

Robust Linear Regression by Subquantile Minimization

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Presentation Overview

① Robust Regression

Huber Contamination Model

Blocks

Preliminaries

② Empirical Results

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③ Mathematics

Huber Contamination Model

Problem

The *Huber Contamination Model* is the following:

$$\hat{P} = (1 - \varepsilon)P + \varepsilon Q \text{ where } \varepsilon \in (0, 0.5)$$

where P and Q represent the general linear models

$$\mathbf{y}_P = \boldsymbol{\beta}_P^\top \mathbf{P} + \epsilon_P$$

$$\mathbf{y}_Q = \boldsymbol{\beta}_Q^\top \mathbf{Q} + \epsilon_Q$$

$\boldsymbol{\beta}_P$ and $\boldsymbol{\beta}_Q$ are oracle regressors and ϵ_P and ϵ_Q represent 0-centered gaussian noise.

Our goal is to learn a model that learns a good distribution of P from \hat{P}

Motivation

Definition

Oblivious Noise is noise which is independent of the input data

Adaptive Noise is noise which is dependent upon the input data

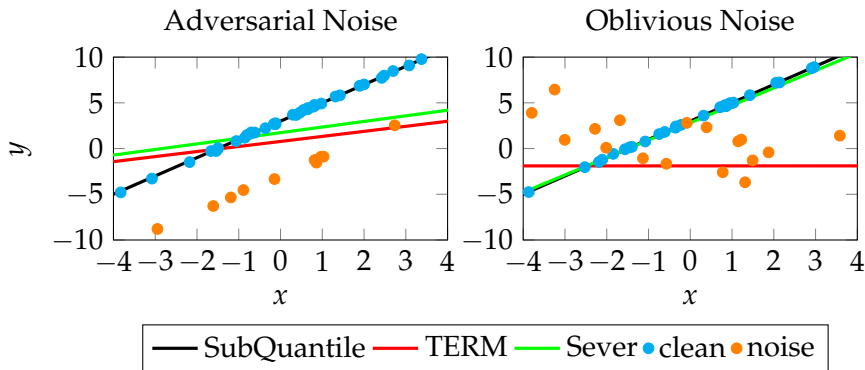


Figure: Sub-Quantile Performance on Adaptive Outliers

Theorem

The expected optimal parameters of the corrupted model \hat{P}

$$\mathbb{E} \left[\mathbf{X}^\dagger \mathbf{y} \right] = (1 - \varepsilon) \boldsymbol{\beta}_P + \varepsilon \boldsymbol{\beta}_Q$$

where $\mathbf{X}^\dagger \triangleq (\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top$, i.e. the Moore-Penrose Inverse.

This theorem motivates our reasoning for optimizing over the Subquantile. We want a method to reduce ε .

Statistical Preliminaries of the Subquantile

- ① The quantile is given as the following:

$$\mathcal{Q}_p = \inf \{x \in \mathbb{R} : p \leq F(x)\}$$

- ② Let ℓ be the loss functions. We can now define risk as:

$$\mathcal{U} = \mathbb{E} [\ell(f(\mathbf{x}; \boldsymbol{\theta}), y)]$$

- ③ The p -Quantile of the Empirical Risk is given by:

$$\mathbb{L}_p = \frac{1}{p} \int_0^p \mathcal{Q}_q(\mathcal{U}) dq = \mathbb{E} [\mathcal{U} | \mathcal{U} \leq \mathcal{Q}_p(\mathcal{U})] = \max_{t \in \mathbb{R}} \left\{ t - \frac{1}{p} \mathbb{E} [(t - \mathcal{U})^+] \right\}$$

- ④ For the least squares regression case:

$$\mathbb{L}_p = \max_{t \in \mathbb{R}} \left\{ t - \frac{1}{np} \sum_{i=1}^n \left(t - \left(\boldsymbol{\theta}^\top \mathbf{x}_i - y_i \right) \right)^+ \right\}$$

Subquantile Optimization Problem

We are now able to define the optimization problem we will solve:

$$\boldsymbol{\theta}_{SM} = \operatorname{argmin}_{\boldsymbol{\theta} \in \mathbb{R}^d} \max_{t \in \mathbb{R}} \left\{ t - \frac{1}{np} \sum_{i=1}^n \left(t - \left(\boldsymbol{\theta}^\top \mathbf{x}_i - y_i \right) \right)^+ \right\}$$

Algorithm:

$$t_{(k+1)} = \operatorname{argmax}_{t \in \mathbb{R}} g(t, \boldsymbol{\theta}_{(k)})$$

$$\boldsymbol{\theta}_{(k+1)} = \boldsymbol{\theta}_{(k)} - \alpha \nabla_{\boldsymbol{\theta}} g(t_{(k+1)}, \boldsymbol{\theta}_{(k)})$$

Lemma

$$\nabla_{\boldsymbol{\theta}} g(t, \boldsymbol{\theta}_{(k)}) = \frac{1}{np} \sum_{i=1}^{np} 2\mathbf{x}_i \left(\boldsymbol{\theta}_{(k)}^{\top} \mathbf{x}_i - y_i \right)$$

where $\{(\mathbf{x}_i, y_i)\}_{i=1}^{np}$ represent the np points in the dataset with the lowest loss.

Lemma

$$\operatorname{argmax}_{t_{k+1} \in \mathbb{R}} g(t, \boldsymbol{\theta}_{(k)}) = y_{np}$$

where y_{np} represents the np th highest loss in the dataset.

Here we are able to see the true nature of Subquantile Optimization. Each iteration we are optimizing over the points within the lowest np errors.

Definition

(t^*, θ^*) is a **Local Nash Equilibrium** of g if there exists $\delta > 0$ such that for any t, θ satisfying $\|t - t^*\| \leq \delta$ and $\|\theta - \theta^*\| \leq \delta$

Lemma

Any Local Nash Equilibrium satisfies $\nabla_{\theta} g(t_{(k)}, \theta_{(k)}) = \mathbf{0}$ and $\nabla_t g(t_{(k)}, \theta_{(k)}) = 0$

We first give intuition on what it means to be at a Local Nash Equilibrium. It means we have a θ that gives minimizes ERM over the points within the lowest np errors.

- [1] Sever computes the gradient of losses in each iteration. This incurs a $\mathcal{O}(dn^2)$ time complexity per iteration. Furthermore, points thrown out in early iterations are not resampled.
- Our work computes a np partition of the loss in each iteration. This incurs only a $\mathcal{O}(n)$ time complexity per iteration. Subquantile Minimization is novel in that it resamples points that may not have been in previous Subquantiles.

Objectives	Test RMSE (Drug Discovery)			
	$\epsilon = 0.1$	$\epsilon = 0.2$	$\epsilon = 0.3$	$\epsilon = 0.4$
ERM	1.303 _(0.0665)	1.790 _(0.0849)	2.198 _(0.0645)	2.623 _(0.1010)
CRR [2]	1.079 _(0.0899)	1.125 _(0.0832)	1.385 _(0.1372)	1.725 _(0.1136)
STIR [4]	1.087 _(0.1256)	1.167 _(0.0750)	1.403 _(0.0987)	1.668 _(0.1142)
Robust Risk [3]	1.176 _(0.1110)	1.336 _(0.1882)	1.437 _(0.1723)	1.800 _(0.0820)
SMART [5]	1.094 _(0.1065)	1.323 _(0.0758)	1.578 _(0.0799)	1.984 _(0.2020)
TERM [6]	1.326 _(0.0757)	1.357 _(0.0990)	1.310 _(0.0670)	1.302 _(0.0851)
SEVER [1]	1.111 _(0.0924)	1.067 _(0.0457)	1.071 _(0.0807)	1.138 _(0.1162)
Huber [7]	1.412 _(0.0474)	1.501 _(0.2918)	2.231 _(0.9054)	2.247 _(1.0399)
RANSAC [8]	1.238 _(0.0529)	1.643 _(0.1331)	2.092 _(0.1935)	2.679 _(0.1365)
SubQuantile($p = 1 - \epsilon$)	0.887 _(0.1046)	0.936 _(0.1051)	0.927 _(0.0729)	1.015 _(0.0978)
Genie ERM	0.959 _(0.0669)	0.955 _(0.0698)	1.038 _(0.0886)	1.004 _(0.0791)

Table: Drug Discovery Dataset. Empirical Risk over P with oblivious noise



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Definitions & Examples

Definition

A **prime number** is a number that has exactly two divisors.

Example

- 2 is prime (two divisors: 1 and 2).
- 3 is prime (two divisors: 1 and 3).
- 4 is not prime (**three** divisors: 1, 2, and 4).

You can also use the theorem, lemma, proof and corollary environments.

Example (Theorem Slide Code)

```
\begin{frame}  
\frametitle{Theorem}  
\begin{theorem}[Mass--energy equivalence]  
$E = mc^2$  
\end{theorem}  
\end{frame}
```

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The End

Questions? Comments?