

## Statistica Sinica Preprint No: SS-2019-0045

<b>Title</b>	A New Principle for Tuning-Free Huber Regression
<b>Manuscript ID</b>	SS-2019-0045
<b>URL</b>	<a href="http://www.stat.sinica.edu.tw/statistica/">http://www.stat.sinica.edu.tw/statistica/</a>
<b>DOI</b>	10.5705/ss.202019.0045
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## A New Principle for Tuning-Free Huber Regression

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*Abstract:* The robustification parameter, which balances bias and robustness, plays a critical role in the construction of subGaussian estimators for heavy-tailed and/or skewed data. Although the parameter can be tuned using cross-validation, in large-scale statistical problems such as high-dimensional covariance matrix estimation and large-scale multiple testing, the number of robustification parameters increases with the dimensionality causing cross-validation to become computationally prohibitive. We propose a new data-driven principle for choosing the robustification parameter for Huber-type subGaussian estimators in three fundamental problems: mean estimation, linear regression, and sparse regression in high dimensions. Our proposal is guided by a nonasymptotic deviation analysis, and is conceptually different from cross-validation, which relies on the mean squared error to assess the fit. Extensive numerical experiments and a real-data analysis further illustrate the efficacy of the proposed methods.

*Key words and phrases:* Data adaptive, heavy tails, Huber loss,  $M$ -estimator, tuning parameters

### 1. Introduction

Data subject to heavy-tailed and/or skewed distributions are frequently observed in various disciplines (Cont, 2001; Purdom and Holmes, 2005). A random variable  $X$  is heavy-tailed if its tail probability  $\mathbb{P}(|X| > t)$  decays to zero polynomially in  $1/t$

as  $t \rightarrow \infty$ , or equivalently, if  $X$  has finite polynomial-order moments. The connection between the moments and the tail probability is revealed by the property that  $\mathbb{E}(|X|^k) = k \int_0^\infty t^{k-1} \mathbb{P}(|X| > t) dt$ , for any  $k \geq 1$ . Here the sampling distribution has only a small number of finite moments, with a high chance that some observations deviate significantly from the population mean. Such observations are known as outliers, and are caused by heavy-tailed noise. In contrast, samples generated from a Gaussian or subGaussian distribution (Vershynin, 2012) are strongly concentrated around the expected value, making the chance of having extreme observations much smaller.

Heavy-tailed data bring new challenges to conventional statistical methods. For linear models, regression estimators based on the least squares loss are suboptimal, both theoretically and empirically, in the presence of heavy-tailed errors. See Catoni (2012) for a deviation analysis showing that the deviation of the empirical mean can be much worse for nonGaussian samples than it is for Gaussian ones. More broadly, this study exemplifies the pitfalls of asymptotic studies in statistics, and inspires new notions of optimality commonly used to assess the performance of estimators. In particular, the minimax optimality under the mean squared error does not quite capture the influence of estimators' extreme behaviors. However, these rare events may have severe negative effects in practice, leading to wrong conclusions or false discoveries. Since the work of Catoni (2012), nonasymptotic deviation analyses have drawn considerable attention, and are becoming increasingly important in the construction of subGaussian estimators (see Section S1.2 in the Supplementary Materials) for heavy-tailed data; see, for

example, Brownlees, Joly, and Lugosi (2015), Minsker (2015, 2018), Hsu and Sabato (2016), Devroye et al. (2016), Lugosi and Mendelson (2016), Fan, Li, and Wang (2017), Lugosi and Mendelson (2019), Lecué and Lerasle (2017), and Zhou et al. (2018), among others.

For linear models, Fan, Li, and Wang (2017) and Zhou et al. (2018) proposed Huber-type estimators in both low- and high-dimensional settings' and derived nonasymptotic deviation bounds for the estimation error. To implement either Catoni's or a Huber-type method, a tuning parameter  $\tau$  needs to be specified in advance to balance the robustness and bias of the estimation. A deviation analysis suggests that this tuning parameter, which we refer to as the robustification parameter, should adapt to the sample size, dimension, variance of the noise, and confidence level. Calibration schemes are typically based on cross-validation or Lepski's method, which can be computationally intensive, especially for large-scale inference and high-dimensional estimation problems, where the number of parameters may be exponential in the number of observations.

For example, Avella-Medina et al. (2018) proposed adaptive Huber estimators for estimating high-dimensional covariance and precision matrices. For a  $d \times d$  covariance matrix, although every entry can be robustly estimated using a Huber-type estimator with  $\tau$  chosen via cross-validation, the overall procedure involves as many as  $d^2$  tuning parameters. As a result, the cross-validation method soon becomes computationally intractable as  $d$  grows. Efficient tuning is important, not only for the problem, but also for applications in a broader context.

First, we develop data-driven Huber-type methods for mean estimation, linear regression, and sparse regression in high dimensions. For each problem, we first provide subGaussian concentration bounds for the Huber-type estimator under a minimal moment condition on the errors. These nonasymptotic results guide the choice of key tuning parameters. Some are of independent interest, and improve existing results by weakening the sample size scaling. Second, we propose a novel data-driven principle to calibrate the robustification parameter  $\tau > 0$  in the Huber loss

$$\ell_\tau(x) = \begin{cases} x^2/2 & \text{if } |x| \leq \tau, \\ \tau|x| - \tau^2/2 & \text{if } |x| > \tau. \end{cases} \quad (1.1)$$

Huber proposed using  $\tau = 1.345\sigma$  to retain 95% of the asymptotic efficiency of the estimator for normally distributed data, and to guarantee the estimator's performance for arbitrary contamination in a neighborhood of the true model (Huber, 1981; Huber and Ronchetti, 2009). This default setting is useful in high-dimensional statistics, even though the asymptotic efficiency is no longer well defined; see, for example, Lambert-Lacroix and Zwald (2011), Elsener and van de Geer (2018), and Loh (2017). Guided by the nonasymptotic deviation analysis, our proposed  $\tau$  grows with the sample size for the bias-robustness trade-off. For linear regressions under different regimes, the optimal  $\tau$  depends on the dimension  $d$ :  $\tau \sim \sigma\sqrt{(n/d)}$  in the low-dimensional setting with small  $d/n$ , and  $\tau \sim \sigma\sqrt{n/\log(d)}$  in high dimensions. Lastly, we design simple and fast algorithms to implement our method for calibrating  $\tau$ .

We focus on the notion of tail robustness (Catoni, 2012; Minsker, 2018; Zhou et

al., 2018; Fan, Li, and Wang, 2017; Avella-Medina et al., 2018), which is characterized by tight nonasymptotic deviation guarantees for the estimators under weak moment assumptions, and is evidenced by better finite-sample performance in the presence of heavy-tailed and/or highly skewed noise. This is inherently different from the traditional definition of robustness under Huber's  $\epsilon$ -contamination model (Huber and Ronchetti, 2009). Following the introduction of the finite-sample breakdown point by Donoho and Huber (1983), traditional robust statistics have focused, in part, on the development of high breakdown point estimators. Informally, the breakdown-point of an estimator is defined as the largest proportion of contaminated samples in the data that an estimator can tolerate before it produces arbitrarily large estimates (Hampel, 1971; Hampel et al., 1986; Maronna et al., 2018). A high breakdown point does not necessarily shed light on an estimator's convergence properties, efficiency, and stability. Refer to Portnoy and He (2000) for a review of classical robust statistics. In contrast, a tail robust estimator is resilient to outliers caused by heavy-tailed noise. Intuitively, the breakdown point describes the worst-case robustness, whereas our focus corresponds to the average-case robustness.

The remainder of this paper is organized as follows. In Section 2, we revisit Catoni's method on robust mean estimation. Motivated by a careful analysis of the truncated sample mean, we introduce a novel data-driven adaptive Huber estimator. We extend this data-driven tuning scheme to robust regression in Section 3 under both low- and high-dimensional settings. Extensive numerical experiments are reported in Section 4

to demonstrate the finite-sample performance of the proposed procedures. Section 5 concludes the paper. All proofs, together with technical details and real-data examples, are relegated to the Supplementary Material.

## 2. Robust data-adaptive mean estimation

### 2.1 Motivation

To motivate our proposed data-driven scheme for Huber-type estimators, we start by revisiting the mean estimation problem. Let  $X_1, \dots, X_n$  ( $n \geq 2$ ) be independent and identically distributed (*i.i.d.*) copies of  $X$  with mean  $\mu$  and finite variance  $\sigma^2 > 0$ . The sample mean, denoted as  $\bar{X}_n$ , is the most natural estimator for  $\mu$ . However, it suffers severely from not being robust to heavy-tailed sampling distributions (Catoni, 2012).

In order to cancel, or at least dampen, the erratic fluctuations in  $\bar{X}_n$ , which are more likely to occur if the distribution of  $X$  is heavy-tailed, we consider the truncated sample mean  $m_\tau = n^{-1} \sum_{i=1}^n \psi_\tau(X_i)$ , for some  $\tau > 0$ , where

$$\psi_\tau(x) = \text{sign}(x) \min(|x|, \tau) \quad (2.1)$$

is a truncation function on  $\mathbb{R}$ . Here, the tuning parameter  $\tau$  controls the bias and tail robustness of  $m_\tau$ . To see this, note that the bias term  $\text{Bias} := \mathbb{E}(m_\tau) - \mu$  satisfies  $|\text{Bias}| = |\mathbb{E}\{X - \text{sign}(X)\tau\}I(|X| > \tau)| \leq \tau^{-1}\mathbb{E}(X^2)$ . For tail robustness, the following result shows that  $m_\tau$  with a properly chosen  $\tau$  is a subGaussian estimator, as long as the second moment of  $X$  is finite.

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**Proposition 2.1.** Assume that  $v_2 := \sqrt{\mathbb{E}(X^2)}$  is finite. For any  $z > 0$ ,

- (i)  $m_\tau$  with  $\tau = v\sqrt{n/z}$ , for some  $v \geq v_2$ , satisfies  $\mathbb{P}\{|m_\tau - \mu| \geq 2v\sqrt{z/n}\} \leq 2e^{-z}$ ;
- (ii)  $m_\tau$  with  $\tau = cv_2\sqrt{n/z}$ , for some  $0 < c \leq 1$ , satisfies  $\mathbb{P}\{|m_\tau - \mu| \geq 2(v_2/c)\sqrt{z/n}\} \leq 2e^{-z/c^2}$ .

Proposition 2.1 shows how  $m_\tau$  performs under various idealized scenarios, thus providing guidance on the choice of  $\tau$ . Here,  $z > 0$  is a user-specified parameter that controls the confidence level; see the discussion before Remark 2.2. Given a properly tuned  $\tau$ , the subGaussian performance is achieved; conversely, if the resulting estimator performs well, the data are truncated at the right level and can be further exploited. An ideal  $\tau$  is such that the sample mean of the truncated data  $\psi_\tau(X_1), \dots, \psi_\tau(X_n)$  serves as a good estimator of  $\mu$ . The influence of outliers caused by heavy-tailed noise is weakened owing to the proper truncation. At the same time, we may expect that the empirical second moment for the same truncated data will provide a reasonable estimate of  $v_2^2$ .

Motivated by this, we propose choosing  $\tau > 0$  by solving  $\tau = \{\sum_{i=1}^n \psi_\tau^2(X_i)\}^{1/2} \sqrt{n/z}$ , which is equivalent to

$$\frac{1}{n} \sum_{i=1}^n \frac{\psi_\tau^2(X_i)}{\tau^2} = \frac{z}{n}, \quad \tau > 0. \quad (2.2)$$

We show that, under mild conditions, (2.2) has a unique solution  $\hat{\tau}_z$ , which gives rise to the following data-driven mean estimator:

$$m_{\hat{\tau}_z} = \frac{1}{n} \sum_{i=1}^n \min(|X_i|, \hat{\tau}_z) \text{sign}(X_i). \quad (2.3)$$

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To understand the property of  $\hat{\tau}_z$ , consider the population version of (2.2):

$$\frac{\mathbb{E}\{\psi_\tau^2(X)\}}{\tau^2} = \frac{\mathbb{E}\{\min(X^2, \tau^2)\}}{\tau^2} = \frac{z}{n}, \quad \tau > 0. \quad (2.4)$$

The following result establishes the existence and uniqueness of the solution to (2.4).

**Proposition 2.2.** Assume that  $v_2 = \sqrt{\mathbb{E}(X^2)}$  is finite.

(i) Provided  $0 < z < n\mathbb{P}(|X| > 0)$ , (2.4) has a unique solution  $\tau_z$  that satisfies

$$[\mathbb{E}\{\min(X^2, q_{z/n}^2)\}]^{1/2} \sqrt{n/z} \leq \tau_z \leq v_2 \sqrt{n/z}, \text{ where } q_\alpha := \inf\{t : \mathbb{P}(|X| > t) \leq \alpha\}$$

is the upper  $\alpha$ -quantile of  $|X|$ .

(ii) Let  $z = z_n > 0$  satisfy  $z_n \rightarrow \infty$  and  $z = o(n)$ . Then,  $\tau_z \rightarrow \infty$  and  $\tau_z \sim v_2 \sqrt{n/z}$ ,

as  $n \rightarrow \infty$ .

As a direct consequence of Proposition 2.2, the following result ensures the existence and uniqueness of the solution to (2.2), the empirical counterpart of (2.4).

**Proposition 2.3.** Provided  $0 < z < \sum_{i=1}^n I(|X_i| > 0)$ , (2.2) admits a unique solution.

Throughout, denote  $\hat{\tau}_z$  as the solution to (2.2), which is unique and positive whenever  $z < \sum_{i=1}^n I(|X_i| > 0)$ . For completeness, we set  $\hat{\tau}_z = 0$  on  $\{z \geq \sum_{i=1}^n I(|X_i| > 0)\}$ . If  $\mathbb{P}(X = 0) = 0$  and  $0 < z < n$ , then  $\hat{\tau}_z > 0$  with probability one. With both  $\tau_z$  and  $\hat{\tau}_z$  well defined, we investigate the property of  $\hat{\tau}_z$  below.

**Theorem 2.1.** Assume  $\mathbb{E}(X^2) < \infty$  and  $\mathbb{P}(X = 0) = 0$ . For any  $1 \leq z < n$  and  $0 < r < 1$ , we have

$$\mathbb{P}(|\hat{\tau}_z/\tau_z - 1| \geq r) \leq e^{-a_1^2 r^2 z^2 / (2z + 2a_1 r z / 3)} + e^{-a_2^2 r^2 z / 2} + 2e^{-(a_1 \wedge a_2)^2 z / 8}, \quad (2.5)$$

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where

$$a_1 = a_1(z, r) = \frac{P(\tau_z)}{2Q(\tau_z)} \frac{2+r}{(1+r)^2} \quad \text{and} \quad a_2 = a_2(z, r) = \frac{P(\tau_z - \tau_z r)}{2Q(\tau_z)} \frac{2-r}{1-r}, \quad (2.6)$$

with  $P(t) = \mathbb{E}\{X^2 I(|X| \leq t)\}$  and  $Q(t) = \mathbb{E}\{\psi_t^2(X)\}$ .

**Remark 2.1.** Here, we give some direct implications of Theorem 2.1.

(i) Let  $z = z_n \geq 1$  satisfy  $z = o(n)$  and  $z \rightarrow \infty$  as  $n \rightarrow \infty$ . By Proposition 2.2,  $\tau_z \rightarrow \infty$  and  $\tau_z \sim v_2 \sqrt{n/z}$ , which implies  $P(\tau_z) \rightarrow v_2^2$  and  $Q(\tau_z) \rightarrow v_2^2$  as  $n \rightarrow \infty$ .

(ii) With  $r = 1/2$  and  $z = \log^\kappa(n)$ , for some  $\kappa \geq 1$ , in (2.5), the constants  $a_1 = a_1(z, 1/2)$  and  $a_2 = a_2(z, 1/2)$  satisfy  $a_1 \rightarrow 5/9$  and  $a_2 \rightarrow 3/2$ , respectively, as  $n \rightarrow \infty$ . The resulting  $\hat{\tau}_z$  satisfies that, with probability approaching one,  $\tau_z/2 \leq \hat{\tau}_z \leq 3\tau_z/2$ .

We conclude this section with a uniform deviation bound for  $m_\tau$ . The uniformity of the rate over a neighborhood of the optimal tuning scale requires an additional  $\log(n)$ -factor. As a result, we show that the data-driven estimator  $m_{\hat{\tau}_z}$  is tightly concentrated around the mean with high probability.

**Theorem 2.2.** For  $z \geq 1$ , let  $\tau_z^* = v_2 \sqrt{n/z}$ . Then, with probability at least  $1 - 2ne^{-z}$ ,

$$\sup_{\tau_z^*/2 \leq \tau \leq 3\tau_z^*/2} |m_\tau - \mu| \leq 4v_2(z/n)^{1/2} + v_2 n^{-1/2}. \quad (2.7)$$

Letting  $z = 2 \log(n)$  and  $\hat{\tau}_z$  be the solution to (2.2), we obtain the following concentration inequality for the mean estimator  $m_{\hat{\tau}_z}$  given in (2.3).

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**Corollary 2.1.** With probability at least  $1 - c_1 n^{-c_2}$  for all sufficiently large  $n$ , we have

$$|m_{\hat{\tau}_z} - \mu| \leq 4v_2 \sqrt{2 \log(n)/n} + v_2 n^{-1/2}, \quad (2.8)$$

where  $c_1, c_2 > 0$  are absolute constants.

### 2.2 Adaptive Huber estimator

For the truncation method, even with the theoretically desirable tuning parameter  $\tau = v_2 \sqrt{n/z}$ , the deviation of the resulting estimator scales only with  $v_2$ , rather than with the standard deviation  $\sigma$ . The optimal deviation, which is enjoyed by the sample mean with subGaussian data, is of order  $\sigma \sqrt{z/n}$ . To achieve such an optimal order, [Fan, Li, and Wang \(2017\)](#) modified Huber's method to construct an estimator that exhibits fast (subGaussian type) concentration under a finite-variance condition.

The Huber loss in (1.1) is continuously differentiable with  $\ell'_\tau(x) = \psi_\tau(x)$ , where  $\psi_\tau(\cdot)$  is defined in (2.1). The Huber estimator is obtained as  $\hat{\mu}_\tau = \operatorname{argmin}_{\theta \in \mathbb{R}} \sum_{i=1}^n \ell_\tau(X_i - \theta)$ , or equivalently,  $\hat{\mu}_\tau$  is the unique solution to

$$0 = \sum_{i=1}^n \psi_\tau(X_i - \theta) = \sum_{i=1}^n \min(|X_i - \theta|, \tau) \operatorname{sign}(X_i - \theta). \quad (2.9)$$

Refer to [Catoni \(2012\)](#) for a general class of robust mean estimators. The following result from Theorem 5 in [Fan, Li, and Wang \(2017\)](#) shows the exponential-type concentration of  $\hat{\mu}_\tau$  when  $\tau$  is properly calibrated.

**Proposition 2.4.** Let  $z > 0$  and  $v \geq \sigma$ . Provided  $n \geq 8z$ ,  $\hat{\mu}_\tau$  with  $\tau = v \sqrt{n/z}$  satisfies the bound  $|\hat{\mu}_\tau - \mu| \leq 4v \sqrt{z/n}$  with probability at least  $1 - 2e^{-z}$ .

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Proposition 2.4 indicates that a theoretically desirable tuning parameter for the Huber estimator is  $\tau \sim \sigma\sqrt{n/z}$ . Motivated by the data-driven approach proposed in Section 2.1, we consider the following modification of (2.4):

$$\frac{\mathbb{E}\{\psi_\tau^2(X - \mu)\}}{\tau^2} = \frac{\mathbb{E}[\min\{(X - \mu)^2, \tau^2\}]}{\tau^2} = \frac{z}{n}, \quad \tau > 0. \quad (2.10)$$

According to Proposition 2.2, provided  $0 < z < n\mathbb{P}(X \neq \mu)$ , (2.10) admits a unique solution  $\tau_{z,\mu}$  that satisfies  $\sqrt{\mathbb{E}[\min\{(X - \mu)^2, \bar{q}_{z/n}\}]} \sqrt{n/z} \leq \tau_{z,\mu} \leq \sigma\sqrt{n/z}$ , where  $\bar{q}_\alpha = \inf\{t : \mathbb{P}(|X - \mu| > t) \leq \alpha\}$ . From a large-sample perspective, if  $z = z_n$  satisfies  $z \rightarrow \infty$  and  $z = o(n)$ , then  $\tau_{z,\mu} \rightarrow \infty$  and  $\tau_{z,\mu} \sim \sigma\sqrt{n/z}$  as  $n \rightarrow \infty$ .

Based on (2.9) and (2.10), a clearly motivated data-driven estimate of  $\mu$  can be obtained by solving the following system of equations:

$$\begin{cases} f_1(\theta, \tau) := \sum_{i=1}^n \psi_\tau(X_i - \theta) = 0, \\ f_2(\theta, \tau) := n^{-1} \sum_{i=1}^n \min\{(X_i - \theta)^2, \tau^2\}/\tau^2 - n^{-1}z = 0, \end{cases} \quad \theta \in \mathbb{R}, \tau > 0. \quad (2.11)$$

Observe that for any given  $\tau > 0$ ,  $f_1(\cdot, \tau) = 0$  always admits a unique solution, and for any given  $\theta$ ,  $f_2(\theta, \cdot) = 0$  has a unique solution, provided that  $z < \sum_{i=1}^n I(X_i \neq \theta)$ . With initial values  $\theta^{(0)} = \bar{X}_n$  and  $\tau^{(0)} = \hat{\sigma}_n\sqrt{n/z}$ , where  $\hat{\sigma}_n^2$  denotes the sample variance, we can solve (2.11) successively by computing a sequence of solutions  $\{(\theta^{(k)}, \tau^{(k)})\}_{k \geq 1}$  that satisfy  $f_2(\theta^{(k-1)}, \tau^{(k)}) = 0$  and  $f_1(\theta^{(k)}, \tau^{(k)}) = 0$ , for  $k \geq 1$ . For a predetermined tolerance  $\epsilon$ , the algorithm terminates within the  $\ell$ th iteration when  $\max\{|\theta^{(\ell)} - \theta^{(\ell-1)}|, |\tau^{(\ell)} - \tau^{(\ell-1)}|\} \leq \epsilon$ , and uses  $\theta^{(\ell)}$  as a robust estimator of  $\mu$ .

In the case of  $z = 1$ , the algorithm stops in the first iteration and delivers the

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solution  $\bar{X}_n$ . According to the results in Section 2.1, for fixed  $z \geq 1$ , there is no net improvement in terms of tail robustness; instead, we should let  $z = z_n$  grow slowly with the sample size to achieve tail robustness without introducing extra bias. Specifically, we choose  $z = \log(n)$  throughout our numerical experiments.

**Remark 2.2.** The proposed estimator is obtained by iteratively solving (2.11), which mimics (1.6) in Bickel (1975), and can be viewed as a variant of (6.28) and (6.29) in Huber and Ronchetti (2009) for joint location and scale estimation. The estimator in Bickel (1975) solves the equation  $\sum_{i=1}^n \psi_{\hat{\sigma}}(X_i - \theta) = 0$ , where  $\hat{\sigma}$  is chosen independently as the normalized interquartile range  $\hat{\sigma}^{(1)} = \{X_{(n-[n/4]+1)} - X_{([n/4])}\}/2\Phi^{-1}(3/4)$  or the symmetrized interquartile range  $\hat{\sigma}^{(2)} = \text{median}\{|X_i - m|\}/\Phi^{-1}(3/4)$ , where  $X_{(1)} < \dots < X_{(n)}$  are the order statistics and  $m$  is the sample median. The consistency of  $\hat{\sigma}^{(1)}$  or  $\hat{\sigma}^{(2)}$  is established under the symmetry assumption of  $X$ , but remains unclear for general distributions. On the other hand, similarly to Bickel (1975), our proposed estimators of  $\theta$  and  $\tau$  are also location and scale equivariant (see Sections S1.7 and S1.8 in the Supplementary Materials).

Unlike this classical approach, we waive the symmetry requirement by allowing the robustification parameter to diverge in order to reduce the bias induced by the Huber loss when the distribution is asymmetric. Another difference is that Bickel's proposal is a two-step method that estimates the scale and location separately, whereas our procedure estimates  $\mu$  and calibrates  $\tau$  simultaneously by solving a system of equations. In fact, as a direct extension of the idea in Section 2.1, we can also tune  $\tau$  independently

from the estimation by solving  $\binom{n}{2}^{-1} \sum_{1 \leq i < j \leq n} \tau^{-2} \min\{(X_i - X_j)^2/2, \tau^2\} = zn^{-1}$ , for  $z > 0$ . Let  $X'$  be an independent copy of  $X$ . Then, the population version of this equation is  $\mathbb{E}[\min\{(X - X')^2/2, \tau^2\}] \tau^{-2} = z/n$ , the solution of which is unique under mild conditions and scales as  $\sigma\sqrt{n/z}$ .

**Remark 2.3.** We assume a finite variance of errors. For more subtle scenarios with a finite  $(1 + \delta)$ th moment and  $0 < \delta < 1$ , the phase transition phenomenon discovered by [Devroye et al. \(2016\)](#) and [Sun, Zhou, and Fan \(2020\)](#) suggests that Huber's  $M$ -estimator no longer admits subGaussian-type deviation bounds. Developing the corresponding data-driven principle to tune Huber's method when  $\delta < 1$  is nontrivial, and thus is left as a topic for future investigation.

### 3. Robust data-adaptive linear regression

In this section, we extend the proposed data-driven method for robust mean estimation to regression problems. Consider the linear regression model

$$Y_i = \beta_0^* + \mathbf{X}_i^\top \boldsymbol{\beta}^* + \varepsilon_i, \quad i = 1, \dots, n, \quad (3.1)$$

where  $Y_i$  is a response variable,  $\mathbf{X}_i$  is a  $d$ -dimensional vector of covariates,  $\beta_0^*$  and  $\boldsymbol{\beta}^* \in \mathbb{R}^d$  denote the intercept and vector, respectively, of the regression coefficients, and  $\varepsilon_1, \dots, \varepsilon_n$  are independent regression errors with zero mean and finite variance. For ease of presentation, we write  $\mathbf{Z}_i = (1, \mathbf{X}_i^\top)^\top$  and  $\boldsymbol{\theta}^* = (\beta_0^*, \boldsymbol{\beta}^*)^\top$ . The goal is to estimate  $\boldsymbol{\theta}^*$  from the observed data  $\{(Y_i, \mathbf{X}_i)\}_{i=1}^n$ .

### 3.1 Adaptive Huber regression in low dimensions

We start with the low-dimensional regime, where  $d \ll n$ . In the presence of heavy-tailed errors, the finite-sample properties of the least squares method are suboptimal, both theoretically and empirically. Under such heavy-tailed models, refer to [Audibert and Catoni \(2011\)](#) and [Sun, Zhou, and Fan \(2020\)](#) for a nonasymptotic analysis of Huber-type robust regressions; the former focuses on the excess risk bounds, and the latter provides deviation bounds for the estimator, along with nonasymptotic Bahadur representations.

Given  $\tau > 0$ , Huber's  $M$ -estimator is defined as

$$\widehat{\boldsymbol{\theta}}_\tau = (\widehat{\beta}_{0,\tau}, \widehat{\boldsymbol{\beta}}_\tau^\top)^\top \in \operatorname{argmin}_{\boldsymbol{\theta} \in \mathbb{R}^{d+1}} \sum_{i=1}^n \ell_\tau(Y_i - \mathbf{Z}_i^\top \boldsymbol{\theta}), \quad (3.2)$$

where  $\ell_\tau(\cdot)$  is given in (1.1). By the convexity of the Huber loss, the solution to (3.2) is determined uniquely using the first-order condition  $\sum_{i=1}^n \psi_\tau(Y_i - \mathbf{Z}_i^\top \widehat{\boldsymbol{\theta}}_\tau) \mathbf{Z}_i = \mathbf{0}$ . Most desirable features of Huber's method are established under the assumption that the error distribution is symmetric around zero. In the absence of symmetry, the bias induced by the Huber loss becomes non-negligible. To make this statement precise, note that  $\widehat{\boldsymbol{\theta}}_\tau = (\widehat{\beta}_{0,\tau}, \widehat{\boldsymbol{\beta}}_\tau^\top)^\top$  is a natural  $M$ -estimator of

$$\boldsymbol{\theta}_\tau^* = (\beta_{0,\tau}^*, \boldsymbol{\beta}_\tau^{*\top})^\top = \operatorname{argmin}_{(\beta_0, \boldsymbol{\beta}^\top) \in \mathbb{R}^{d+1}} \sum_{i=1}^n \mathbb{E}\{\ell_\tau(Y_i - \beta_0 - \mathbf{X}_i^\top \boldsymbol{\beta})\}, \quad (3.3)$$

whereas the true parameters  $\beta_0^*$  and  $\boldsymbol{\beta}^*$  are identified as  $\operatorname{argmin}_{\beta_0, \boldsymbol{\beta}} \sum_{i=1}^n \mathbb{E}\{(Y_i - \beta_0 - \mathbf{X}_i^\top \boldsymbol{\beta})^2\}$ . For any fixed  $\tau > 0$ , although  $\widehat{\beta}_{0,\tau}$  and  $\widehat{\boldsymbol{\beta}}_\tau$  are robust estimates of  $\beta_{0,\tau}^*$  and  $\boldsymbol{\beta}_\tau^*$ , respectively,  $(\beta_{0,\tau}^*, \boldsymbol{\beta}_\tau^*)$  differs from  $(\beta_0^*, \boldsymbol{\beta}^*)$ , in general. The following proposition

### 3.1 Adaptive Huber regression in low dimensions 15

provides an explicit bound on the bias, complementing the results in Section 4.9.2 of [Maronna et al. \(2018\)](#).

**Proposition 3.1.** Assume that  $\varepsilon$  and  $\mathbf{X}$  are independent, and that the function  $\alpha \mapsto \mathbb{E}\{\ell_\tau(\varepsilon - \alpha)\}$  has a unique minimizer  $\alpha_\tau = \operatorname{argmin}_{\alpha \in \mathbb{R}} \mathbb{E}\{\ell_\tau(\varepsilon - \alpha)\}$ , which satisfies

$$\mathbb{P}(|\varepsilon - \alpha_\tau| \leq \tau) > 0. \quad (3.4)$$

Assume further that  $\mathbb{E}(\mathbf{Z}\mathbf{Z}^\top)$  is positive definite. Then, we have  $\beta_{0,\tau}^* = \beta_0^* + \alpha_\tau$  and  $\boldsymbol{\beta}_\tau^* = \boldsymbol{\beta}^*$ . Moreover,  $\alpha_\tau$  with  $\tau > \sigma$  satisfies the bound

$$|\alpha_\tau| \leq \frac{\sigma^2 - \mathbb{E}\{\psi_\tau^2(\varepsilon)\}}{1 - \tau^{-2}\sigma^2} \frac{1}{\tau}. \quad (3.5)$$

Note too that the Huber loss minimization is equivalent to the penalized least squares problem ([She and Owen, 2011](#)),  $(\hat{\boldsymbol{\mu}}_\tau, \hat{\boldsymbol{\theta}}_\tau) = \operatorname{argmin}_{\boldsymbol{\mu} \in \mathbb{R}^n, \boldsymbol{\theta} \in \mathbb{R}^{d+1}} \{ \frac{1}{2} \sum_{i=1}^n (Y_i - \mu_i - \mathbf{Z}_i^\top \boldsymbol{\theta})^2 + \tau \sum_{i=1}^n |\mu_i| \}$ , where  $\boldsymbol{\mu} = (\mu_1, \dots, \mu_n)^\top$  and  $\hat{\boldsymbol{\theta}}_\tau$  coincide with those in (3.2). This loss function can be written as  $\sum_{i=1}^n (Y_i - \mu_i - \beta_0 - \mathbf{X}_i^\top \boldsymbol{\beta})^2 / 2 + \tau \sum_{i=1}^n |\mu_i|$ , which explains from a different perspective that the bias arises only at the intercept. The larger  $\tau$  is, the sparser  $\hat{\boldsymbol{\mu}}_\tau$  is and, therefore, the smaller the estimation bias is.

Proposition 3.1 draws attention to intercept estimation, a problem of independent interest that needs to be treated with greater caution. If the distribution of  $\varepsilon$  is asymmetric,  $\alpha_\tau$  is typically nonzero, for any  $\tau > 0$ ; here, a smaller  $\tau$  results in a larger bias and, thus, a larger prediction error. To balance the bias and the tail robustness, we propose two modifications to Huber's method (a one-step method, and a two-step

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method) that are robust against heavy-tailed and asymmetric errors, while maintaining high efficiency for normal data.

#### 3.1.1 One-step method

As noted in Zhou et al. (2018), there is an inherent bias–robustness trade-off in the choice of  $\tau$ , which should adapt to the sample size, dimension, and the variance of the noise; see Theorem 3.1. To begin with, we impose the following moment conditions.

**Condition 3.1.** The covariates  $\mathbf{X}_1, \dots, \mathbf{X}_n$  are *i.i.d.* random vectors from  $\mathbf{X}$ . There exists  $A_0 > 0$ , such that for any  $\mathbf{u} \in \mathbb{R}^{d+1}$  and  $t \in \mathbb{R}$ ,  $\mathbb{P}(|\langle \mathbf{u}, \mathbf{z} \rangle| \geq A_0 \|\mathbf{u}\|_2 \cdot t) \leq e^{-t}$ , where  $\mathbf{z} = \mathbf{S}^{-1/2} \mathbf{Z}$  and  $\mathbf{S} = \mathbb{E}(\mathbf{Z} \mathbf{Z}^\top)$  is positive definite. The regression errors  $\varepsilon_i$  are independent and satisfy  $\mathbb{E}(\varepsilon_i | \mathbf{X}_i) = 0$  and  $\mathbb{E}(\varepsilon_i^2 | \mathbf{X}_i) \leq \sigma^2$  almost surely.

**Theorem 3.1.** Assume Condition 3.1 holds. For any  $z > 0$  and  $v \geq \sigma$ , the estimator  $\widehat{\boldsymbol{\theta}}_\tau$  in (3.2), with  $\tau = v \sqrt{n/(d+z)}$ , satisfies the bound  $\|\mathbf{S}^{1/2}(\widehat{\boldsymbol{\theta}}_\tau - \boldsymbol{\theta}^*)\|_2 \leq c_1 v \sqrt{(d+z)/n}$  with probability at least  $1 - 2e^{-z}$ , provided  $n \geq c_2(d+z)$ , where  $c_1, c_2 > 0$  are constants depending only on  $A_0$ .

This theorem establishes a subGaussian concentration bound for  $\widehat{\boldsymbol{\theta}}_\tau$  under the optimal sampling size scaling. Compared with Theorem 2.1 in Zhou et al. (2018), there are two technical improvements: first, the moment condition on the random predictor is relaxed from subGaussian to sub-exponential; and second, the sample size requirement is improved to  $n \gtrsim d$ , which is in line with the classical asymptotic consistency result

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that requires  $d = o(n)$ . To achieve subGaussian performance under the finite-variance condition, the key observation is that the robustification parameter  $\tau$  should adapt to the sample size, dimension, variance of the noise, and confidence level for an optimal trade-off between bias and robustness. Extending our proposal for mean estimation, for  $\boldsymbol{\theta} \in \mathbb{R}^{d+1}$  and  $\tau > 0$ , we estimate  $\boldsymbol{\theta}^*$  and calibrate  $\tau$  simultaneously by solving the system of equations

$$\begin{cases} g_1(\boldsymbol{\theta}, \tau) := \sum_{i=1}^n \psi_\tau(Y_i - \mathbf{Z}_i^\top \boldsymbol{\theta}) \mathbf{Z}_i = \mathbf{0}, \\ g_2(\boldsymbol{\theta}, \tau) := (\tau^2 n)^{-1} \sum_{i=1}^n \min\{(Y_i - \mathbf{Z}_i^\top \boldsymbol{\theta})^2, \tau^2\} - n^{-1}(d + z) = 0. \end{cases} \quad (3.6)$$

With initial values  $\boldsymbol{\theta}^{(0)} := \widehat{\boldsymbol{\theta}}_{\text{ols}} = (\sum_{i=1}^n \mathbf{Z}_i \mathbf{Z}_i^\top)^{-1} \sum_{i=1}^n Y_i \mathbf{Z}_i$  and  $\tau^{(0)} = \widehat{\sigma}_n \sqrt{n/(d+z)}$ , where  $\widehat{\sigma}_n^2 = (1/n) \sum_{i=1}^n (Y_i - \mathbf{Z}_i^\top \widehat{\boldsymbol{\theta}}_{\text{ols}})^2$ , for  $k \geq 1$ , solve  $g_2(\boldsymbol{\theta}^{(k-1)}, \tau^{(k)}) = 0$  to obtain  $\tau^{(k)}$ , and then compute  $\boldsymbol{\theta}^{(k)}$  as the solution to  $g_1(\boldsymbol{\theta}^{(k)}, \tau^{(k)}) = 0$ . Iterate until convergence, and set  $\widehat{\boldsymbol{\theta}}^I := \widehat{\boldsymbol{\theta}}_{\widehat{\tau}}$  as the one-step estimator, where  $(\widehat{\boldsymbol{\theta}}, \widehat{\tau})$  is the final output.

The main advantage of the proposed adaptive Huber regression over the traditional one with  $\tau = 1.345\sigma$  is that the estimation bias with respect to the intercept is alleviated. Examining the proof of Proposition 3.1, we find that the bias is of order  $1/\tau$  when the second moment is finite, and is quadratic in  $1/\tau$  if the third moment is finite. The statistical error, on the other hand, is determined by the  $\ell_2$ -norm of the score function evaluated at  $\boldsymbol{\theta}^*$ , which is of order  $\sigma \sqrt{d/n} + \tau d/n$ ; see Theorem 3.2. The overall error is then optimized at  $\tau \asymp \sigma \sqrt{n/d}$ . For the normal model, because  $\max_{1 \leq i \leq n} |\varepsilon_i| \sim \sigma \sqrt{2 \log(2n)} \lesssim \sigma \sqrt{n/d}$ , the adaptive Huber estimator is almost identical to the least squares estimator. The numerical results in Section 4 provide strong

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support for the tail-adaptivity of our proposed data-driven Huber regression.

When  $\tau$  scales as a constant, such as  $c\sigma$ , the corresponding Huber loss is Lipschitz with a bounded score function, and because  $\beta_\tau^* = \beta^*$  for any  $\tau > 0$ , there is no sacrifice in bias when estimating the slope  $\beta^*$ . Again, a constant  $c$  is typically tuned to ensure a given level of asymptotic efficiency. The asymptotic properties of general robust  $M$ -estimators have been well studied in the literature; see [Avella-Medina and Ronchetti \(2015\)](#) for a selective overview. The next result further complements Theorem 3.1 by establishing the deviations of the Huber estimator with fixed  $\tau$  from a nonasymptotic viewpoint.

**Theorem 3.2.** Suppose Condition 3.1 and the assumptions in Proposition 3.1 hold. Assume further that  $\rho_\tau := \mathbb{P}(|\varepsilon - \alpha_\tau| \leq \tau/2) > 0$ . Then, the estimator  $\hat{\theta}_\tau$  in (3.2) satisfies  $\|\mathbf{S}^{1/2}(\hat{\theta}_\tau - \theta_\tau^*)\|_2 \lesssim \rho_\tau^{-1} A_0 \{\sigma \sqrt{(d+z)/n} + \tau(d+z)/n\}$ , for any  $z > 0$ , with probability at least  $1 - 2e^{-z}$ , provided  $n \geq c_3(d+z)$ , where  $c_3 > 0$  is a constant depending only on  $(A_0, \rho_\tau)$ .

#### 3.1.2 Two-step method

Motivated by our bias-robustness analysis and the results of the finite-sample investigation, we introduce a two-step procedure that estimates the regression coefficients and the intercept successively.

In the first step, we compute the Huber estimator  $\hat{\theta}_\tau = (\hat{\beta}_{0,\tau}, \hat{\beta}_\tau^{*\top})^\top$  by solving (3.2) with  $\tau = c\sigma$ . We take  $c = 1.345$  to retain the 95% efficiency for the normal model.

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Here,  $\sigma$  can be estimated simultaneously with  $\boldsymbol{\theta}^*$  by solving a system of equations, as in Huber's "Proposal 2" (Huber, 1964; Huber and Ronchetti, 2009), or we can fix  $\sigma$  at an initial robust estimate, and then optimize over  $\boldsymbol{\theta}$  (Hampel et al., 1986). We follow the former route and consider an iterative procedure. Start with an initial estimate  $\boldsymbol{\theta}^{(0)}$ . At iteration  $k = 0, 1, 2, \dots$ , we employ a simple procedure to obtain  $\widehat{\sigma}^{(k)}$ , based on which we update  $\boldsymbol{\theta}^{(k+1)}$ . This step involves two procedures.

*Procedure 1: Scale estimation.* Using the current estimate  $\boldsymbol{\theta}^{(k)}$ , we compute the vector of residuals  $\mathbf{r}^{(k)} = (r_1^{(k)}, \dots, r_n^{(k)})^\top$  and the robustification parameter  $\tau^{(k)} = 1.345\widehat{\sigma}^{(k)}$ , where  $\widehat{\sigma}^{(k)}$  denotes the median absolute deviation (MAD) estimator  $\text{median}\{|r_i^{(k)} - \text{median}(r_i^{(k)})|\}/\Phi^{-1}(3/4)$ .

*Procedure 2: Weighted least squares.* Compute the  $n \times n$  diagonal matrix  $\mathbf{W}^{(k)} = \text{diag}((1 + w_1^{(k)})^{-1}, \dots, (1 + w_n^{(k)})^{-1})$ , where  $w_i^{(k)} = |r_i^{(k)}|/\tau^{(k)} - 1$  if  $|r_i^{(k)}| > \tau^{(k)}$ , and  $w_i^{(k)} = 0$  if  $|r_i^{(k)}| \leq \tau^{(k)}$ . Then, we update  $\boldsymbol{\theta}^{(k)}$  to produce  $\boldsymbol{\theta}^{(k+1)}$  using the weighted least squares; that is,

$$\boldsymbol{\theta}^{(k+1)} = \underset{\boldsymbol{\theta} \in \mathbb{R}^{d+1}}{\text{argmin}} \sum_{i=1}^n \frac{(Y_i - \mathbf{Z}_i^\top \boldsymbol{\theta})^2}{1 + w_i^{(k)}} = (\mathbf{Z}^\top \mathbf{W}^{(k)} \mathbf{Z})^{-1} \mathbf{Z}^\top \mathbf{W}^{(k)} \mathbf{Y},$$

where  $\mathbf{Z} = (\mathbf{Z}_1, \dots, \mathbf{Z}_n)^\top \in \mathbb{R}^{n \times (d+1)}$  and  $\mathbf{Y} = (Y_1, \dots, Y_n)^\top$ .

Starting with  $\boldsymbol{\theta}^{(0)} = \widehat{\boldsymbol{\theta}}_{\text{ols}}$ , we repeat the above two procedures until convergence. Denote  $\widehat{\boldsymbol{\beta}}^{\text{II}} \in \mathbb{R}^d$  as the vector of coefficient estimates extracted from the final solution.

In the second step, observe that  $\beta_0^* = \mathbb{E}(\delta_i)$ , where  $\delta_i = Y_i - \mathbf{X}_i^\top \boldsymbol{\beta}^* = \beta_0^* + \varepsilon_i$  are the residuals. To estimate  $\beta_0^*$ , define the fitted residuals  $\widehat{\delta}_i = Y_i - \mathbf{X}_i^\top \widehat{\boldsymbol{\beta}}^{\text{II}}$ , and solve the

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system of equations

$$\begin{cases} f_1(\beta_0, \tau) := (\tau^2 n)^{-1} \sum_{i=1}^n \min\{(\hat{\delta}_i - \beta_0)^2, \tau^2\} - n^{-1} \log(n) = 0, \\ f_2(\beta_0, \tau) := \sum_{i=1}^n \psi_\tau(\hat{\delta}_i - \beta_0) = 0 \end{cases} \quad (3.7)$$

in the same way as (2.11) to obtain  $\hat{\beta}_0^{\text{II}}$ . Then,  $\hat{\boldsymbol{\theta}}^{\text{II}} = (\hat{\beta}_0^{\text{II}}, \hat{\boldsymbol{\beta}}^{\text{II}})$  is the two-step estimator of  $\boldsymbol{\theta}^*$ .

The two-step procedure leverages the fact that for the asymmetric regression errors with potentially heavy tails, the Huber loss with a fixed  $\tau$  introduces bias to the intercept estimation, but not to the estimation of the slope coefficients. To alleviate the influence of skewness in the error, in the second step, we use the adaptive Huber method with a divergent  $\tau$  to re-estimate the intercept. The two-step estimator therefore achieves both a high degree of tail robustness and unbiasedness.

### 3.2 Adaptive Huber regression in high dimensions

We now move to the high-dimensional setting, where  $d \gg n$  and  $\boldsymbol{\beta}^* = (\beta_1^*, \dots, \beta_d^*)^\top \in \mathbb{R}^d$  is sparse, with  $\|\boldsymbol{\beta}^*\|_0 := \sum_{j=1}^d I(\beta_j^* \neq 0) = s \ll n$ . Since the invention of the Lasso (Tibshirani, 1996), a variety of variable selection methods have been developed for finding a small group of response-associated covariates from a large pool; refer to Bühlmann and van de Geer (2011) and Hastie, Tibshirani, and Wainwright (2015) for a comprehensive review along this line.

The Lasso estimator is  $\hat{\boldsymbol{\beta}}_{\text{lasso}}(\lambda) \in \operatorname{argmin}_{\beta_0 \in \mathbb{R}, \boldsymbol{\beta} \in \mathbb{R}^d} \{(2n)^{-1} \sum_{i=1}^n (Y_i - \beta_0 - \mathbf{X}_i^\top \boldsymbol{\beta})^2 + \lambda \|\boldsymbol{\beta}\|_1\}$ , where  $\lambda > 0$  is a regularization parameter. Viewing the noise variable as

### 3.2 Adaptive Huber regression in high dimensions 21

Gaussian, this can be interpreted as a penalized maximum likelihood estimate, where the  $\ell_1$ -penalty encourages sparsity in the estimation. However, least squares fitting is sensitive to the tails of the error distributions, particularly for ultrahigh-dimensional covariates, because their spurious correlations with the noise can be large. Therefore, this method is not ideal in the presence of heavy-tailed noise.

Recently, [Fan, Li, and Wang \(2017\)](#) modified Huber's procedure to obtain an  $\ell_1$ -regularized robust estimator that admits the desirable concentration bound under a finite-variance condition on the regression errors. According to Section 3.1, the intercept, albeit often ignored in the literature, plays an important role in studies of robust methods. To take into account the effect of the intercept, we consider the regularized Huber estimator of the form

$$\widehat{\boldsymbol{\theta}}_H(\tau, \lambda) \in \underset{\boldsymbol{\theta} = (\beta_0, \boldsymbol{\beta}^\top)^\top \in \mathbb{R}^{d+1}}{\operatorname{argmin}} \{ \mathcal{L}_\tau(\boldsymbol{\theta}) + \lambda \|\boldsymbol{\beta}\|_1 \}, \quad (3.8)$$

where  $\mathcal{L}_\tau(\boldsymbol{\theta}) := (1/n) \sum_{i=1}^n \ell_\tau(Y_i - \mathbf{Z}_i^\top \boldsymbol{\theta}) = (1/n) \sum_{i=1}^n \ell_\tau(Y_i - \beta_0 - \mathbf{X}_i^\top \boldsymbol{\beta})$ , and  $\tau$  and  $\lambda$  are the robustification and regularization parameters, respectively.

Given  $\varepsilon_i$  with finite variance, Theorem 3.3 reveals that the  $\ell_1$ -regularized Huber regression with properly tuned  $(\tau, \lambda)$  gives rise to consistent estimators, with  $\ell_1$ - and  $\ell_2$ -errors scaling as  $s\sqrt{\log(d)/n}$  and  $\sqrt{s\log(d)/n}$ , respectively, under the sample size scaling  $n \gtrsim s\log(d)$ . These rates are exactly the minimax rates enjoyed by the Lasso with subGaussian errors.

**Theorem 3.3.** Assume Condition 3.1 holds, and denote by  $\underline{\lambda}_S$  the minimal eigenvalue

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of  $\mathbf{S}$ . Assume further that the unknown  $\boldsymbol{\beta}^*$  is sparse with  $s = \|\boldsymbol{\beta}^*\|_0$ . Let  $\sigma_{jj} = \mathbb{E}(X_j^2)$ , for  $j = 1, \dots, d$ . Then, the estimator  $\widehat{\boldsymbol{\theta}}_H(\tau, \lambda)$  given in (3.8), with  $\tau = \sigma\sqrt{n/\log(d)}$  and  $\lambda$  scaling with  $A_0 \max_{1 \leq j \leq d} \sigma_{jj}^{1/2} \sigma \sqrt{\log(d)/n}$ , satisfies

$$\|\widehat{\boldsymbol{\theta}}_H(\tau, \lambda) - \boldsymbol{\beta}^*\|_2 \lesssim \frac{s^{1/2}\lambda}{\underline{\lambda}_S} \quad \text{and} \quad \|\widehat{\boldsymbol{\theta}}_H(\tau, \lambda) - \boldsymbol{\beta}^*\|_1 \lesssim \frac{s\lambda}{\underline{\lambda}_S} \quad (3.9)$$

with probability at least  $1 - 3d^{-1}$ , as long as  $n \geq c_1 s \log(d)$ , where  $c_1 > 0$  is a constant depending only on  $(A_0, \max_{1 \leq j \leq d} \sigma_{jj}, \underline{\lambda}_S)$ .

Theorem 3.3 complements Theorem 3 in [Fan, Li, and Wang \(2017\)](#). The latter provides convergence rates for an  $\ell_1$ -penalized Huber  $M$ -estimator under the weakly sparse setting that  $\|\boldsymbol{\beta}^*\|_q \leq R_q$ , for some  $0 < q \leq 1$ . Their results, however, do not directly apply to the sparse regime where  $q = 0$ . Moreover, the subGaussian condition imposed in [Fan, Li, and Wang \(2017\)](#) is now relaxed to the sub-exponential condition.

**Remark 3.1.** The main purpose of using the Huber loss for data fitting is to gain robustness against outliers from the contamination models ([Huber, 1973](#)) or the heavy-tailed models considered here. For other purposes, different loss functions have been proposed to replace the squared loss, such as the nonconvex Tukey and Cauchy losses, quantile loss, and asymmetric quadratic loss, among others. Refer to [Owen \(2007\)](#), [Loh and Wainwright \(2015\)](#), [Loh \(2017\)](#), [Zhou et al. \(2018\)](#), [Mei, Bai, and Montanari \(2018\)](#), [Alquier, Cottet, and Lecué \(2019\)](#), and [Pan, Sun, and Zhou \(2019\)](#) for discussions on the regularized  $M$ -estimator with different loss functions.

In practice, it is computationally demanding to choose the optimal values of  $\tau$  and

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$\lambda$  using a two-dimensional grid search and cross-validation. We consider the following procedure that estimates  $\boldsymbol{\theta}^*$  and tunes  $\tau$  simultaneously. Given a random sample of size  $n$ , we use a cross-validated Lasso as an initialization  $\widehat{\boldsymbol{\theta}}^{(0)}$ . At iteration  $k = 1, 2, \dots$ , using the previous estimate  $\widehat{\boldsymbol{\theta}}^{(k-1)}$ , we compute  $\tau^{(k)}$  as the solution to

$$\frac{1}{\{n - \widehat{s}^{(k-1)}\}} \sum_{i=1}^n \frac{\min\{(Y_i - \mathbf{Z}_i^\top \widehat{\boldsymbol{\theta}}^{(k-1)})^2, \tau^2\}}{\tau^2} = \frac{\log(nd)}{n}, \quad (3.10)$$

where  $\widehat{s}^{(k-1)} = \|\widehat{\boldsymbol{\beta}}^{(k-1)}\|_0$ . Next, take  $\tau = \tau^{(k)}$ , and compute  $\widehat{\boldsymbol{\theta}}^{(k)}$  by solving

$$\min_{\boldsymbol{\theta}} \left\{ \frac{1}{n} \sum_{i=1}^n \ell_\tau(Y_i - \mathbf{Z}_i^\top \boldsymbol{\theta}) + \lambda \|\boldsymbol{\beta}\|_1 \right\}, \quad (3.11)$$

where  $\lambda$  is chosen using cross-validation. Repeat the above two steps until convergence, or until the maximum number of iterations is reached.

To implement the data-driven Huber regression in high dimensions, starting with some initial guess, we iteratively solve (3.10) and (3.11). For the convex optimization problems in (3.11), the minimizer satisfies the Karush–Kuhn–Tucker conditions, and therefore can be found by solving the following system of nonsmooth equations:

$$\begin{cases} -n^{-1} \sum_i \psi_\tau(Y_i - \mathbf{Z}_i^\top \widehat{\boldsymbol{\theta}}) = 0, \\ -n^{-1} \sum_i \psi_\tau(Y_i - \mathbf{Z}_i^\top \widehat{\boldsymbol{\theta}}) X_{ij} + \lambda \widehat{\eta}_j = 0, \quad j = 1, \dots, d \\ \widehat{\beta}_j - S(\widehat{\beta}_j + \widehat{\eta}_j) = 0, \quad j = 1, \dots, d, \end{cases} \quad (3.12)$$

where  $\widehat{\boldsymbol{\theta}} = (\widehat{\beta}_0, \widehat{\boldsymbol{\beta}}^\top)^\top \in \mathbb{R}^{d+1}$  with  $\widehat{\boldsymbol{\beta}} = (\widehat{\beta}_1, \dots, \widehat{\beta}_d)^\top$ ,  $\widehat{\eta}_j \in \partial|\widehat{\beta}_j|$ , and  $S(z) = \text{sign}(z)(|z| - 1)_+$  is the soft-thresholding operator. Instead of directly applying the semismooth Newton algorithm (SNA) to the entire system of equations, we adapt the semismooth

### 3.2 Adaptive Huber regression in high dimensions 24

Newton coordinate descent (SNCD) algorithm proposed by [Yi and Huang \(2017\)](#), which combines the SNA with a cyclic coordinate descent to solve (3.12). More specifically, in the SNCD, we divide (3.12) into two parts in order to avoid the cumbersome matrix operations involved in solving the entire system. In a cyclic fashion, we update the intercept using only the first equation, and update the coefficients with its subgradients using the last two equations. Therefore, we reduce the computational cost from  $O(nd^2)$  to  $O(nd)$  at each iteration. The gain in computational scalability and efficiency is substantial for large  $d$ . After obtaining a solution path of (3.11), we employ the cross-validation method to select  $\lambda$  and then the associated  $\widehat{\boldsymbol{\theta}}^{(k)}$ .

**Remark 3.2.** The above regularized data-adaptive Huber (DA-Huber) regression method is a direct extension of the one-step method proposed in Section 3.1 to high dimensions. Furthermore, note that Proposition 3.1 holds in high dimensions, as long as the population Gram matrix  $\mathbf{S}$  is positive definite. Therefore, to further reduce the estimation bias of the intercept, we suggest using the two-step procedure that estimates the regression coefficients using the standard regularized Huber regression, and then estimates the intercept by applying the adaptive-Huber method to the fitted residuals, as in (3.7). Section 4.3 provides numerical studies of both the one- and the two-step regularized adaptive Huber estimators.

#### 4. Empirical analysis

In this section, we examine numerically the finite-sample performance of the proposed DA-Huber methods for mean estimation and linear regressions. In the Supplementary Material, using three real data sets, we also demonstrate the desirable performance of the proposed DA-Huber methods in terms of their prediction accuracy.

We consider the following four distribution settings to investigate the robustness and efficiency of the proposed method:

- (1) Normal distribution  $\mathcal{N}(0, \sigma^2)$  with mean zero and variance  $\sigma^2 > 0$ ;
- (2) Skewed generalized  $t$  distribution (Theodossiou, 1998)  $\text{sgt}(\mu, \sigma^2, \lambda, p, q)$ , where mean  $\mu = 0$ , variance  $\sigma^2 = q/(q-2)$  with  $q > 2$ , shape  $p = 2$ , and skewness  $\lambda = 0.75$ ;
- (3) Lognormal distribution  $\text{LN}(\mu, \sigma)$  with  $\mu = 0$  and  $\sigma > 0$ ; and
- (4) Pareto distribution  $\text{Par}(x_m, \alpha)$  with scale  $x_m = 1$  and shape  $\alpha > 0$ .

All of the above settings except (1) are skewed and might be very heavy-tailed for some choice of the distribution parameters, such as  $\alpha < 2$  for the Pareto distribution.

##### 4.1 Mean estimation

For each setting, we generate an independent sample of size  $n = 100$  and compute three mean estimators: the sample mean, the Huber estimator with  $\tau$  chosen using five-fold cross-validation (CV-Huber), and the proposed DA-Huber mean estimator. Figure 1 displays the  $\alpha$ -quantile of the estimation error, with  $\alpha$  ranging from 0.5 to 1 based on 2000 simulations. Figure S1 in the Supplementary Material shows box plots of the

#### 4.1 Mean estimation 26

estimation error. The DA-Huber estimator and sample mean perform almost identically for the normal data. For the heavy-tailed skewed distributions, the deviation of the sample mean from the population mean grows rapidly with the confidence level, in striking contrast to the CV- and DA-Huber estimators.

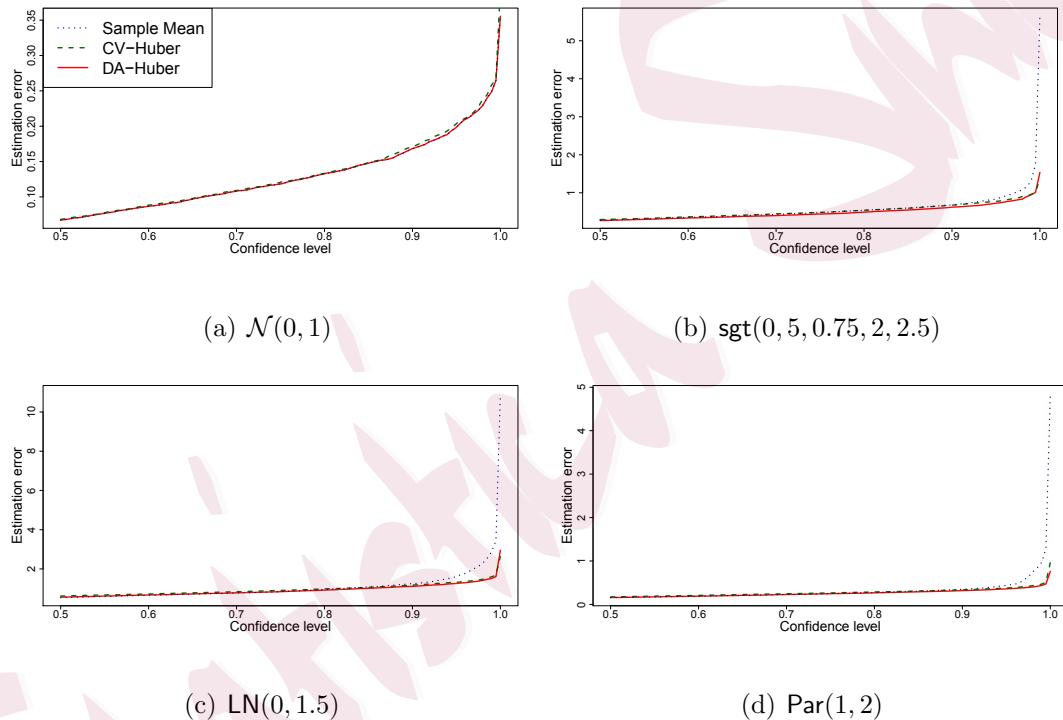


Figure 1: Estimation error versus confidence level for the sample mean, CV-Huber, and DA-Huber estimators based on 2000 simulations.

In Figure 2, we examine the 99%-quantile of the estimation error versus a distribution parameter measuring the tail behavior and the skewness. That is, for normal data we let  $\sigma$  vary between 1 and 4; for skewed generalized  $t$  distributions, we increase the shape parameter  $q$  from 2.5 to 4; for the lognormal and Pareto distributions, the shape

#### 4.1 Mean estimation 27

parameters  $\sigma$  and  $\alpha$  vary from 0.25 to 2 and 1.5 to 3, respectively. The Huber-type

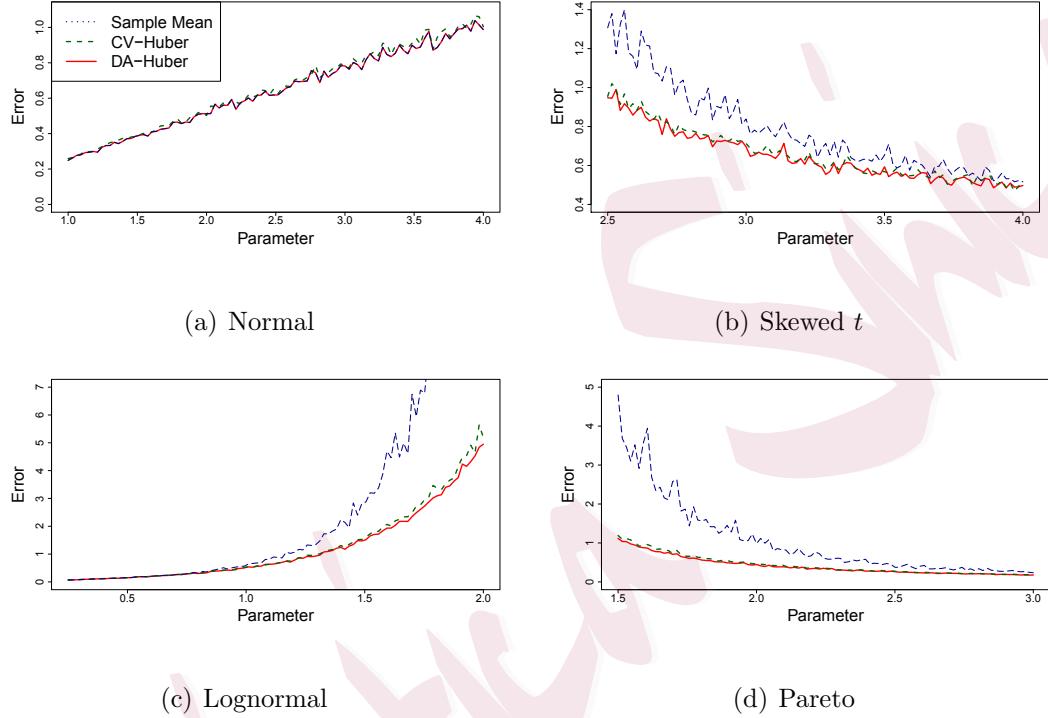


Figure 2: Empirical 99%-quantile of the estimation error versus a parameter measuring the tails and skewness for the sample mean, CV-Huber, and DA-Huber estimators.

estimators show substantial improvement in the deviations from the population mean because the distribution tends to have heavier tails and becomes more skewed. In summary, the most attractive feature of our method is its adaptivity: (i) it is as efficient as the sample mean for the normal model and is more robust for asymmetric and/or heavy-tailed data; (ii) it performs as well as the cross-validation method but with a much lower computational cost. The latter is particularly important for large-scale inferences, in which a myriad of parameters need to be estimated simultaneously.

## 4.2 Linear regression

We generate data  $\{(Y_i, \mathbf{X}_i)\}_{i=1}^n$  from the linear model in (3.1), with  $n = 500$  and  $d = 5$ . The intercept and the vector of regression coefficients are taken as  $\beta_0 = 5$  and  $\beta^* = (1, -1, 1, -1, 1)^\top$ , respectively. The covariates  $\mathbf{X}_i$  are *i.i.d.* random vectors that consist of independent coordinates from a uniform distribution  $\text{Unif}(-1.5, 1.5)$ .

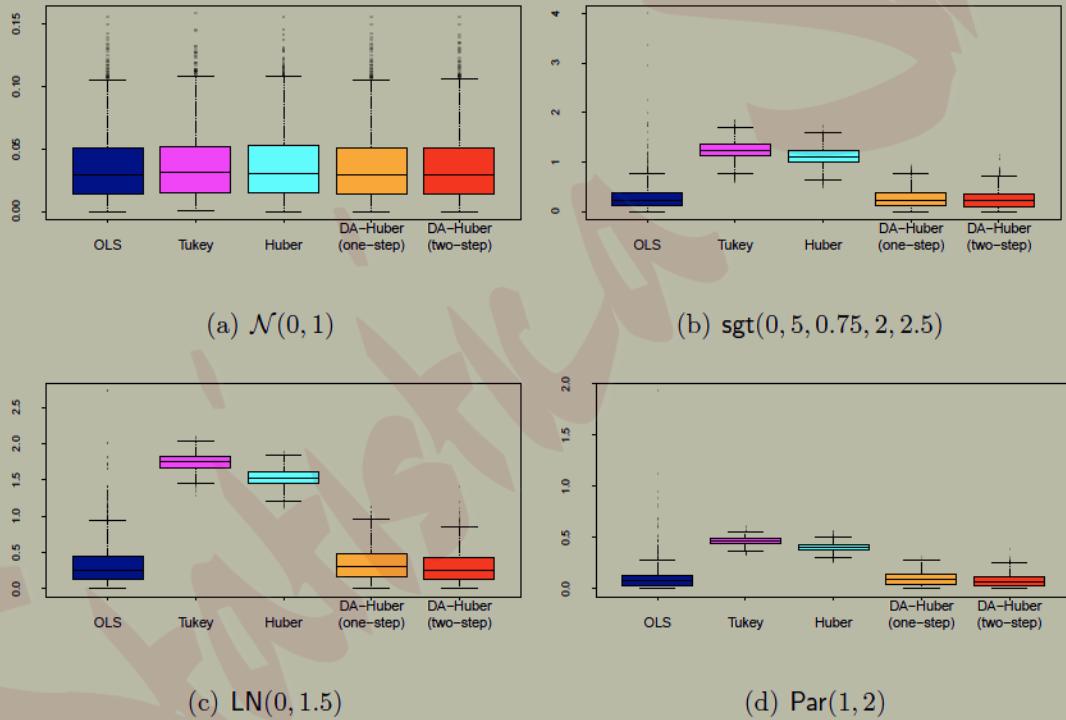


Figure 3: Estimation errors of intercept under different settings.

We compare the DA-Huber regression estimator with the ordinary least squares (OLS) estimator, and with classical robust  $M$ -estimators with a Huber loss  $\ell_\tau(\cdot)$ , as in (1.1), and Tukey's biweight loss  $\ell_\tau^T(x) = \{1 - (1 - x^2/\tau^2)^3\}\mathbb{I}(|x| \leq \tau) + \mathbb{I}(|x| > \tau)$ .

The tuning parameter  $\tau$  in  $\ell_\tau^T(\cdot)$  and  $\ell_\tau(\cdot)$  is taken as 4.685 and 1.345, respectively, according to the 95% efficiency rule. We carry out 1000 Monte Carlo simulations to (1) evaluate the overall performance of the DA-Huber methods by comparing it with that of the three competitors, OLS, Tukey, and Huber (see Figures 3 and 4), and (2) explore the robustness of different methods with varying degrees of heavy-tailedness and skewness (see Figures 5 and 6).

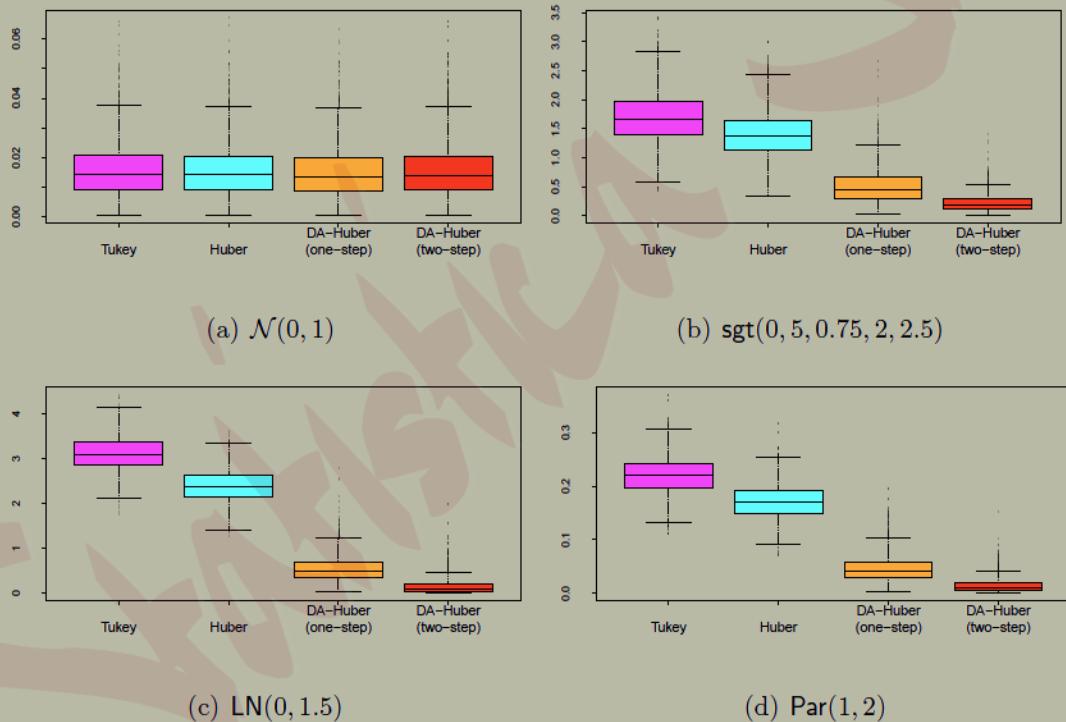


Figure 4: Total  $\ell_2$ -errors under different settings.

Figures 3 and 4 display box plots of the estimation error of the intercept  $|\hat{\beta}_0 - \beta_0^*|$  and the total  $\ell_2$ -error  $\|\hat{\theta} - \theta^*\|_2^2$ , respectively, for a fixed distribution parameter, as in Section 4.1. The one-step and two-step DA-Huber estimators both outperform

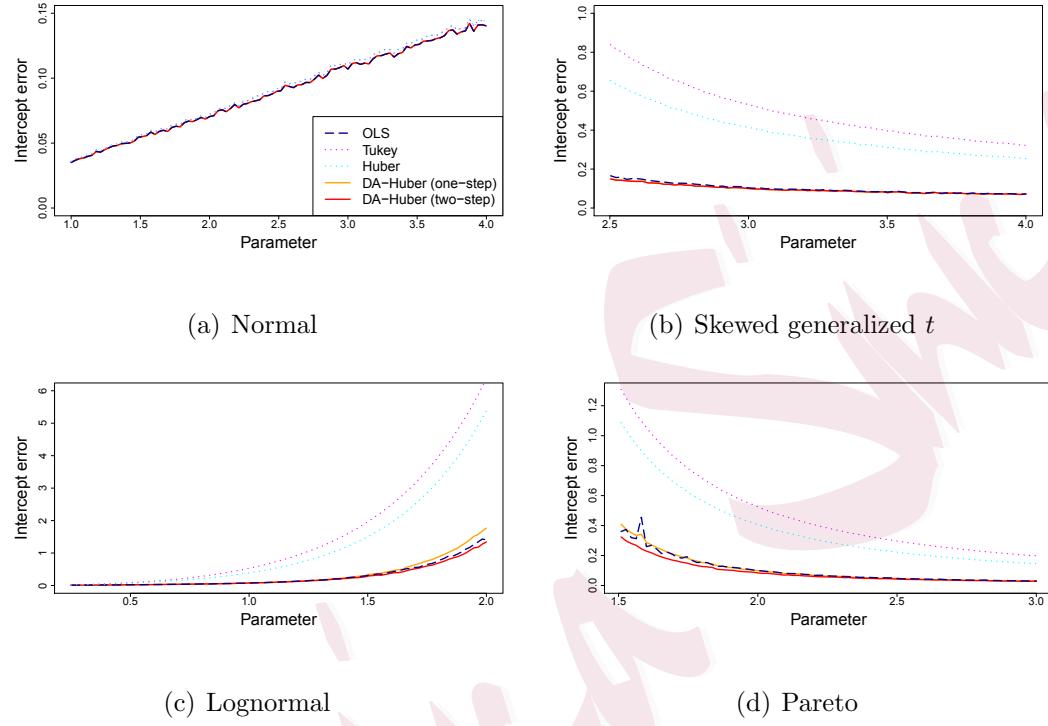


Figure 5: Average estimation error of the intercept versus the distribution parameters controlling the tails for the OLS estimator, standard Tukey and Huber estimators, and data-adaptive Huber estimators (one-step and two-step).

the other methods across all examples. When estimating the intercept, DA-Huber rectifies the non-negligible bias in the traditional robust  $M$ -estimator, as predicted by the theory. In the normal case, the DA-Huber estimator performs almost identically to the OLS estimator, and is therefore highly efficient. The  $\ell_2$ -error of the OLS tends to spread out (due to outliers), and thus is not reported. Figures 5 and 6 show the average estimation error of the intercept and the total  $\ell_2$ -error versus the distribution parameters controlling the shape of the tails, respectively. In the normal case, the

one-step DA-Huber and OLS slightly outperform the others. With heavy-tailed and skewed errors, the DA-Huber methods enjoy a notable advantage. However, the two-step approach is the most desirable because it strikes a good balance between bias and tail robustness. Overall, the numerical results confirm that the proposed methods have substantial advantages in the presence of asymmetric and heavy-tailed errors, while maintaining high efficiency for the normal model.

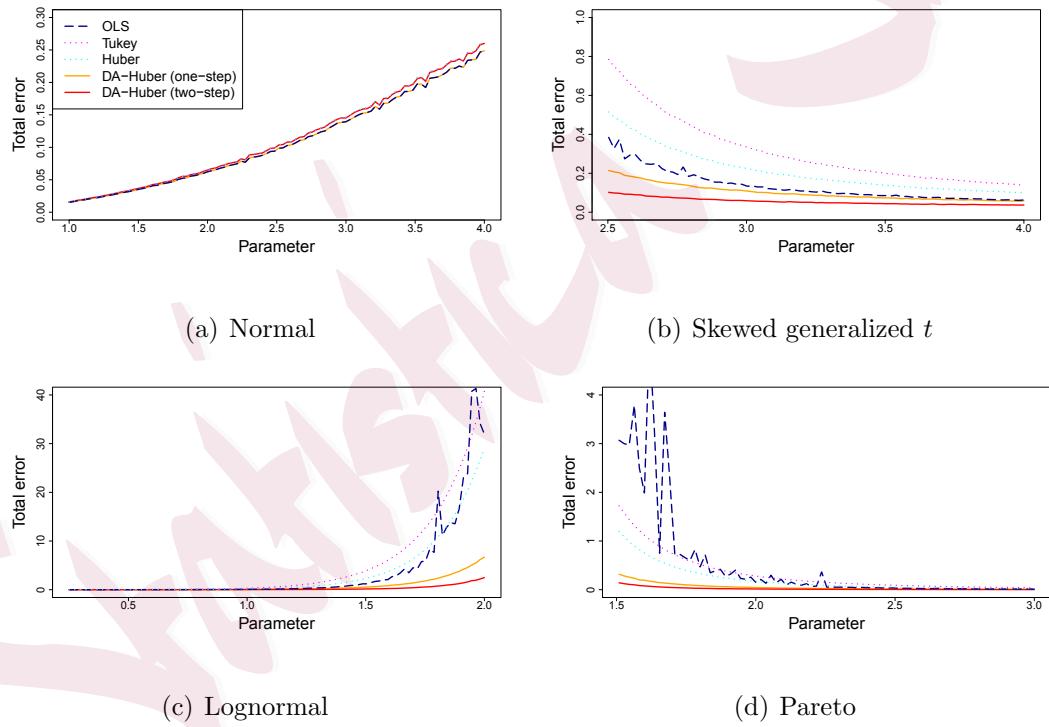


Figure 6: Average  $\ell_2$ -errors versus the distribution parameters controlling the tails for the OLS estimator, standard Tukey and Huber estimators, and data-adaptive Huber estimators (one-step and two-step).

### 4.3 Sparse linear regression32

#### 4.3 Sparse linear regression

Now, we consider the sparse linear regression,  $Y_i = \beta_0^* + \mathbf{X}_i^\top \boldsymbol{\beta}^* + \varepsilon_i$ , with  $i = 1, \dots, n$ , where  $\boldsymbol{\beta}^* \in \mathbb{R}^d$  is sparse, with  $s = \|\boldsymbol{\beta}^*\|_0 \ll n$  and  $d \gg n$ . In our simulations, we take  $n = 250$ ,  $d = 1000$ , and  $s = 20$ . We set  $\beta_0^* = 3$  and  $\boldsymbol{\beta}^* = (3, \dots, 3, 0, \dots, 0)^\top$ , where the first  $s = 20$  nonzero entries of  $\boldsymbol{\beta}^*$  are all equal to three. As before, the covariates  $\mathbf{X}_i$  are *i.i.d.* random vectors whose independent coordinates are from  $\text{Unif}(-1.5, 1.5)$ , and  $\varepsilon_i$  follows one of four distributions: normal, skewed generalized  $t$ , lognormal, and Pareto.

To implement the iterative procedure proposed in Section 3.2, at the  $k$ th iteration, we use five-fold cross-validation to choose  $\lambda_1^{(k)}$  and  $\lambda_2^{(k)}$  in the optimization programs in (3.11), producing  $\widehat{\boldsymbol{\theta}}_1^{(k)}$  and  $\widehat{\boldsymbol{\theta}}_2^{(k)}$ , respectively. We evaluate the proposed regularized DA-Huber estimators using the following measurements: RG, the relative gain of the DA-Huber estimator with respect to the Lasso in terms of the  $\ell_1$ - and  $\ell_2$ -errors;  $\text{RG}_q = \|\widehat{\boldsymbol{\theta}}_H - \boldsymbol{\theta}\|_q / \|\widehat{\boldsymbol{\theta}}_{\text{lasso}} - \boldsymbol{\theta}\|_q$ , with  $q = 1, 2$ ; FP, the number of false positives (selected noise covariates); and FN, the number of false negatives (missed signal covariates).

Table 1 summarizes the relative gains of the DA-Huber estimators under the  $\ell_1$ - and  $\ell_2$ -errors and the numbers of false positive and false negative discoveries. Across all four models, both DA-Huber estimators outperform the Lasso, with smaller  $\ell_1$ -errors and fewer false positive discoveries. Therefore, they are less greedy in terms of model selection. For the normal model, the proposed robust methods and the Lasso perform equally well. In the presence of heavy-tailed skewed errors, the DA-Huber methods lead

to remarkably better outputs in regard of both estimation and model selection. Similar results are observed in Figure S2 in the Supplementary Material, which displays the empirical distributions of the  $\ell_2$ -errors for all estimators.

Table 1: RG, FP, and FN and their standard errors (in parenthesis) of the Lasso and DA-Huber estimators under different models. The results are based on 200 simulations.

		DA-Huber		DA-Huber		
Lasso		(one-step)	(two-step)	Lasso	(one-step)	(two-step)
Normal, $\mathcal{N}(0, 1)$						sgt(0, 5, 0.75, 2, 2.5)
RG <sub>1</sub> × 100	100	93.4 (0.6)	91.4 (0.9)	100	87.5 (1.0)	86.2 (0.9)
RG <sub>2</sub> × 100	100	100.3 (0.2)	102.7 (0.3)	100	98.3 (0.5)	98.1 (0.5)
FP	87.9 (1.7)	77.6 (1.4)	73.5 (2.0)	86.1 (1.8)	63.1 (1.8)	60.7 (1.5)
FN	0 (0)	0 (0)	0 (0)	0 (0)	0 (0)	0 (0)
Lognormal, LN(0, 1.5)						Pareto, Par(1, 2)
RG <sub>1</sub> × 100	100	34.7 (0.7)	22.7 (0.5)	100	65.3 (1.1)	41.7 (0.8)
RG <sub>2</sub> × 100	100	49.5 (1.0)	30.5 (0.7)	100	84.5 (0.9)	51.2 (0.9)
FP	80.8 (2.0)	21.9 (0.6)	26.6 (0.7)	85.1 (1.9)	34.5 (1.6)	44.2 (0.9)
FN	0.26 (0.1)	0 (0)	0 (0)	0 (0)	0 (0)	0 (0)

## 5. Conclusion

We have proposed a new principle for choosing a robustification parameter adaptively from data for a variety of fundamental statistical problems, including mean estimations, a linear regression, and a sparse regression in high dimensions. Inspired by the censored moment equation approach, the proposed principle is tuning-free and data-adaptive. It is conceptually different from the traditional practice of selecting the robustification parameter using cross-validation, which is not only computationally demanding, but also lacks the underpinning mathematical guarantees. The proposed principle is guided by nonasymptotic deviation analysis, providing a unified method for choosing a robustification parameter for tail-robust estimation and inference. In particular, the analysis guiding our method can be extended easily to a broader class of robust convex loss functions, including the pseudo-Huber loss functions. The key is the global Lipschitz and local quadratic geometry of the loss function  $\ell_\tau(x) = \tau^2 \ell(x/\tau)$ . In light of the numerical evidence from both synthetic and real data, our proposal outperforms those widely known procedures in terms of estimation, variable selection, and prediction in the presence of heavy-tailed and skewed errors. Finally, an R package that implements the DA-Huber method can be found at <https://github.com/XiaoouPan/tfHuber>.

## Supplementary Material

The online Supplementary Material contains proofs of all theoretical results in the main text, as well as additional empirical studies.

### Acknowledgements

We thank the editor, the associate editor, and two referees for their insightful comments and suggestions. Wang's research was partially supported by NSFC grant 11701509 and the First Class Discipline of Zhejiang-A (Zhejiang Gongshang University-Statistics). Chao Zheng's research was supported by EPSRC Grant EP/N031938/1. Wen Zhou's research was partially supported by the Department of Energy. Wen-Xin Zhou's research was supported by NSF Grant DMS-1811376.

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