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# STAT 444/844 Course Notes

STATISTICAL LEARNING: FUNCTION ESTIMATION

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# **Table of Contents**

| 1 | January 8, 2019                                  |    |  |  |  |  |  |  |  |
|---|--|----|--|--|--|--|--|--|--|
|   | 1.1 What is a function?                          |    |  |  |  |  |  |  |  |
|   | 1.2 Advertising data example                     |    |  |  |  |  |  |  |  |
|   | 1.3 Notation                                     |    |  |  |  |  |  |  |  |
|   | 1.4 Definitions and properties                   | 4  |  |  |  |  |  |  |  |
| 2 | January 10, 2019                                 | Ę  |  |  |  |  |  |  |  |
|   | .1 Linear models                                 |    |  |  |  |  |  |  |  |
|   | 2.2 Piecewise linear                             | (  |  |  |  |  |  |  |  |
|   | 2.3 Piecewise quadratic                          | ,  |  |  |  |  |  |  |  |
|   | 2.4 Weighted least squares                       | ,  |  |  |  |  |  |  |  |
| 3 | January 15, 2019                                 | 8  |  |  |  |  |  |  |  |
|   | 3.1 Weight least squares applications            | 8  |  |  |  |  |  |  |  |
|   | 3.2 Types of errors                              | 8  |  |  |  |  |  |  |  |
| 4 |  | _  |  |  |  |  |  |  |  |
|   | January 17, 2019                                 | ,  |  |  |  |  |  |  |  |
|   | 4.1 Notes on terminology and 1m in R             | (  |  |  |  |  |  |  |  |
|   | 4.2 Notes on model selection                     | (  |  |  |  |  |  |  |  |
|   | 4.3 Geometric interpretation of linear models    | 10 |  |  |  |  |  |  |  |
| 5 | January 24, 2019                                 | 10 |  |  |  |  |  |  |  |
|   | 5.1 Discrepancy function                         | 1( |  |  |  |  |  |  |  |
|   | 5.2 Discrepancy function and log-likelihood      | 1  |  |  |  |  |  |  |  |
|   | 5.3 Iteratively re-weighted least squares (IRLS) | 1  |  |  |  |  |  |  |  |
|   | 5.4 Why IRLS?                                    | 12 |  |  |  |  |  |  |  |
|   | Robust regression                                | 12 |  |  |  |  |  |  |  |
| 6 | January 29, 2019                                 | 14 |  |  |  |  |  |  |  |
|   | 6.1 Remark on robust regression and constants    | 14 |  |  |  |  |  |  |  |
|   | 6.2 Sensitivity curve and breakdown point        | 14 |  |  |  |  |  |  |  |
|   | 6.3 Least median squares (LMS)                   | 15 |  |  |  |  |  |  |  |
|   | 6.4 Least trimmed average sum of squares (LTS)   | 16 |  |  |  |  |  |  |  |

#### Abstract

These notes are intended as a resource for myself; past, present, or future students of this course, and anyone interested in the material. The goal is to provide an end-to-end resource that covers all material discussed in the course displayed in an organized manner. These notes are my interpretation and transcription of the content covered in lectures. The instructor has not verified or confirmed the accuracy of these notes, and any discrepancies, misunderstandings, typos, etc. as these notes relate to course's content is not the responsibility of the instructor. If you spot any errors or would like to contribute, please contact me directly.

# 1 January 8, 2019

## 1.1 What is a function?

Suppose we have some measured **response** variate y and we have one or more **explanatory** variables  $x_1, \ldots, x_p$ . The response and explanatory ariables are approximately related through an unknown function  $\mu(x)$  (to be estimated/learned) where

$$y = \mu(x) + r$$

where r is residual that cannot be explained by  $\mu(x)$ .

Some other names for response and explanatory variables include:

| response   | explanatory |  |  |
|------------|-------------|--|--|
| response   | predictor   |  |  |
| response   | design      |  |  |
| output     | input       |  |  |
| dependent  | independent |  |  |
| endogenous | exogenous   |  |  |

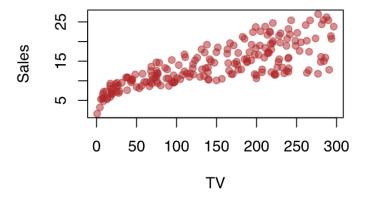
#### 1.2 Advertising data example

Suppose we want to predict Sales (response) from how much companies spend on TV, Radio, and Newspaper advertising (explanatory).

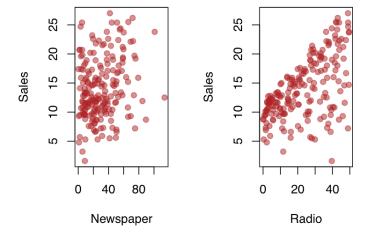
The dataset is

| ## |   | X | TV    | Radio | Newspaper | Sales |
|----|---|---|-------|-------|-----------|-------|
| ## | 1 | 1 | 230.1 | 37.8  | 69.2      | 22.1  |
| ## | 2 | 2 | 44.5  | 39.3  | 45.1      | 10.4  |
| ## | 3 | 3 | 17.2  | 45.9  | 69.3      | 9.3   |
| ## | 4 | 4 | 151.5 | 41.3  | 58.5      | 18.5  |
| ## | 5 | 5 | 180.8 | 10.8  | 58.4      | 12.9  |
| ## | 6 | 6 | 8.7   | 48.9  | 75.0      | 7.2   |

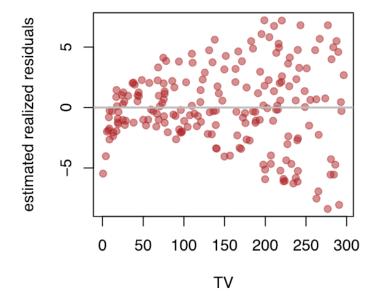
if we plot sales against TV



we see there is some positive correlation. Similarly against Newspaper and Radio

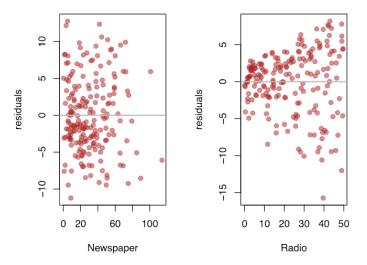


What if we tried a simple linear model where  $\hat{\mu}(x_1) = \hat{\alpha} + \hat{\beta}x_1$  where  $x_1$  is the TV advertising? We obtain estimates  $\hat{\alpha} = 7.03$  and  $\hat{\beta} = 0.05$  which are interpretable. However if we take a look at the residuals



we see that the residuals are not independently distributed accordingly to  $x_1$ , which violates our Markov-Gaussian assumptions.

The residuals of the model with Newspaper and Radio are



we observe that we do not observe constant variance across the explanatory variables.

Therefore a linear model does not seem to work (we could of course introduce scaling e.g. log-scaling for the Radio variate or polynomial terms).

### 1.3 Notation

Some notes on notation:

- Capital letters are matrices or vectors:  $A, X, \Sigma$
- Lower letters are scalars:  $a, x, \sigma$

- Arrows on letters are vectors:  $\vec{a}, \vec{x}$
- All vectors are column vectors
- The transpose of any matrix A is  $A^T$  (ocassionally A')

## 1.4 Definitions and properties

Quadratic form Suppose  $A = (a_{ij})_{n \times n}$  is symmetric i.e.  $a_{ij} = a_{ji} \ \forall i, j$ . Then

$$f = Y^T A Y$$
$$= \sum_{i} \sum_{j} a_{ij} y_i y_j$$

is called a quadratic form.

**Trace** For a matrix, the **trace** is defined as

$$\operatorname{tr}(A_{m \times m}) = \sum_{i=1}^{m} a_{ii}$$

Note that tr(BC) = tr(CB).

**Rank** The **rank** of a matrix denoted rank(A) is the maximum number of *linearly independent columns* (or rows) of A.

Note that vectors  $Y_1, \ldots, Y_n$  are linearly independent iff

$$c_1Y_1 + \ldots + c_nY_n = 0$$

implies  $c_1 = \ldots = c_n = 0$  (i.e. no non-trivial solution).

Eigenvector and eigenvalue A non-zero vector  $\vec{v}_i$  is an eigenvector of  $A_{m\times m}$  if

$$A\vec{v}_i = \lambda_i \vec{v}_i$$
  $i = 1, 2, \dots, m$ 

where  $\lambda_i$  is the corresponding *i*th **eigenvalue**.

**Idempotent** A matrix A is **idempotent** if AA = A.

Some notable results:

- 1. If A is idempotent, then all its eigenvalues are either 0 or 1.
- 2. If A is idempotent, there exists an orthogonal matrix P such that  $A = P\Lambda P^T$  where

$$\Lambda = \begin{bmatrix} 1 & \dots & 0 & 0 & 0 & 0 \\ 0 & \ddots & 0 & 0 & 0 & 0 \\ 0 & \dots & 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & \dots & 0 \\ 0 & 0 & 0 & 0 & \ddots & 0 \\ 0 & 0 & 0 & 0 & \dots & 0 \end{bmatrix}$$

and  $tr(A) = rank(A) = tr(\Lambda)$  which is equivalent to the number of eigenvalues being 1.

# 2 January 10, 2019

#### 2.1 Linear models

A linear model is generally in the form of

$$y_i = \beta_0 + \beta_1 x_{i1} + \ldots + \beta_p x_{ip} + \epsilon_i \qquad i = 1, \ldots, n$$

which holds under the assumptions that

- $E(\epsilon_i) = 0$
- $Var(\epsilon_i) = \sigma^2$  (constant variance)
- $\epsilon_1, \ldots, \epsilon_n$  are independent
- $\epsilon_1, \ldots, \epsilon_n \stackrel{iid}{\sim} N(0, \sigma^2)$

In matrix form we have

$$\begin{bmatrix} y_1 \\ \vdots \\ y_n \end{bmatrix}_{n \times 1} = \begin{bmatrix} 1 & x_{11} & \dots & x_{1p} \\ \vdots & \vdots & \ddots & \vdots \\ 1 & x_{n1} & \dots & x_{np} \end{bmatrix}_{n \times (p+1)} \begin{bmatrix} \beta_0 \\ \beta_1 \\ \vdots \\ \beta_p \end{bmatrix}_{(p+1) \times 1} + \begin{bmatrix} \epsilon_1 \\ \vdots \\ \epsilon_n \end{bmatrix}_{n \times 1}$$

or in short matrix form  $Y = X\vec{\beta} + \vec{\epsilon}$ .

The Least Squares Estimator (LSE) of  $\vec{\beta}$  minimizing the discrepancy function

$$S(\vec{\beta}) = (Y - X\vec{\beta})^T (Y - X\vec{\beta})$$

has a closed form solution

$$\hat{\vec{\beta}} = (X^T X)^{-1} X^T Y$$

The fitted values are thus

$$\hat{Y} = X\hat{\vec{\beta}} = X(X^TX)^{-1}X^TY$$
$$= HY$$

where  $H = X(X^TX)^{-1}X^T$  (hat matrix). Note that H is idempotent and symmetric.

Geometric interpretation of LSE:  $\hat{Y}$  is the projection of Y onto C(X), the column space of X (we can thus see that the fitted errors should be orthogonal to our fitted values in LSE).

The **degrees of freedom** of our model is n - (p + 1) where p + 1 is the number of free parameters in our model. This is equivalent to  $n - \operatorname{tr}(H)$  i.e.  $\operatorname{tr}(H) = p + 1$ . Under normality

- $\hat{\vec{\beta}} = MVN(\vec{\beta}, \sigma^2(X^TX)^{-1})$
- $\hat{\vec{\beta}}$  and  $\hat{\sigma}^2$  are independent (Note  $\hat{\sigma}^2 = \frac{SSE}{df}$ ).
- $\bullet \ \frac{(n-p-1)\hat{\sigma}^2}{\sigma^2} \sim \chi^2_{n-p-1}$

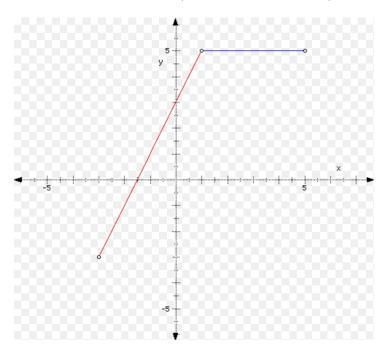
Let  $\vec{a}_p = (1, x_1, \dots, x_p)^T$  (observation  $\vec{x}$  extended with intercept term). The  $(1 - \alpha)$  prediction interval at  $\vec{a}_p$  is

$$\vec{a}_p^T \hat{\vec{\beta}} \pm t_{n-p-1,\alpha/2} \hat{\sigma} \sqrt{1 + \vec{a}_p^T (X^T X)^{-1} \vec{a}_p}$$

We can also estimate **confidence intervals** as well (drop  $1 + \dots$  term above).

#### 2.2 Piecewise linear

We can specify the following piecewise linear function (with discontinuity at a)



as two linear functions

$$y = \begin{cases} \beta_0 + \beta_1 x & x \le a \\ \beta_2 + \beta_3 x & x \ge a \end{cases}$$

subject to  $\beta_0 + \beta_1 a = \beta_2 + \beta_3 a$ .

A more convenient way to express the above

$$y = \beta_0 + \beta_1 x + \beta_2 (x - a) I(x > a)$$

where I is the indicator function. Note the above is linear in terms of  $\vec{\beta}$  BUT NOT in terms of x. However we can simply construct a new variate  $(x-a)I(x \ge a)$  from x.

Note that  $\beta_2$  is the *change in slope right of a* for samples where  $x \geq a$ .

Extension to more than one interesting point (knot) is straightforward.

## 2.3 Piecewise quadratic

Similar to piecewise linear models, we can specify

$$y = \begin{cases} \beta_0 + \beta_1 x + \beta_2 x^2 & x \le a \\ \beta_3 + \beta_4 x + \beta_5 x^2 & x \ge a \end{cases}$$

subject to  $\beta_0 + \beta_1 a + \beta_2 a^2 = \beta_3 + \beta_4 a + \beta_5 a^2$  (continuity) and  $\beta_1 + 2\beta_2 a = \beta_4 + 2\beta_5 a$  (differentiable at a). Alternatively we can express this as one linear function

$$y = \beta_0 + \beta_1 x + \beta_2 x^2 + \beta_3 (x - a)^2 I(x \ge a)$$

continuity is trivially satisfied. Note the 1st derivative is

$$\frac{dy}{dx} = \beta_1 + \beta_2 x + 2\beta_3 (x - a)I(x \ge a)$$

where the last term is 0 when x = a, thus our additional indicator term does not affect the derivative.

## 2.4 Weighted least squares

Sometimes we would like to give more importance to some observations than others. Instead of minimizing  $(Y - X\vec{\beta})^T (Y - X\vec{\beta})$  we can minimize

$$(Y - X\vec{\beta})^T W (Y - X\vec{\beta})$$

where

$$W = \begin{bmatrix} w_1 & 0 & \dots & 0 \\ 0 & w_2 & \dots & 0 \\ 0 & 0 & \ddots & 0 \\ 0 & 0 & \dots & w_n \end{bmatrix}_{n \times n}$$

a diagonal matrix.  $w_i$  corresponds to the weight assigned to observation i (a higher  $w_i$  the more important that observation is).

Claim. The closed form solution is

$$\hat{\vec{\beta}}_{WLS} = (X^T W X)^{-1} X^T W Y$$

Proof. Note that

$$S(\vec{\beta}) = (Y - X\vec{\beta})^T W (Y - X\vec{\beta})$$
  
=  $Y^T W Y - Y^T W X \vec{\beta} - \vec{\beta}^T X^T W Y + \vec{\beta}^T X^T W X \vec{\beta}$ 

Note that  $Y^TWX\vec{\beta}=(\vec{\beta}^TX^TWY)^T$  which is a scalar, so  $Y^TWX\vec{\beta}=\vec{\beta}^TX^TWY$  (transposes of scalars are equivalent). thus

$$S(\vec{\beta}) = Y^T W Y - 2Y^T W X \vec{\beta} + \vec{\beta}^T X^T W X \vec{\beta}$$

where  $-2Y^TWX\vec{\beta}$  is the "linear term" and  $\vec{\beta}^TX^TWX\vec{\beta}$  is of quadratic form.

Recall that

$$\frac{d\vec{c}^T Y}{dY} = \vec{c}^T$$

$$\frac{dY^T AY}{dY} = 2Y^T A^T$$

SO

$$\frac{dS(\vec{\beta})}{d\vec{\beta}} = -2Y^T W X + 2\vec{\beta}^T X^T W X$$

$$\Rightarrow \vec{\beta}^T X^T W X = Y^T W X$$

$$\Rightarrow (X^T W X) \vec{\beta} = X^T W Y$$

$$\Rightarrow \vec{\beta} = (X^T W X)^{-1} X^T W Y$$

$$\frac{dS(\vec{\beta})}{d\vec{\beta}} = 0$$

$$W^T = W$$

as claimed.  $\Box$ 

Here is an alternative proof:

*Proof.* Let  $Y^* = W^{\frac{1}{2}}Y$  and  $X^* = W^{\frac{1}{2}}X$ .

Note that minimizing  $(Y - X\vec{\beta})^T W(Y - X\vec{\beta})$  is equivalent to minimizing  $(Y^* - X^*\vec{\beta})^T (Y^* - X^*\vec{\beta})$  (simply expand out  $X^*$  and  $Y^*$ ).

Thus the LSE of  $\vec{\beta}$  with  $X^*, Y^*$  is

$$\vec{\beta} = (X^{*^T}X)^{-1}X^{*^T}Y^*$$

$$= ((X^TW^{\frac{1}{2}})(W^{\frac{1}{2}}X))^{-1}(X^TW^{\frac{1}{2}})(W^{\frac{1}{2}}Y)$$

$$= (X^TWX)^{-1}X^TWY$$

which is equivalent to our previous derivation.

# 3 January 15, 2019

#### 3.1 Weight least squares applications

**Example 3.1.** We can apply weighted least squares to do **local regression** where we downweight observations farther away from a given observation.

**Example 3.2.** Suppose that  $Var(\epsilon_i) = \sigma_i^2$  (i.e. not all observations are drawn with the same variance). If we want to overweight observations that have lower variance, we can set  $w_i = \frac{1}{\sigma_i^2}$  to obtain an **unbiased estimator** of  $\vec{\beta}$  with the *smallest variance* (**Best Linear Unbiased Estimator** or **BLUE**).

#### 3.2 Types of errors

We will use the example of 790 Facebook posts published by a comestic company to illustrate.

The population being examined is called the **study population** (the 790 posts). The analysis of these posts may be applied to a larger population (whether it's future Facebook posts for this company or Facebook posts for *any* company) which we call the **targt population**.

The difference between the study and target population is called the **study error**.

In a paper by Soros et al., they ended up using a **sample** of only 500 posts for confidentiality reasons. The difference between the sample and study population is called **sample error**.

# 4 January 17, 2019

## 4.1 Notes on terminology and 1m in R

- When using 1m the intercept term is included by default. To remove it simply specify  $Y \sim X 1$ .
- Factors are like categorical variables in R: there are a finite number of categories (called factor levels).
- In 1m almost any function of variates may appear in the formula e.g.  $Y \sim X + \sin(X)$  or  $Y \sim X + \sin(X) + \sin(X)$

To specify  $Y = X \cdot Z$ , we need to use  $Y \sim I(X \star Z)$  or  $Y \sim X:Z$  instead of  $X \star Z$  since  $X \star Z$  represents interaction in 1m and translates to the model  $y = \alpha x + \beta z + \gamma xz + r$ .

• Some arithmetic operations e.g. +, -, \*, ^ are interpreted as formula operators rather than arithmetic operators in lm. One should wrap them in I(·).

#### 4.2 Notes on model selection

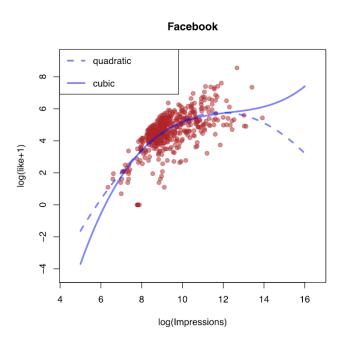


Figure 4.1: Quadratic and cubic polynomial linear models on Facebook data.

In the above figure we see that while both quadratic and cubic models are **global** (predict any value of x) the quadratic model seems to predict likes returning to 0 as impressions approach infinity.

The cubic function on the contrary continues to increase: this makes more sense intuitively, thus examining a model often requires human understanding of the data and problem.

## 4.3 Geometric interpretation of linear models

A linear model is a linear combination of functions called **generators** e.g.

$$\mu(x) = \beta_1 g_1(x) + \beta_2 g_2(x) + \beta_3 g_3(x) + \beta_4 g_4(x)$$

where  $g_1, \ldots, g_4$  could be arbitrary continuous functions of x.

All possible linear combinations of the generators forms a **subspace** (the functions generate the subspace). The functions are a **basis** for this subspace.  $\mu(x)$  lies in the subspace whose dimension equals to the number of basis functions.

The functions should be *linearly independent* of each other: otherwise the solution to parameters will be ill-defined.

# 5 January 24, 2019

## 5.1 Discrepancy function

Let the **discrepancy function** for a fit of parameters  $\vec{\beta}$  be denoted

$$S(\vec{\beta}) = \sum_{i=1}^{n} \rho(y_i - \vec{x}_i^T \vec{\beta}) = \sum_{i=1}^{n} \rho(r_i)$$

where  $\rho$  is a real-valued loss function (in the OLS case, this was simply the square function).  $r_i$  is our residual for observation i.

Taking the derivative

$$\frac{dS(\vec{\beta})}{d\vec{\beta}} = \sum_{i=1}^{n} \rho'(y_i - \vec{x}_i^T \vec{\beta})(-1)\vec{x}_i^T$$
$$= -\sum_{i=1}^{n} \rho'(r_i)\vec{x}_i^T$$

Solving for the extremum we have

$$\sum_{i=1}^{n} \phi(r_i) \vec{x}_i^T = \vec{0}^T$$

where  $\phi(r) = \rho'(r)$  (derivative with respect to  $\vec{\beta}$ ).

**Remark 5.1** (LSE). If  $\rho(r) = r^2$  we get LSE, that is  $(\phi(r) = 2r)$ 

$$\vec{0} = \sum_{i=1}^{n} 2r\vec{x}_i$$

$$= 2\sum_{i=1}^{n} (\vec{x}_i y_i - \vec{x}_i^T \vec{x}_i \vec{\beta})$$

$$= 2(X^T Y - X^T X \vec{\beta})$$

which exactly solves to our LSE closed form solution.

## 5.2 Discrepancy function and log-likelihood

Let us compare our discrepancy function with the log-likelihood for linear models:

$$l(\vec{\beta}) = \sum_{i=1}^{n} l_i(\vec{\beta})$$
$$= \sum_{i=1}^{n} l(r_i)$$

where  $l(r_i) = \frac{-r_i^2}{2\sigma^2}$ , a function only of  $r_i$ .

The second equality follows from the following remark:

**Remark 5.2.** Note  $l_i(\vec{\beta})$  is the *i*th observation's contribution to  $l(\vec{\beta})$  i.e.

$$l_i(\vec{\beta}) = \log f(y_i \mid \vec{x}_i^T \vec{\beta}, \sigma^2)$$

For a linear model we have

$$f(y_i \mid \vec{x}_i^T \vec{\beta}, \sigma^2) \sim N(\vec{x}_i^T \vec{\beta}, \sigma^2)$$

so  $l_i(\vec{\beta}) = -\frac{r_i^2}{2\sigma^2} + C$  where  $C = -\frac{1}{2}\log(2\pi) - \log\sigma$  a constant.

We let  $l(r_i) = -\frac{r_i^2}{2\sigma^2}$ . Since the constant does not change with respect to  $\beta$  we can omit it from our objective function.

From above we observe that minimizing the discrepancy function is the same as maximizing the log likelihood where  $\rho(r) = -l(r)$  in the discrepancy function.

**Definition 5.1** (M-estimator). We call the estimator  $\vec{\beta}$  that minimizes  $\sum_{i=1}^{n} \rho(r_i)$  the M-estimator or the maximum-likelihood type estimator.

## 5.3 Iteratively re-weighted least squares (IRLS)

Note that the solution turns out to be a WLS estimator:

$$\vec{0} = \sum_{i=1}^{n} \phi(r_i) \vec{x}_i$$

$$= \sum_{i=1}^{n} \frac{\phi(r_i)}{r_i} r_i \vec{x}_i$$

$$= \sum_{i=1}^{n} w(r_i) (y_i - \vec{x}_i^T \vec{\beta}) \vec{x}_i$$

$$= \sum_{i=1}^{n} w_i (y_i - \vec{x}_i^T \vec{\beta}) \vec{x}_i$$

where we let  $w_i = w(r_i) = \frac{\phi(r_i)}{r_i}$ . If we solve this we see that the solution is WLS where

$$\hat{\vec{\beta}} = (X^T W X)^{-1} X^T W Y$$

with  $W = \operatorname{diag}(w_1, \ldots, w_n)$ .

**However**, the weights of this WLS depend on the residuals which in turn depends on  $\vec{\beta}$ . If we are given an initial estimate of  $\vec{\beta}^{(0)}$ , we could *iteratively update residuals* and  $\vec{\beta}$  to converge to a solution. We proceed as follows:

nitialization Initialization: set j = 0

Step 1 Compute residuals

$$\hat{r}_i^{(j)} = y_i - \vec{x}_i^T \hat{\vec{\beta}}^{(j)}$$
  $i = 1, \dots, n$ 

Step 2 Update weights

$$w_i^{(j)} = \frac{\phi(\hat{r}_i^{(j)})}{\hat{r}_i^{(j)}}$$

and let  $W^{(j)} = \text{diag}(w_1^{(j)}, \dots, w_n^{(j)}).$ 

Step 3 WLS to estimate next set of  $\hat{\vec{\beta}}^{(j+1)}$ 

$$\hat{\vec{\beta}}^{(j+1)} = (X^T W^{(j)} X)^{-1} X^T W^{(j)} Y$$

Step 4 Set j = j + 1 and return to Step 1 if convergence criterion is not met.

We can this procedure iteratively re-weighted least squares (IRLS).

The convergence criterion is typically

$$\|\hat{\vec{\beta}}^{(j+1)} - \hat{\vec{\beta}}^{(j)}\| \le \epsilon$$

with the L2/Euclidean norm and for some small positive constant  $\epsilon$ .

#### 5.4 Why IRLS?

Question 5.1. Why do we need to use iteratively re-weighted least squares?

In ordinary least squares with Gaussian response and loss function  $r_i^2$ , there is no reason to use IRLS since OLS and IRLS are equivalent (the loss function  $r_i^2$  simplifies IRLS to OLS).

However in **generalized linear models (GLMs)** (STAT 431/831) we may have a different type of response (e.g. Bernoulli 0/1 or categorical) and thus we may define our loss function  $\rho(r_i)$  differently.

We may also want to modify our  $\rho(r_i)$  to de-emphasize huge outliers (see next section).

#### 5.5 Robust regression

Robust regression tries to de-emphasize the influence of large outliers.

**Question 5.2.** What loss function  $\rho$  (and  $\phi$ ) should we use?

In ordinary least squares (OLS) we can use

$$\rho(r) = \frac{1}{2}r^2$$

$$\phi(r) = r$$

$$w(r) = \frac{\phi(r)}{r} = 1$$

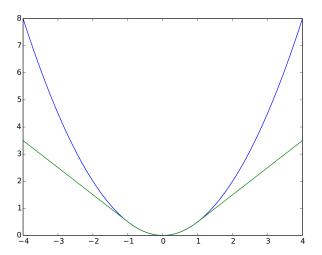
so we essentially have  $W = I_{n \times n}$  which devolves into OLS as expected.

**Remark 5.3.** The residual function for OLS is unbounded and so extreme outliers with large residuals have significantly more influence.

Huber (1964) proposed a modified loss function (**Huber loss**) which de-emphasizes outliers:

$$\rho(r) = \begin{cases} \frac{1}{2}r^2 & \text{if } |r| \le c\\ c(|r| - \frac{1}{2}c) & \text{if } |c| > c \end{cases}$$

The modified loss function essentially makes the loss function linear after a certain threshold c:



We also let

$$\phi(r) = \begin{cases} r & \text{if } |r| \le c \\ c \text{sign}(r) & \text{if } |c| > c \end{cases}$$

and thus

$$w(r) = \begin{cases} 1 & \text{if } |r| \le c \\ \frac{c}{|r|} & \text{if } |c| > c \end{cases}$$

The  $\phi$  and weight w functions look like



Figure 5.1: Left:  $\phi(r)$ . Right: w(r) for Huber's loss function.

How do we decide c? Huber suggested c = 1.345 and showed it achieved 95% of LSE asymptotically when the true distribution is normal (95% efficiency essentially means the true variance of the betas from OLS is 95% that of the variance of the betas using Huber's loss).

Question 5.3. Since c is fixed, what if our residuals are scaled to very large or small values (e.g. O(1e5) or O(1e-4))? We would have to scale our data beforehand to make it within a sensible range so that c = 1.345 makes sense.

Sometimes we prefer the  $\phi$  function to "redescend" i.e.  $\phi(r) \to 0$  when |r| is large (that is: we fully de-emphasize outliers). Other  $\phi$  functions include

#### Redescending M-estimator (Hampel)

$$\phi(r) = \begin{cases} r & \text{if } 0 \le |r| \le a \\ a \text{sign}(r) & \text{if } a \le |r| \le b \\ a \frac{c - |r|}{c - b} \text{sign}(r) & \text{if } b \le |r| \le c \\ 0 & \text{if } |r| > c \end{cases}$$

The recommended settings are a = 2, b = 4, c = 8 (with appropriately scaled data and residuals).

#### Tukey's biweight

$$\phi(r) = \begin{cases} r \left(1 - \left(\frac{r}{c}\right)^2\right)^2 & \text{if } |r| \le c \\ 0 & \text{if } |r| > c \end{cases}$$

where c = 4.685 is typically used. This is designed to have 95% efficiency as well for a true normal distribution.

# 6 January 29, 2019

#### 6.1 Remark on robust regression and constants

**Remark 6.1.** All recommended constants in the various robust regression methods (Huber, Hampel, Tukey) are based on the assumption that Var(r) = 1. Therefore in practice we typically need to scale the residuals i.e.  $r'_i = \frac{r_i}{s}$  where s is a scale parameter.

One simple solution is to estimate the **median absolute deviation (MAD)**:

$$MAD = median(|r_i|)$$

and let  $\hat{s} = \frac{\text{MAD}}{0.6745}$ . For the standard normal distribution we note that MAD = 0.6745.

#### 6.2 Sensitivity curve and breakdown point

Let  $T_n(y_1, \ldots, y_n)$  be a population attribute (that is a function of the same points). To see how **sensitive**  $T_n$  is to an individual data point, define

$$SC(y) = \frac{T_n(y_1, \dots, y_{n-1}, y) - T_{n-1}(y_1, \dots, y_{n-1})}{\frac{1}{n}}$$

which is the difference between  $T_n(\cdot)$  (with all n points) and  $T_{n-1}(\cdot)$  (with one point y omitted) compare to the contamination size  $\frac{1}{n}$ .

**Example 6.1.** Let  $T_n(y_1, \ldots, y_n) = \frac{1}{n} \sum_{i=1}^n y_i = \bar{y}_n$  (sample mean). Note that

$$T_n = \sum_{i=1}^{n-1} y_i + y = \frac{n-1}{n} \bar{y}_{n-1} + y$$

Note that SC(y) is simply

$$SC(y) = n(T_n - T_{n-1}) = (n-1)\bar{y}_{n-1} + y - n\bar{y}_{n-1}$$
  
=  $y - \bar{y}_{n-1}$ 

**Definition 6.1** (Breakdown point). *Informally*, the **breakdown point** of a statistic is the largest proportion of contamination before the statistic breaks down.

Formally, let  $\vec{z}_i = (x_{i1}, x_{i2}, \dots, x_{ip}, y_i)^T$  for  $i = 1, \dots, n$  be the *i*th data vector.

Let  $Z = (\vec{z_1}, \dots, \vec{z_n})$  be the whole set. Let T be the statistic of intrest. The worst error for swapping  $m \ z_i$ 's is

$$e(m;T,Z) = \sup_{Z_m^*} \lVert T(Z_m^*) - T(Z) \rVert$$

where  $Z_m^*$  is Z with any of its m data vectors replaced.

The **breakdown point** is then defined as

$$\min\left\{\frac{m}{n}\mid e(m;T,Z)=\infty\right\}$$

**Remark 6.2.** That is: the breakdown point measures the minimum **proportion** of points required to influence the statistic *significantly*.

Some breakdown point examples:

**Sample mean** Note we can simply swap out m=1 point arbitrarily such that  $e(1;T,Z)\to\infty$  thus the breakdown point is  $\frac{1}{n}\to 0$  as  $n\to\infty$ .

**Median** The breakdown point is  $\frac{1}{2}$  as  $n \to \infty$ : we need to change at least half of them to aribitrarily influence the median e.g. make it go to infinity.

k% trimmed mean The k% trimmed mean is defined as the mean after discarding the lowest k% and highest k% of  $y_i$ 's.

Breakdown point is k% (we swap out the top k% + 1 points).

### 6.3 Least median squares (LMS)

Recall for regression, the LSE of  $\vec{\beta}$  is

$$\operatorname{argmin}_{\vec{\beta}} \sum_{i=1}^{n} (y_i \vec{x}_i^T \vec{\beta})^2$$

or equivalently

$$\operatorname{argmin}_{\vec{\beta}} \operatorname{average}(y_i \vec{x}_i^T \vec{\beta})^2$$

To make it robust for "outliers" or contaminations i.e. to ensure we have a *high breakdown point* we could consider the **least median squares (LMS) estimator**:

$$\vec{\beta}_{LMS} = \operatorname{argmin}_{\vec{\beta}} \operatorname{median}(y_i \vec{x}_i^T \vec{\beta})^2$$

which has a breakdown point of  $\frac{1}{2}$  (compared to a breakdown point of  $\frac{1}{n}$  for OLS).

## 6.4 Least trimmed average sum of squares (LTS)

Similar to how we made our objective function for OLS more robust by considering the median in LMS, we can also consider the (least) trimmed average sum of squares (LTS) estimator:

$$\vec{\beta}_{LTS} = \operatorname{argmin}_{\vec{\beta}} \sum_{i=1}^{k} r_{(i)}^2$$

where  $r_{(i)}^2$  is the *i*th smallest squared residual.

Note the breakdown point for LTS is  $\frac{n-k+1}{n}$  (compared to a breakdown point of  $\frac{1}{n}$  for OLS).