y := \sum_{i=1}^p X_i^2 \sim \chi_p^2$$



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Machine Learning / 3. Multiple Regression

Parameter Variance



 $\hat{\beta} = (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{Y}$ is an unbiased estimator for β (i.e., $E(\hat{\beta}) = \beta$). Its variance is

$$V(\hat{\beta}) = (\mathbf{X}^T \mathbf{X})^{-1} \sigma^2$$

proof:

$$\hat{\beta} = (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \mathbf{Y} = (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T (\mathbf{X}\beta + \epsilon) = \beta + (\mathbf{X}^T \mathbf{X})^{-1} \mathbf{X}^T \epsilon$$

As
$$E(\epsilon) = 0$$
: $E(\hat{\beta}) = \beta$

$$V(\hat{\beta}) = E((\hat{\beta} - E(\hat{\beta}))(\hat{\beta} - E(\hat{\beta}))^{T})$$

= $E((\mathbf{X}^{T}\mathbf{X})^{-1}\mathbf{X}^{T}\epsilon\epsilon^{T}\mathbf{X}(\mathbf{X}^{T}\mathbf{X})^{-1})$
= $(\mathbf{X}^{T}\mathbf{X})^{-1}\sigma^{2}$

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Parameter Variance



An unbiased estimator for σ^2 is

$$\hat{\sigma}^2 = \frac{1}{n-p} \sum_{i=1}^n \hat{\epsilon}_i^2 = \frac{1}{n-p} \sum_{i=1}^n (y_i - \hat{y}_i)^2$$

If $\epsilon \sim \mathcal{N}(0,\sigma^2)$, then

$$\hat{\beta} \sim \mathcal{N}(\beta, (\mathbf{X}^T \mathbf{X})^{-1} \sigma^2)$$

Furthermore

$$(n-p)\hat{\sigma}^2\sim\sigma^2\chi^2_{n-p}$$

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Machine Learning / 3. Multiple Regression



Parameter Variance / Standardized coefficient

standardized coefficient ("z-score"):

$$z_i := \frac{\beta_i}{\widehat{\mathbf{se}}(\hat{\beta}_i)}, \quad \text{with } \widehat{\mathbf{se}}^2(\hat{\beta}_i) \text{ the } i \text{-th diagonal element of } (\mathbf{X}^T \mathbf{X})^{-1} \hat{\sigma}^2$$

 z_i would be $z_i \sim \mathcal{N}(0, 1)$ if σ is known (under $H_0 : \beta_i = 0$). With estimated $\hat{\sigma}$ it is $z_i \sim t_{n-p}$.

The Wald test for $H_0: \beta_i = 0$ with size α is:

reject
$$H_0$$
 if $|z_i| = |\frac{\hat{\beta}_i}{\widehat{\mathbf{se}}(\hat{\beta}_i)}| > F_{t_{n-p}}^{-1}(1-\frac{\alpha}{2})$

i.e., its *p*-value is

$$p-\mathsf{value}(H_0:\beta_i=0) = 2(1 - F_{t_{n-p}}(|z_i|)) = 2(1 - F_{t_{n-p}}(|\frac{\beta_i}{\widehat{\mathsf{se}}(\hat{\beta}_i)}|))$$

and small p-values such as 0.01 and 0.05 are good.

Parameter Variance / Confidence interval

The $1 - \alpha$ confidence interval for β_i :

$$\beta_i \pm F_{t_{n-p}}^{-1}(1-\frac{\alpha}{2})\widehat{\mathbf{se}}(\hat{\beta}_i)$$

For large *n*, $F_{t_{n-p}}$ converges to the standard normal cdf Φ .

As $\Phi^{-1}(1-\frac{0.05}{2})\approx 1.95996\approx 2,$ the rule-of-thumb for a 5% confidence interval is

$$\beta_i \pm 2\widehat{\mathbf{se}}(\hat{\beta}_i)$$

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Machine Learning / 3. Multiple Regression

Parameter Variance / Example

We have already fitted

$$\hat{Y} = \hat{\beta}_0 + \hat{\beta}_1 X_1 + \hat{\beta}_2 X_2$$

= 5.583 + 0.779X₁ - 1.699X₂

x_1	x_2	~	\hat{y}	$\hat{\epsilon}^2 = (y - \hat{y})^2$
1	2	3	2.965	0.00122
2	3	2	2.045	0.00207
4	1	7	7.003	0.0000122 0.000196
5	5	1	0.986	0.000196
RSS				0.00350

$\hat{\sigma}^2 = \frac{1}{n-p}$	$\sum_{i=1}^{n} \hat{\epsilon}_i^2 =$	$=\frac{1}{4-3}($).00350 =	= 0.00350
$(X^T X)^{-1} \hat{\sigma}^2 =$	$ \left(\begin{array}{c} 0.0\\ -0.0\\ -0.0 \end{array}\right) $	$\begin{array}{r} 0520 \ - \\ 0075 \ 0076 \ - \end{array}$	0.00075 0.00043 0.00020	-0.00076 -0.00020 0.00049
covariate	\hat{eta}_i	$\widehat{\mathbf{Se}}(\hat{\beta}_i)$	z-score	p-value
(intercept)	5.583	0.0721	77.5	0.0082
X_1	0.779	0.0207	37.7	0.0169
X_2	-1.699	0.0221	-76.8	0.0083

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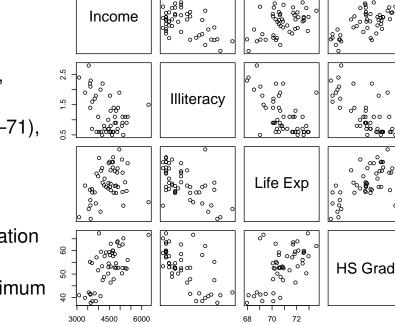


Parameter Variance / Example 2

Example: sociographic data of the 50 US states in 1977.

state dataset:

- income (per capita, 1974),
- illiteracy (percent of population, 1970),
- life expectancy (in years, 1969-71),
- percent high-school graduates (1970).
- population (July 1, 1975)
- murder rate per 100,000 population (1976)
- mean number of days with minimum temperature below freezing (1931–1960) in capital or large city
- land area in square miles



0.5

1.5

2.5

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Machine Learning / 3. Multiple Regression

Parameter Variance / Example 2



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 $\begin{aligned} \mathsf{Murder} = & \beta_0 + \beta_1 \mathsf{Population} + \beta_2 \mathsf{Income} + \beta_3 \mathsf{Illiteracy} \\ & + \beta_4 \mathsf{LifeExp} + \beta_5 \mathsf{HSGrad} + \beta_6 \mathsf{Frost} + \beta_7 \mathsf{Area} \end{aligned}$

n = 50 states, p = 8 parameters, n - p = 42 degrees of freedom.

Least squares estimators:

	Estimate	Std. Error	t value	Pr(> t)	
(Intercept)	1.222e+02	1.789e+01	6.831	2.54e-08	***
Population	1.880e-04	6.474e-05	2.905	0.00584	**
Income	-1.592e-04	5.725e-04	-0.278	0.78232	
Illiteracy	1.373e+00	8.322e-01	1.650	0.10641	
`Life Exp`	-1.655e+00	2.562e-01	-6.459	8.68e-08	* * *
` HS Grad`	3.234e-02	5.725e-02	0.565	0.57519	
Frost	-1.288e-02	7.392e-03	-1.743	0.08867	•
Area	5.967e-06	3.801e-06	1.570	0.12391	

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Machine Learning / 4. Variable Interactions





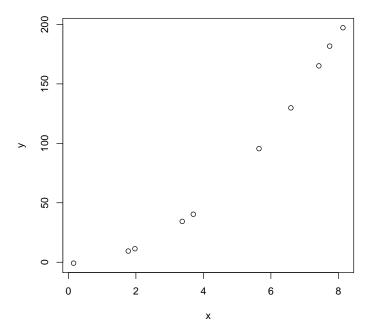
Assume a target variable does not depend linearly on a predictor variable, but say quadratic.

Example: way length vs. duration of a moving object with constant acceleration *a*.

$$s(t) = \frac{1}{2}at^2 + \epsilon$$

Can we catch such a dependency?

Can we catch it with a linear model?



Need for general transformations



To describe many phenomena, even more complex functions of the input variables are needed.

Example: the number of cells n vs. duration of growth t:

 $n = \beta e^{\alpha t} + \epsilon$

n does not depend on *t* directly, but on $e^{\alpha t}$ (with a known α).

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Machine Learning / 4. Variable Interactions

Need for variable interactions

In a linear model with two predictors

 $Y = \beta_0 + \beta_1 X_1 + \beta_2 X_2 + \epsilon$

Y depends on both, X_1 and X_2 .

But changes in X_1 will affect Y the same way, regardless of X_2 .

There are problems where X_2 mediates or influences the way X_1 affects Y, e.g. : the way length s of a moving object vs. its constant velocity v and duration t:

 $s = vt + \epsilon$

Then an additional 1s duration will increase the way length not in a uniform way (regardless of the velocity), but a little for small velocities and a lot for large velocities.

v and t are said to interact: y does not depend only on each predictor separately, but also on their product.

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Derived variables

All these cases can be handled by looking at **derived variables**, i.e., instead of

$$Y = \beta_0 + \beta_1 X_1^2 + \epsilon$$
$$Y = \beta_0 + \beta_1 e^{\alpha X_1} + \epsilon$$
$$Y = \beta_0 + \beta_1 X_1 \cdot X_2 + \epsilon$$

one looks at

$$Y = \beta_0 + \beta_1 X_1' + \epsilon$$

with

$$X'_{1} := X_{1}^{2}$$
$$X'_{1} := e^{\alpha X_{1}}$$
$$X'_{1} := X_{1} \cdot X_{2}$$

Derived variables are computed before the fitting process and taken into account either additional to the original variables or instead of.

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Machine Learning



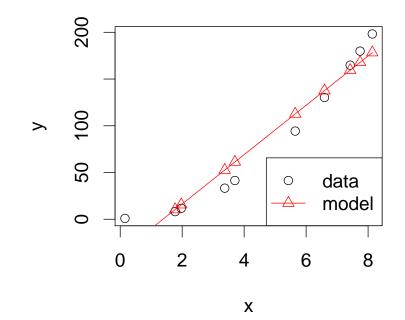
- 1. The Regression Problem
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- 6. Case Weights

Underfitting



sitä

2003

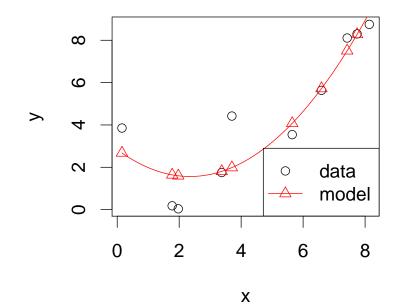


If a model does not well explain the data, e.g., if the true model is quadratic, but we try to fit a linear model, one says, the model **underfits**.

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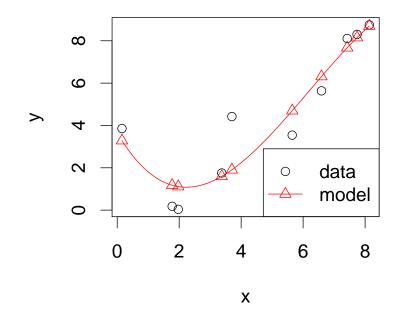
Machine Learning / 5. Model Selection

Overfitting / Fitting Polynomials of High Degree



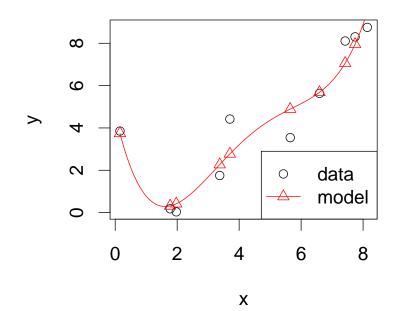
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Overfitting / Fitting Polynomials of High Degree



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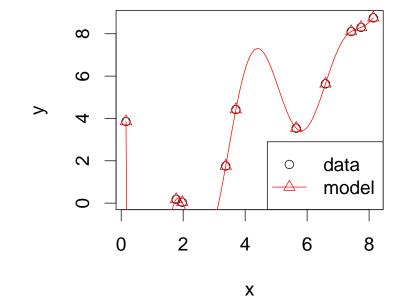








Overfitting / Fitting Polynomials of High Degree



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Machine Learning / 5. Model Selection

Overfitting / Fitting Polynomials of High Degree

If to data

$$(x_1, y_1), (x_2, y_2), \dots, (x_n, y_n)$$

consisting of n points we fit

 $X = \beta_0 + \beta_1 X_1 + \beta_2 X_2 + \dots + \beta_{n-1} X_{n-1}$

i.e., a polynomial with degree n-1, then this results in an interpolation of the data points

(if there are no repeated measurements, i.e., points with the same X_1 .)

As the polynomial

$$r(X) = \sum_{i=1}^{n} y_i \prod_{j \neq i} \frac{X - x_j}{x_i - x_j}$$

is of this type, and has minimal RSS = 0.





Model Selection Measures

Model selection means: we have a set of models, e.g.,

$$Y = \sum_{i=0}^{p-1} \beta_i X_i$$

indexed by p (i.e., one model for each value of p), make a choice which model **describes** the data best.

If we just look at losses / fit measures such as RSS, then

the larger p, the better the fit

or equivalently

the larger p, the lower the loss

as the model with p parameters can be **reparametrized** in a model with p' > p parameters by setting

$$\beta_i' = \begin{cases} \beta_i, \text{ for } i \leq p \\ 0, \text{ for } i > p \end{cases}$$

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Machine Learning / 5. Model Selection

Model Selection Measures



model selection measure = fit - complexity

or equivalently

model selection measure = loss + complexity

The smaller the loss (= lack of fit), the better the model.

The smaller the complexity, the simpler and thus better the model.

The model selection measure tries to find a trade-off between fit/loss and complexity.





Model Selection Measures



Akaike Information Criterion (AIC): (maximize)

 $\mathsf{AIC} := \log L - p$

or (minimize)

$$AIC := -2\log L + 2p = -2n\log(RSS/n) + 2p$$

Bayes Information Criterion (BIC) / Bayes-Schwarz Information Criterion: (maximize)

$$\mathsf{BIC} := \log L - \frac{p}{2} \log n$$

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Machine Learning / 5. Model Selection

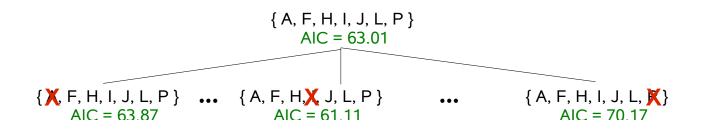
Variable Backward Selection



{ A, F, H, I, J, L, P } AIC = 63.01

Variable Backward Selection

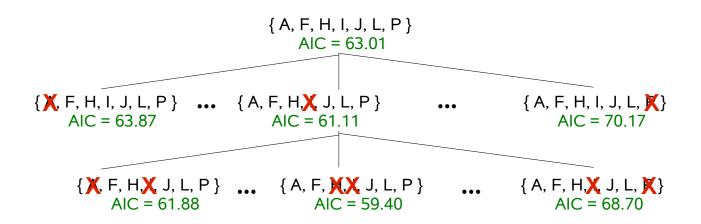




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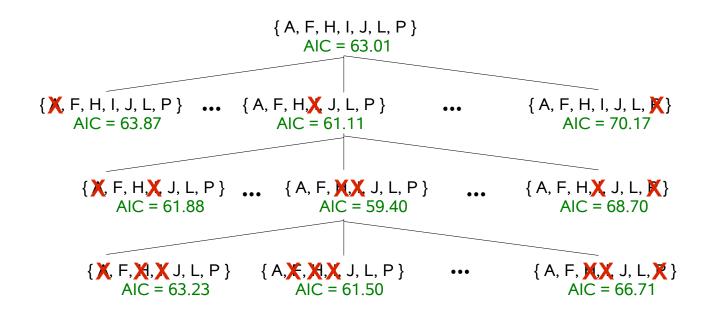
Machine Learning / 5. Model Selection

Variable Backward Selection



Variable Backward Selection





X removed variable

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Machine Learning / 5. Model Selection

Variable Backward Selection

full model:

	Estimate	Std. Error	t value	Pr(> t)	
(Intercept)	1.222e+02	1.789e+01	6.831	2.54e-08	* * *
Population	1.880e-04	6.474e-05	2.905	0.00584	* *
Income	-1.592e-04	5.725e-04	-0.278	0.78232	
Illiteracy	1.373e+00	8.322e-01	1.650	0.10641	
`Life Exp`	-1.655e+00	2.562e-01	-6.459	8.68e-08	* * *
'HS Grad'	3.234e-02	5.725e-02	0.565	0.57519	
Frost	-1.288e-02	7.392e-03	-1.743	0.08867	•
Area	5.967e-06	3.801e-06	1.570	0.12391	

AIC optimal model by backward selection:

	Estimate	Std. Error	t value	Pr(> t)	
(Intercept)	1.202e+02	1.718e+01	6.994	1.17e-08	***
Population	1.780e-04	5.930e-05	3.001	0.00442	* *
Illiteracy	1.173e+00	6.801e-01	1.725	0.09161	•
`Life Exp`	-1.608e+00	2.324e-01	-6.919	1.50e-08	* * *
Frost	-1.373e-02	7.080e-03	-1.939	0.05888	•
Area	6.804e-06	2.919e-06	2.331	0.02439	*

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How to do it in R



```
library(datasets);
library(MASS);
st = as.data.frame(state.x77);
mod.full = lm(Murder ~ ., data=st);
summary(mod.full);
mod.opt = stepAIC(mod.full);
summary(mod.opt);
```

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Machine Learning / 5. Model Selection

Shrinkage



Model selection operates by

- fitting models for a set of models with varying complexity and then picking the "best one" ex post,
- omitting some parameters completely (i.e., forcing them to be 0)

shrinkage operates by

- including a penalty term directly in the model equation and
- favoring small parameter values in general.

Shrinkage / Ridge Regression [?]



Ridge regression: minimize

$$\begin{split} \mathsf{RSS}_{\lambda}(\hat{\beta}) = &\mathsf{RSS}(\hat{\beta}) + \lambda \sum_{j=1}^{p} \hat{\beta}_{j}^{2} \\ = &\langle \mathbf{y} - \mathbf{X}\hat{\beta}, \mathbf{y} - \mathbf{X}\hat{\beta} \rangle + \lambda \sum_{j=1}^{p} \hat{\beta}_{j}^{2} \\ \Rightarrow \hat{\beta} = &(\mathbf{X}^{T}\mathbf{X} + \lambda I)^{-1}\mathbf{X}^{T}\mathbf{y} \end{split}$$

with $\lambda \ge 0$ a complexity parameter / regularization parameter.

As

- solutions of ridge regression are not equivariant under scaling of the predictors, and as
- it does not make sense to include a constraint for the parameter of the intercept

data is normalized before ridge regression:

$$x_{i,j}' := \frac{x_{i,j} - \bar{x}_{.,j}}{\hat{\sigma}(x_{.,j})}$$

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Machine Learning / 5. Model Selection

Shrinkage / Ridge Regression (2/3)



Ridge regression is a combination of

$$\sum_{i=1}^{n} (y_i - \hat{y}_i)^2 + \lambda \sum_{j=1}^{p} \beta_j^2$$

$$=$$
 L2 loss $+\lambda$ L2 regularization

Shrinkage / Ridge Regression (3/3) / Tikhonov Regularization (1/2)

L2 regularization / Tikhonov regularization can be derived for linear regression as follows:

Treat the true parameters θ_j as random variables Θ_j with the following distribution (**prior**):

$$\Theta_j \sim \mathcal{N}(0, \sigma_{\Theta}), \quad j = 1, \dots, p$$

Then the joint likelihood of the data and the parameters is

$$L_{\mathcal{D},\Theta}(\theta) := \left(\prod_{i=1}^{n} p(x_i, y_i \mid \theta)\right) \prod_{j=1}^{p} p(\Theta_j = \theta_j)$$

and the conditional joint log likelihood of the data and the parameters accordingly

$$\log L_{\mathcal{D},\Theta}^{\mathsf{cond}}(\theta) := \left(\sum_{i=1}^{n} \log p(y_i \mid x_i, \theta)\right) + \sum_{j=1}^{p} \log p(\Theta_j = \theta_j)$$

and

$$\log p(\Theta_j = \theta_j) = \log \frac{1}{\sqrt{2\pi}\sigma_{\Theta}} e^{-\frac{\theta_j^2}{2\sigma_{\Theta}^2}} = -\log(\sqrt{2\pi}\sigma_{\Theta}) - \frac{\theta_j^2}{2\sigma_{\Theta}^2}$$

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Machine Learning / 5. Model Selection

Shrinkage / Ridge Regression (3/3) / Tikhonov Regularization (2/2)

Dropping the terms that do not depend on θ_j yields:

$$\log L_{\mathcal{D},\Theta}^{\mathsf{cond}}(\theta) := \left(\sum_{i=1}^{n} \log p(y_i \mid x_i, \theta)\right) + \sum_{j=1}^{p} \log p(\Theta_j = \theta_j)$$
$$\propto \left(\sum_{i=1}^{n} \log p(y_i \mid x_i, \theta)\right) - \frac{1}{2\sigma_{\Theta}^2} \sum_{j=1}^{p} \theta_j^2$$

This also gives a semantics to the complexity / regularization parameter λ :

$$\lambda = \frac{1}{2\sigma_{\Theta}^2}$$

but σ_{Θ}^2 is unknown. (We will see methods to estimate λ later on.)

The parameters θ that maximize the joint likelihood of the data and the parameters are called **Maximum Aposteriori Estimators (MAP estimators)**.

Putting a prior on the parameters is called Bayesian approach.





 $Y = \beta_0 + \beta_1 X_1 + \beta_2 X_2 + \epsilon$

How to compute ridge regression / Example



Fit

to the data:

x_1	x_2	y
1	2	3
2	3	2
4	1	7
5	5	1

$$X = \begin{pmatrix} 1 & 1 & 2 \\ 1 & 2 & 3 \\ 1 & 4 & 1 \\ 1 & 5 & 5 \end{pmatrix}, \quad Y = \begin{pmatrix} 3 \\ 2 \\ 7 \\ 1 \end{pmatrix}, \qquad I := \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix},$$
$$X^{T}X = \begin{pmatrix} 4 & 12 & 11 \\ 12 & 46 & 37 \\ 11 & 37 & 39 \end{pmatrix}, \quad X^{T}X + 5I = \begin{pmatrix} 9 & 12 & 11 \\ 12 & 51 & 37 \\ 11 & 37 & 44 \end{pmatrix}, \quad X^{T}Y = \begin{pmatrix} 13 \\ 40 \\ 24 \end{pmatrix}$$

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Machine Learning



- 1. The Regression Problem
- 2. Simple Linear Regression
- 3. Multiple Regression
- 4. Variable Interactions
- 5. Model Selection
- 6. Case Weights

Cases of Different Importance

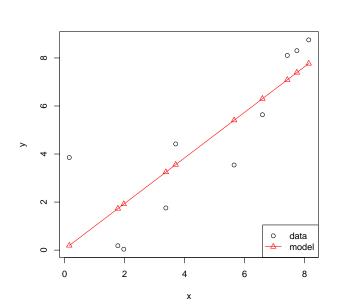
Young 2003

Sometimes different cases are of different importance, e.g., if their measurements are of different accurracy or reliability.

Example: assume the left most point is known to be measured with lower reliability.

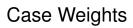
Thus, the model does not need to fit to this point equally as well as it needs to do to the other points.

I.e., residuals of this point should get lower weight than the others.



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Machine Learning / 6. Case Weights





In such situations, each case (x_i, y_i) is assigned a **case weight** $w_i \ge 0$:

- the higher the weight, the more important the case.
- cases with weight 0 should be treated as if they have been discarded from the data set.

Case weights can be managed as an additional pseudo-variable w in applications.

Weighted Least Squares Estimates

Formally, one tries to minimize the **weighted residual sum of** squares

$$\sum_{i=1}^{n} w_i (y_i - \hat{y}_i)^2 = ||\mathbf{W}^{\frac{1}{2}}(\mathbf{y} - \hat{\mathbf{y}})||^2$$

with

$\mathbf{W} := \begin{pmatrix} w_1 & & 0 \\ & w_2 & & \\ & & \ddots & \\ 0 & & & w_n \end{pmatrix}$

The same argument as for the unweighted case results in the **weighted least squares estimates**

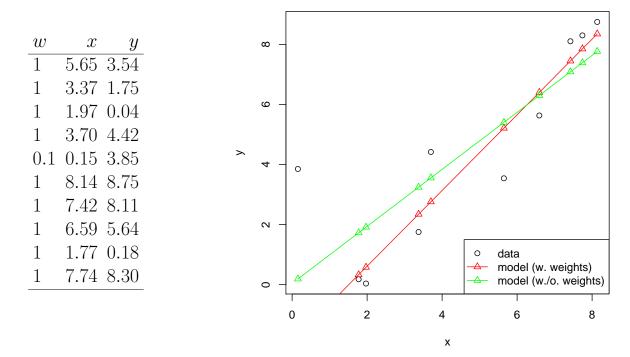
$$\mathbf{X}^T \mathbf{W} \mathbf{X} \hat{\boldsymbol{\beta}} = \mathbf{X}^T \mathbf{W} \mathbf{y}$$

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Machine Learning / 6. Case Weights

Weighted Least Squares Estimates / Example

To downweight the left most point, we assign case weights as follows:



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Summary



- For regression, **linear models** of type $Y = \langle X, \beta \rangle + \epsilon$ can be used to predict a quantitative *Y* based on several (quantitative) *X*.
- The ordinary least squares estimates (OLS) are the parameters with minimal residual sum of squares (RSS). They coincide with the maximum likelihood estimates (MLE).
- OLS estimates can be computed by solving the system of linear equations X^TXβ = X^TY.
- The variance of the OLS estimates can be computed likewise $((\mathbf{X}^T\mathbf{X})^{-1}\hat{\sigma}^2)$.
- For deciding about inclusion of predictors as well as of powers and interactions of predictors in a model, model selection measures (AIC, BIC) and different search strategies such as forward and backward search are available.

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