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In this paper, we consider the problem of identifying a linear map from measurements which are subject to intermittent and arbitarily large errors. This is a fundamental problem in many estimation-related applications such as fault detection, state estimation in lossy networks, hybrid system identification, robust estimation, etc. The problem is hard because it exhibits some intrinsic combinatorial features. Therefore, obtaining an effective solution necessitates relaxations that are both solvable at a reasonable cost and effective in the sense that they can return the true parameter vector. The current paper discusses a nonsmooth convex optimization approach and provides a new analysis of its behavior. In particular, it is shown that under appropriate conditions on the data, an exact estimate can be recovered from data corrupted by a large (even infinite) number of gross errors.

Analysis of A Nonsmooth Optimization Approach to Robust

Estimation \star

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Introduction

1.1Problem and motivations

We consider a linear measurement model of the form

$$y_t = x_t^\top \theta^o + f_t + e_t \tag{1}$$

where $y_t \in \mathbb{R}$ is the measured signal, $x_t \in \mathbb{R}^n$ the regression vector, $\{e_t\}$ a sequence of zero-mean and bounded errors (e.q., measurement noise, model mismatch, uncertainties, etc.) and $\{f_t\}$ a sequence of intermittent and arbitrarily large errors. Assume that we observe the sequences $\{x_t\}_{t=1}^N$ and $\{y_t\}_{t=1}^N$ and would like to compute the parameter vector θ^o from these observations. We are interested in doing so without knowing any of the sequences $\{f_t\}$ and $\{e_t\}$. We do however make the following assumptions:

• $\{e_t\}$ is a bounded sequence.

• $\{f_t\}$ is a sequence containing zeros and intermittent gross errors with (possibly) arbitrarily large magnitudes.

This is an important estimation problem arising in many situations such as fault detection [30,11], hybrid system identification [16], subspace clustering [40,2], error correction in communication networks [7]. The case when $\{f_t\}$ is zero and $\{e_t\}$ is a Gaussian process has been well-studied in linear system identification theory (see, e.q., the text books [23,38]). A less studied, but very relevant scenario in the system identification community, is when the additional perturbation $\{f_t\}$ in (1) is nonzero and contains intermittent and arbitrarily large errors [7,37,26,42]. It is worth noticing the difference with the problem studied in the field of compressive sensing [7,13,10]. In compressive sensing, the sought parameter vector is assumed sparse and the measurement noise $\{e_t\}$, often Gaussian or bounded. Here, no assumptions are made concerning sparsity of θ^{o} . We will, in this contribution, study essentially the case when the data is noise-free (*i.e.*, $e_t = 0$ for all t) and $\{f_t\}$ is a sequence with intermittent gross errors. We will derive conditions

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for perfect recovery and point to effective algorithms for computing θ^o . In the second part of the paper, the model assumption is relaxed to allow both e_t and f_t to be simultaneously nonzero. Note that this might be a more realistic scenario since most applications have measurement noise.

For illustrative purposes, let us discuss briefly some applications where a model of the form (1) is of interest.

Switched linear system identification. A discretetime Multi-Input Single-Output (MISO) Switched Linear System (SLS) can be written in the form

$$y_t = x_t^\top \theta_{\sigma_t}^o + e_t, \qquad (2)$$

where $x_t \in \mathbb{R}^n$ is the regressor at time $t \in \mathbb{Z}_+$ defined by

$$x_t = \begin{bmatrix} y_{t-1} \cdots y_{t-n_a} & u_t^\top & u_{t-1}^\top \cdots & u_{t-n_b}^\top \end{bmatrix}^\top, \quad (3)$$

where $u_t \in \mathbb{R}^{n_u}$ and $y_t \in \mathbb{R}$ denote respectively the input and the output of the system. The integers n_a and n_b in (3) are the maximum output and input lags (also called the orders of the system). $\sigma_t \in \{1, \ldots, s\}$ is the discrete mode (or discrete state) indexing the active subsystem at time t; it is in general assumed *unobserved*. $\theta_{\sigma_t}^o \in \mathbb{R}^n$, $n = n_a + n_b n_u$, is the parameter vector (PV) associated with the mode σ_t . For $\theta^o \in \{\theta_1^o, \ldots, \theta_s^o\}$, the Switched Auto-Regressive eXogenous (SARX) model (2) can be written in the form (1), with unknown f_t of the following structure $f_t = x_t^T (\theta_{\sigma_t}^o - \theta^o)$. For a background on hybrid system identification, we refer to the references [32,16,41,1,25,31,28].

Identification from faulty data. A model of the form (1) also arises when one has to identify a linear dynamic system which is subject to intermittent sensor faults. This is the case in general when the data are transmitted over a communication network [7,30]. Model (1) is suitable for such situations and the sequence $\{f_t\}$ then models occasional data packets losses or potential outliers. More precisely, a dynamic MISO system with process faults can be directly written in the form (1). In the case of sensor faults, the faulty model might be defined by

$$\begin{cases} \bar{y}_t = \bar{x}_t^\top \theta^o + e_t \\ y_t = \bar{y}_t + w_t \end{cases}$$

where $y_t \in \mathbb{R}$ is the observed output which is affected by the fault w_t (assumed to be nonzero only occasionally); \bar{x}_t is defined as in (3) from the known input u_t and the unobserved output \bar{y}_t . We can rewrite the faulty model exactly in the form (1) with $f_t = w_t - [w_{t-1} \cdots w_{t-n_a}]\theta^o$. Sparsity of $\{w_t\}$ induces sparsity of $\{f_t\}$ but in a lesser extent.

State estimation in the presence of intermittent errors. Considering a MISO dynamic system with state dynamics described by $z_{t+1} = Az_t + Bu_t$ and observation equation $\tilde{y}_t = C^{\top} z_t + f_t$, (A, B, C) being known matrices of appropriate dimensions, and $\{f_t\}$ a sparse sequence of possibly very large errors, the finite horizon state estimation problem reduces to the estimation of the initial state $z_0 = \theta$. We get a model of the form (1) by setting $y_t = \tilde{y}_t - C^{\top} \Delta_t \bar{u}_t$ and $x_t = (A^t)^{\top} C$, with $\Delta_t = [A^{t-1}B \cdots AB B], \ \bar{u}_t = [u_0^{\top} \cdots u_{t-1}^{\top}]^{\top}$. Examples of relevant works are those reported in [3,15]. In this latter application, it can however be noted that the dataset $\{x_t\}$ may not be generic enough. ¹

Connection to robust statistics. Indeed, the problem of identifying the parameters from model (1) under the announced assumptions can be viewed as a robust regression problem where the nonzero elements in the sequence $\{f_t\}$ are termed outliers. As such, it has received a lot of attention in the robust statistics literature (see, e.g., [21,35,24] for an overview). Examples of methods to tackle the robust estimation problem include the least absolute deviation [20], the least median of squares [34], the least trimmed squares [35], the M-estimator [21], etc. Most of these estimators come with an analysis in terms of the breakdown point [19,36], a measure of the (asymptotic) minimum proportion of points which cause an estimator to be unbounded if they were to be arbitrarily corrupted by gross errors. The current report focuses on the analysis of a nonsmooth convex optimization approach which includes the least absolute deviation method as a particular case corresponding to the situation when the output in (1) is a scalar. The analysis approach taken in the current paper is different in the following sense.

- In robust statistics the quality of an estimator is measured by its breakdown point. The higher the breakdown point, the better. The available analysis is therefore directed to determining a sort of absolute robustness: how many outliers (expressed in proportion of the total number of samples) cause the estimator to become unbounded.
- Here, the question of robust performance of the estimator is posed differently. We are interested in estimating the maximum number of outliers that a nonsmooth-optimization-based estimator can accommodate while still returning the exact value one would obtain in the absence of any outlier. This is more related to the traditional view developed in compressive sensing.

Contributions of this paper. One promising method for estimating model (1) is by nonsmooth convex optimization as suggested in [7,37,1,26,42]. More precisely, inspired by the recent theory of compressed sensing [7,13,10], the idea is to minimize a nonsmooth (and non differentiable) sum-of-norms objective function involv-

¹ In this paper, the term genericity for a dataset characterizes a notion of linear independence. For example, a set of N > n data points in general linear position in \mathbb{R}^n is more generic than a set of data points contained in one subspace. We will introduce different quantitative measures of data genericity in the sequel (see Definition 2 and Theorem 11).

ing the fitting errors. Under noise-free assumption, such a cost function has the nice property that it is able to provide the true parameter vector in the presence of arbitrarily large errors $\{f_t\}$ provided that the number of nonzero errors is small in some sense. Of course, when the data are corrupted simultaneously by the noise $\{e_t\}$ and the gross errors $\{f_t\}$, the recovery cannot be exact any more. It is however expected (as Proposition 17 and simulations tend to suggest) that the estimate will still be close to the true one.

The current paper intends to present a new analysis of the nonsmooth optimization approach and provide some elements for further understanding its behavior. The line of analysis goes from a full characterization of the nonsmooth optimization based estimator (both for SISO and MIMO systems) to the study of robustness to outliers including in the presence of dense noise. With respect to relevant works [7,37,1,26,42], we derive new bounds on the number of outliers (in the least favorable situations) that the estimator is capable to accommodate. It is emphasized that a quite broad spectrum of such bounds can be derived based on the basic characterization of the nonsmooth identifier. Note however that evaluating numerically the tightest of these bounds is a high computational process while less tight bounds have a more affordable complexity. Some of the bounds developed in this contribution meet both relative tightness requirement and computability in polynomial time (see the bound based on $\xi(X)$ in Theorem 11). Finally, the paper show how the results derived in the first part for ℓ_1 -norm estimator when applied to the estimation of SISO systems are generalizable to multivariable systems.

Outline of this paper. The outline of the paper is as follows. We start in Section 2.1 by viewing the nonsmooth optimization as the convex relaxation of a (ideal) combinatorial ℓ_0 -"norm" formulation. We then derive in Section 2.2 necessary and sufficient conditions for optimality. Based on those conditions we establish in Section 2.4 new sufficient conditions for exact recovery of the true parameter vector in (1). The noisy case is treated in Section 3.2. Section 4 presents a generalization of the earlier discussions to multi-output systems. Finally, numerical experiments are described in Section 5 and concluding remarks are given in Section 6.

1.2 Notations

Let $\mathbb{I} = \{1, \ldots, N\}$ be the index set of the measurements. For any $\theta \in \mathbb{R}^n$, define a partition of the set of indices \mathbb{I} by $\mathbb{I}^-(\theta) = \{t \in \mathbb{I} : \theta^\top x_t - y_t < 0\}$, $\mathbb{I}^+(\theta) = \{t \in \mathbb{I} : \theta^\top x_t - y_t > 0\}$, $\mathbb{I}^0(\theta) = \{t \in \mathbb{I} : \theta^\top x_t - y_t = 0\}$. *Cardinality of a finite set.* Throughout the paper, whenever S is a finite set, the notation |S| will refer to the cardinality of S. However, for a real number x, |x| will denote the absolute value of x.

Submatrices and subvectors. Let $X = \begin{bmatrix} x_1 & x_2 & \cdots & x_N \end{bmatrix} \in \mathbb{R}^{n \times N}$ be the matrix formed with the available regressors $\{x_t\}_{t=1}^N$. If $I \subset \mathbb{I}$, the notation X_I denotes a matrix

in $\mathbb{R}^{n \times |I|}$ formed with the columns of X indexed by I. Likewise, with $\boldsymbol{y} = \begin{bmatrix} y_1 \ y_2 \ \cdots \ y_N \end{bmatrix}^\top \in \mathbb{R}^N$, \boldsymbol{y}_I is the vector in $\mathbb{R}^{|I|}$ formed with the entries of \boldsymbol{y} indexed by I. We will use the convention that $X_I = 0 \in \mathbb{R}^n$ (resp. $\boldsymbol{y}_I = 0 \in \mathbb{R}$) when the index set I is empty.

Vector norms. $\|\cdot\|_p$, $p = 1, 2, ..., \infty$, denote the usual *p*-norms for vectors defined for any vector $z = [z_1 \cdots z_N]^\top \in \mathbb{R}^N$, by $\|z\|_p = (|z_1|^p + \cdots + |z_N|^p)^{1/p}$. Note that $\|z\|_{\infty} = \max_{i=1,...,N} |z_i|$. The ℓ_0 "norm" of z is defined to be the number of nonzero entries in z, i.e., $\|z\|_0 = |\{i : z_i \neq 0\}|$.

Matrix norms. The following matrix norms will be used: $\|\cdot\|_p, p = 1, 2, ..., \infty, \|\cdot\|_{2, \text{col}}, \|\cdot\|_{2, \infty}$. They are defined as follows: for a matrix $A = [a_1 \cdots a_N] \in \mathbb{R}^{n \times N}$ with $a_i \in \mathbb{R}^n$,

$$\begin{aligned} \|A\|_{p} &= \sup_{x \in \mathbb{R}^{N}, \|x\|_{p}=1} \|Ax\|_{p}, \quad \|A\|_{2, \text{col}} = \sum_{i=1}^{N} \|a_{i}\|_{2}, \\ \|A\|_{2, \infty} &= \max_{i=1, \dots, N} \|a_{i}\|_{2}. \end{aligned}$$

2 Nonsmooth optimization for the estimation problem

2.1 Sparse optimization

The main idea for identifying the parameter vector θ^{o} from (1) is by solving a sparse optimization problem, that is, a problem which involves the minimization of the number of nonzeros entries in the error vector. To be more specific, assume for the time being that the error sequence $\{e_t\}$ is identically equal to zero. Consider a candidate parameter vector $\theta \in \mathbb{R}^n$ and let

$$\phi(\theta) = \boldsymbol{y} - \boldsymbol{X}^{\top} \boldsymbol{\theta},$$

where $\boldsymbol{y} = \begin{bmatrix} y_1 \cdots y_N \end{bmatrix}^{\top}$, $X = \begin{bmatrix} x_1 \cdots x_N \end{bmatrix}$, be the fitting error vector induced by θ on the experimental data. Then the vector θ^o can naturally be searched for by minimizing an ℓ_0 objective function,

$$\underset{\theta \in \mathbb{R}^n}{\operatorname{minimize}} \left\| \phi(\theta) \right\|_0 \tag{4}$$

where $\|\cdot\|_0$ denotes the ℓ_0 pseudo-norm which counts the number of nonzero entries. Because problem (4) aims at making the error $\phi(\theta)$ sparse by minimizing the number of nonzero elements (or maximizing the number of zeros), it is sometimes called a sparse optimization problem [1]. As can be intuitively guessed, the recoverability of the true parameter vector θ^o from (4) depends naturally on some properties of the available data. This is outlined by the following lemma.

Lemma 1 (A sufficient condition for ℓ_0 recovery) Assume that $\{e_t\}$ is equal to zero and let $\boldsymbol{f} = [f_1 \cdots f_N]^\top$. Assume that for any $I \subset \mathbb{I}$ with |I| > n, $\boldsymbol{f}_I \notin \operatorname{im}(X_I^\top)$ whenever $\boldsymbol{f}_I \neq 0$, with $\operatorname{im}(\cdot)$ referring here to range space. Then provided $|\mathbb{I}^{0}(\theta^{o})| > n$, it holds that

$$\theta^{o} \in \underset{\theta}{\operatorname{arg\,min}} \|\phi(\theta)\|_{0} \,. \tag{5}$$

PROOF. We proceed by contradiction. Assume that (5) does not hold, *i.e.*, $\min_{\theta} \|\phi(\theta)\|_{0} < \|\phi(\theta^{o})\|_{0}$. Then, by letting θ^{m} be any vector in $\arg\min_{\theta} \|\phi(\theta)\|_{0}$, the above inequality translates into $|\mathbb{I}^{0}(\theta^{m})| > |\mathbb{I}^{0}(\theta^{o})| > n$. It follows that $\mathbf{f}_{\mathbb{I}^{0}(\theta^{m})} \neq 0$ because $|\mathbb{I}^{0}(\theta^{o})| = |\{t \in \mathbb{I} : f_{t} = 0\}|$ is the exact (largest) number of zero elements in the sequence $\{f_{t}\}_{t=1}^{N}$. Note also that we have necessarily $|\mathbb{I}^{0}(\theta^{m})| > n$. On the other hand, with $\mathbf{y}_{\mathbb{I}^{0}(\theta^{m})} = X_{\mathbb{I}^{0}(\theta^{m})}^{\top} \theta^{m}} = X_{\mathbb{I}^{0}(\theta^{m})}^{\top} (\theta^{m} - \theta^{o}) \in \operatorname{im}(X_{\mathbb{I}^{0}(\theta^{m})}^{\top})$. This, together with $\mathbf{f}_{\mathbb{I}^{0}(\theta^{m})} \neq 0$, constitutes a contradiction to the assumption of the Lemma. Hence, (5) holds as claimed. \Box

Lemma 1 specifies a condition involving both X and fand under which θ^o lies in the solution set but it does not ensure that θ^o will be recovered uniquely from data. Before proceeding further, we recall from [1] a sufficient condition under which θ^o is the unique solution to (4). **Definition 2** ([1] An integer measure of genericity)

Let $X \in \mathbb{R}^{n \times N}$ be a data matrix satisfying rank(X) = n. The n-genericity index of X denoted $\nu_n(X)$, is defined as the minimum integer m such that any $n \times m$ submatrix of X has rank n,

$$\nu_n(X) = \min\left\{m : \forall \mathcal{S} \subset I \text{ with } |\mathcal{S}| = m, \operatorname{rank}(X_{\mathcal{S}}) = n\right\}.$$
(6)

Theorem 3 ([1] Sufficient condition for ℓ_0 **recovery)** Assume that the sequence $\{e_t\}$ in (1) is identically equal to zero. If the sequence $\{f_t\}$ in (1) contains enough zero values in the sense that

$$\left|\mathbb{I}^{0}(\theta^{o})\right| = \left|\left\{t \in \mathbb{I} : f_{t} = 0\right\}\right| \ge \frac{N + \nu_{n}(X)}{2}, \quad (7)$$

then θ^{o} is the unique solution to the ℓ_{0} -norm minimization problem (4).

In other words, if the number of nonzero gross errors $\{f_t\}$ affecting the data generated by (1) does not exceed the threshold $(N - \nu_n(X))/2$, then θ^o can be exactly recovered by solving (4). Unfortunately, this problem is a hard combinatorial optimization problem. A tractable solution can be obtained by relaxing the ℓ_0 -norm into its best convex approximant, the ℓ_1 -norm. Doing this substitution in (4) gives

$$\underset{\theta \in \mathbb{R}^n}{\operatorname{minimize}} \|\phi(\theta)\|_1 \tag{8}$$

with $\|\phi(\theta)\|_1 = \sum_{t=1}^N |y_t - \theta^\top x_t|$. The latter problem is termed a nonsmooth convex optimization problem [27, Chap. 3] because the objective function is convex but

non-differentiable. Compared to (4), problem (8) has the advantage of being convex and can hence be efficiently solved by many existing numerical solvers, *e.g.*, [18]. Note further that it can be written as a linear programming problem. The ℓ_1 relaxation process has been intensively used in the compressed sensing literature [14] for approaching the sparsest solution of an underdetermined set of linear equations. In the robust statistics literature as surveyed above, (8) corresponds to a wellknown estimator referred to as the least absolute deviation estimator [34]. As will be shown next, the underlying reason why problem (8) can obtain the true parameter vector despite the presence of gross perturbations $\{f_t\}$ is related to its nonsmoothness.

2.2 Solution to the ℓ_1 problem

There is a wealth of analysis in the literature of compressed sensing investigating under which conditions some problems² of similar structure as (4) and (8) can yield the same solution. This analysis is mainly based on the concepts of mutual coherence [14] and the Restricted Isometry Property [8]. Here, we shall propose a parallel but different analysis for the robust estimation problem. We start by characterizing the solution to the ℓ_1 -norm problem (8).

Theorem 4 (Solution to the ℓ_1 **problem)** A vector $\theta^* \in \mathbb{R}^n$ solves the ℓ_1 -norm problem (8) if and only if any of the following equivalent statements hold:

S1. There exist some numbers $\lambda_t \in [-1, 1], t \in \mathbb{I}^0(\theta^*)$, such that³

$$\sum_{t \in \mathbb{I}^+(\theta^*)} x_t - \sum_{t \in \mathbb{I}^-(\theta^*)} x_t = \sum_{t \in \mathbb{I}^0(\theta^*)} \lambda_t x_t.$$
(9)

S2. For any $\eta \in \mathbb{R}^n$,

$$\left|\sum_{t\in\mathbb{I}^+(\theta^*)}\eta^\top x_t - \sum_{t\in\mathbb{I}^-(\theta^*)}\eta^\top x_t\right| \le \sum_{t\in\mathbb{I}^0(\theta^*)}|\eta^\top x_t|.$$
(10)

S3. The optimal value of the optimization problem

$$\min_{\alpha} \|\alpha\|_{\infty} \quad subject \ to \ z = X_{\mathbb{I}^0(\theta^\star)}\alpha, \tag{11}$$

where
$$z = \sum_{t \in \mathbb{I}^+(\theta^*)} x_t - \sum_{t \in \mathbb{I}^-(\theta^*)} x_t, \ \alpha \in \mathbb{R}^{|\mathbb{I}^0(\theta^*)|}$$
, is less than or equal to 1.

Moreover, the solution θ^* is unique if and only if any of the following statements is true:

 $^{^2}$ Those problems look for the sparsest solution to an underdetermined set of linear equations. As such they are similar but different to the problem studied in the current paper. Note that the process of converting problems (4) and (8) into the format treated in compressed sensing yields a system of linear equations which is much less underdetermined.

 $^{^{3}}$ Eq. (9) should be understood here with the implicit convention that any of the three terms is equal to zero whenever the corresponding index set is empty.

S1'. (9) holds and $\operatorname{rank}(X_S) = n$ where

$$S = \left\{ t \in \mathbb{I}^0(\theta^\star) : |\lambda_t| < 1 \right\}.$$

S2'. (10) holds with strict inequality symbol for all $\eta \in \mathbb{R}^n$, $\eta \neq 0$.

PROOF. Proof of S1. Since $\|\phi(\theta)\|_1$ is a proper convex function, it has a non empty subdifferential [33]. The necessary and sufficient condition for θ^* to be a solution of (8) is then

$$0 \in \partial \|\phi(\theta^{\star})\|_{1}, \qquad (12)$$

where the notation ∂ refers to subdifferential with respect to θ . Indeed, using additivity of subdifferentials, it is straightforward to write

$$\partial \|\phi(\theta^{\star})\|_{1} = \sum_{t \in \mathbb{I}^{+}(\theta^{\star})} x_{t} - \sum_{t \in \mathbb{I}^{-}(\theta^{\star})} x_{t} + \sum_{t \in \mathbb{I}^{0}(\theta^{\star})} \operatorname{conv}\left\{-x_{t}, x_{t}\right\}$$
(13)

where conv refers to the convex hull. Here, the addition symbol is meant in the Minkowski sum sense. It follows that $0 \in \partial \|\phi(\theta^*)\|_1$ is equivalent to the existence of a set of numbers λ_t in $[-1, 1], t \in \mathbb{I}^0(\theta^*)$, such that (9) holds. **Proof of S2.** Define two functions $q, h : \mathbb{R}^n \to \mathbb{R}_{\geq 0}$ by $q(\theta) = \sum_{t \notin \mathbb{I}^0(\theta^*)} |y_t - \theta^\top x_t|$ and $h(\theta) = \sum_{t \in \mathbb{I}^0(\theta^*)} |y_t - \theta^\top x_t|$. Then $\|\phi(\theta)\|_1 = q(\theta) + h(\theta)$ and q is differentiable at θ^* . It follows that $\partial \|\phi(\theta^*)\|_1 = \nabla q(\theta^*) + \partial h(\theta^*)$, where $\nabla q(\theta^*)$ is the gradient of q at θ^* . We can hence write

$$\begin{aligned} \theta^{\star} \text{ minimizes } \|\phi(\theta)\|_1 \Leftrightarrow 0 \in \partial \|\phi(\theta^{\star})\|_1 \\ \Leftrightarrow -\nabla q(\theta^{\star}) \in \partial h(\theta^{\star}). \end{aligned}$$

Note from (13) that $\partial h(\theta^*) = \sum_{t \in \mathbb{I}^0(\theta^*)} \operatorname{conv} \{-x_t, x_t\}$ so that $-\nabla q(\theta^*) \in \partial h(\theta^*)$ if and only if $\pm \nabla q(\theta^*) \in \partial h(\theta^*)$ and this in turn is equivalent to $g^{\top}(\theta - \theta^*) \leq h(\theta) - h(\theta^*) \forall \theta$, for $g \in \{-\nabla q(\theta^*), +\nabla q(\theta^*)\}$. It follows that θ^* minimizes $\|\phi(\theta)\|_1$ if and only if

$$\left|\nabla q(\theta^{\star})^{\top}(\theta - \theta^{\star})\right| \le h(\theta) - h(\theta^{\star}) = \sum_{t \in \mathbb{I}^{0}(\theta^{\star})} \left|(\theta - \theta^{\star})^{\top} x_{t}\right|$$
(14)

for all θ . The last equality is obtained by using the fact that $y_t - x_t^\top \theta^* = 0$ for all t in $\mathbb{I}^0(\theta^*)$. Finally the result follows by setting $\eta = \theta - \theta^*$ and noting that $\nabla q(\theta^*) = \sum_{t \in \mathbb{I}^+(\theta^*)} x_t - \sum_{t \in \mathbb{I}^-(\theta^*)} x_t$.

 $S1 \Leftrightarrow S3$. The proof of the last equivalence is immediate.

Uniqueness. For convenience, we first prove S2'. Along the lines of the proof of S2 (see Eq. (14) and preceding arguments), we can see that strict inequality in (10) is equivalent to the following strict inequality $-\nabla q(\theta^*)^\top (\theta - \theta^*) < h(\theta) - h(\theta^*) \quad \forall \theta \neq \theta^*$. On the other hand, $\nabla q(\theta^*)^\top (\theta - \theta^*) \leq q(\theta) - q(\theta^*) \quad \forall \theta$. Summing the two yields

$$\left\|\phi(\theta^{\star})\right\|_{1} = q(\theta^{\star}) + h(\theta^{\star}) < q(\theta) + h(\theta) = \left\|\phi(\theta)\right\|_{1} \, \forall \theta \neq \theta^{\star}.$$

Hence S2' is proved.

For the proof of S1', we proceed in two steps.

Sufficiency. Assume rank $(X_S) = n$. Then for any nonzero vector $\eta \in \mathbb{R}^n$ there is at least one $t_0 \in S$ such that $\eta^{\top} x_{t_0} \neq 0$. Recall that $|\lambda_{t_0}| < 1$ by definition of S. It follows that by multiplying (9) on the left by η^{\top} with $\eta \in \mathbb{R}^n$ an arbitrary nonzero vector, and taking the absolute value yields (10) with strict inequality symbol. We can therefore apply the proof of S2' to conclude that θ^* is unique.

Necessity. Assume rank $(X_S) < n$. Then pick any nonzero vector η in ker (X_S^{\top}) . Set $\eta_1 = \nu \eta$ with $\nu \neq 0$. Indeed ν can be chosen sufficiently small such that $x_t^{\top}(\theta^* + \eta_1) - y_t$ has the same sign as $x_t^{\top}\theta^* - y_t$ for $t \in \mathbb{I}^-(\theta^*) \cup \mathbb{I}^+(\theta^*)$. For such values of ν we have $\mathbb{I}^+(\theta^*) \subset \mathbb{I}^+(\theta^* + \eta_1)$ and $\mathbb{I}^-(\theta^*) \subset \mathbb{I}^-(\theta^* + \eta_1)$. Moreover, since $\eta_1 \in \ker(X_S^{\top}), x_t^{\top}(\theta^* + \eta_1) - y_t = \eta_1^{\top}x_t = 0$ $\forall t \in S$, so that $S \subset \mathbb{I}^0(\theta^* + \eta_1)$. Finally, it remains to re-assign the indices t contained in $\mathbb{I}^0(\theta^*) \setminus S$ for which $\lambda_t = 1$. We get the following partition $\mathbb{I}^+(\theta^* + \eta_1) = \mathbb{I}^+(\theta^*) \cup \{t \in \mathbb{I}^0(\theta^*) : \eta_1^{\top}x_t > 0\},$ $\mathbb{I}^-(\theta^* + \eta_1) = \mathbb{I}^-(\theta^*) \cup \{t \in \mathbb{I}^0(\theta^*) : \eta_1^{\top}x_t < 0\},$ $\mathbb{I}^0(\theta^* + \eta_1) = S \cup \{t \in \mathbb{I}^0(\theta^*) : \eta_1^{\top}x_t = 0\}$. It follows that $\theta^* + \eta_1 \neq \theta^*$ also satisfies (9) with the sequence $\{\lambda_t\}_{t \in S}$ and is therefore a minimizer. In conclusion, if rank $(X_S) < n$, the minimizer cannot be unique. \Box

A number of important comments follow from Theorem 4.

- One first consequence of the theorem is that θ^{o} can be computed exactly from a finite set of erroneous data (by solving problem (8)) provided it satisfies the conditions S1' or S2' of the theorem. Note that there is no explicit boundedness condition imposed on the error sequence $\{f_t\}$. Hence the nonzero errors in this sequence can have arbitrarily large magnitudes as long as the optimization problem makes sense, *i.e.*, provided $\|\phi(\theta^*)\|_1$ remains finite.
- Second, the true parameter vector θ^o can be exactly recovered in the presence of, say, infinitely many nonzero errors f_t (see also Proposition 6). For example, if the regressors $\{x_t\}$ satisfy

$$\sum_{t \in \mathbb{I}^+(\theta^o)} x_t - \sum_{t \in \mathbb{I}^-(\theta^o)} x_t = 0,$$

and rank $(X_{\mathbb{I}^0(\theta^o)}) = n$, then by condition S2' θ^o is the unique solution to problem (8) regardless of the number of errors affecting the data.

• Third, if problem (8) admits a solution θ^* that satisfies $y_t - x_t^\top \theta^* \neq 0$ for all t = 1, ..., N, then θ^* is non-unique. In effect, $\mathbb{I}^0(\theta^*) = \emptyset$ in this case and so, rank $(X_{\mathbb{I}^0}(\theta^*)) = 0 < n$ which, by Theorem 4, implies non-uniqueness. Indeed this is typically the case whenever the noise $\{e_t\}$ is nonzero.

Another immediate consequence of Theorem 4 can be stated as follows.

Corollary 5 (On the special case of affine model) If the model (1) is affine in the sense that the regressor x_t has the form $x_t = [\tilde{x}_t^\top \ 1]^\top$, with $\tilde{x}_t \in \mathbb{R}^{n-1}$, then a necessary condition for θ^* to solve problem (8) is that

$$\left| \left| \mathbb{I}^+(\theta^*) \right| - \left| \mathbb{I}^-(\theta^*) \right| \right| \le \left| \mathbb{I}^0(\theta^*) \right|.$$
 (15)

Here, the outer bars $|\cdot|$ refer to absolute value while the inner ones which apply to sets refer to cardinality.

PROOF. The proof is immediate by considering the condition (10) and taking $\eta = [\mathbf{0}^{\top} \ \mathbf{1}]^{\top} \in \mathbb{R}^n$. \Box

Eq. (15) implies that if the measurement model is affine and all the f_t 's have the same sign, *i.e.*, if one of the cardinalities $|\mathbb{I}^+(\theta^o)|$ or $|\mathbb{I}^-(\theta^o)|$ is equal to zero, then problem (8) cannot find the true θ^o whenever more than 50% of the elements of the sequence $\{f_t\}$ are nonzero. Next, we discuss a special case in which the true parameter vector θ^o in (1) can, in principle, be obtained asymptotically in the presence of an infinite number of nonzero errors f_t 's.

Proposition 6 (Infinite number of outliers) Assume that the error sequence $\{e_t\}$ in (1) is identically equal to zero. Assume further that the data $\{(x_t, y_t)\}_{t=1}^N$ are generated such that:

- There is a set $I^0 \subset \mathbb{I}$ with $|I^0| \ge n$, such that for any $t \in I^0$, $f_t = 0$ and $\operatorname{rank}(X_{I^0}) = n$,
- For any $t \notin I^0$, f_t is sampled from a distribution which is symmetric around zero.
- The regression vector sequence {x_t} ⊂ ℝⁿ is drawn from a probability distribution having a finite first moment.

Then

$$\lim_{N \to \infty} \operatorname*{arg\,min}_{\theta \in \mathbb{R}^n} \frac{1}{N} \sum_{t=1}^{N} \left| y_t - x_t^\top \theta \right| = \{\theta^o\}$$
(16)

with probability one.

PROOF. Under the conditions of the proposition, we have $\operatorname{Prob}(y_t - x_t^\top \theta^o < 0) = \operatorname{Prob}(y_t - x_t^\top \theta^o > 0) = 1/2$, where Prob denotes probability measure. It follows that $|\mathbb{I}^+(\theta^o)|$ and $|\mathbb{I}^-(\theta^o)|$ go jointly to infinity as the total number of samples N tends to infinity. Hence, the expressions $\frac{1}{|\mathbb{I}^+(\theta^o)|} \sum_{t \in \mathbb{I}^+(\theta^o)} x_t$ and $\frac{1}{|\mathbb{I}^-(\theta^o)|} \sum_{t \in \mathbb{I}^-(\theta^o)} x_t$ are both sample estimates for the expectation of the process $\{x_t\}$. By the law of large numbers, as $N \to \infty$, the two quantities converge to the true expectation of the process $\{x_t\}$ with probability one, so that

$$\lim_{N \to \infty} \left[\frac{1}{|\mathbb{I}^+(\theta^o)|} \sum_{t \in \mathbb{I}^+(\theta^o)} x_t - \frac{1}{|\mathbb{I}^-(\theta^o)|} \sum_{t \in \mathbb{I}^-(\theta^o)} x_t \right] = 0.$$

As a consequence, θ^o satisfies condition S1' of Theorem 4 asymptotically with $\lambda_t = 0$ for any $t \in \mathbb{I}^0(\theta^o) = I^0$. Hence the solution of the ℓ_1 minimization problem tends to θ^o with probability one as the number of samples approaches infinity. \Box

2.3 Worst-case necessary and sufficient conditions

The conditions (9)-(11) derived in Theorem 4 characterize completely the solution to the ℓ_1 -problem. However such conditions depend on which data points (x_t, y_t) are affected by the gross errors and on the sign of the f'_ts . We wish now to find necessary and sufficient conditions that depend solely on the number of gross errors (or, equivalently on the number of zero elements in the sequence $\{f_t\}$).

Corollary 7 (Necessary and sufficient conditions) Let d be an integer. Then the following statements are equivalent: (i)

$$\forall \theta \in \mathbb{R}^{n}, \forall \boldsymbol{y} \in \mathbb{R}^{N}, |\mathbb{I}^{0}(\theta)| \geq d \Rightarrow \theta \in \operatorname*{arg\,min}_{w \in \mathbb{R}^{n}} \|\phi(w)\|_{1}$$
 (17)

$$\max_{\substack{(I,I^c):\\|I|=d}} \max_{\eta \in \mathbb{R}^n} \left\{ \left\| X_{I^c}^\top \eta \right\|_1 \text{ s. t. } \left\| X^\top \eta \right\|_1 = 1 \right\} \le 1/2$$
(18)

(iii)

$$\max_{\substack{(I,I^c):\\|I|=d}} \max_{h\in\{\pm 1\}^{|I^c|}} \min_{\alpha\in\mathbb{R}^{|I|}} \left\{ \|\alpha\|_{\infty} \text{ s. t. } X_{I^c}h = X_I\alpha \right\} \le 1$$
(19)

In (18)-(19) and similar equations in the paper, the leftmost maximum is taken over the set of those partitions (I, I^c) of \mathbb{I} that satisfy |I| = d. Eq. (19) should be read with the implicit assumption that the inequality fails to hold whenever the optimization problem is not feasible.

PROOF. [of Corollary 7] That (ii) and (iii) are equivalent is a statement that results directly from the equivalence of (10) and (11) in Theorem 4. To see this, let $\theta \in \mathbb{R}^n$ be a solution to problem (8) and set $I^c = \mathbb{I}^-(\theta) \cup \mathbb{I}^+(\theta)$, $I = \mathbb{I}^0(\theta)$, $h_{I^c} \in \{-1, +1\}^{|I^c|}$ such that $h_i = +1$ if $i \in \mathbb{I}^+(\theta)$ and $h_i = -1$ if $i \in \mathbb{I}^-(\theta)$. Then Eq. (10) can be written as

$$\begin{aligned} \left|\eta^{\top} X_{I^{c}} h_{I^{c}}\right| &\leq \left\|X_{I}^{\top} \eta\right\|_{1} \quad \forall \eta \in \mathbb{R}^{n} \\ \Leftrightarrow \max_{\eta \in \mathbb{R}^{n}} \left\{ \left|\eta^{\top} X_{I^{c}} h_{I^{c}}\right| + \left\|X_{I^{c}}^{\top} \eta\right\|_{1} \text{s. t. } \left\|X^{\top} \eta\right\|_{1} = 1 \right\} &\leq 1. \end{aligned}$$

$$(20)$$

Similarly Eq. (11) reads as

$$\min_{\alpha_I} \left\{ \left\| \alpha_I \right\|_{\infty} \text{s.t.} X_{I^c} h_{I^c} = X_I \alpha_I \right\} \le 1.$$
 (21)

The equivalence (ii) \Leftrightarrow (iii) then follows by applying the chains of maximums $\max_{(I,I^c):|I|=d} \max_{h_{I^c} \in \{\pm 1\}^{|I^c|}}$ to each of the equations (20) and (21) and noting that $\max_{h_{I^c} \in \{\pm 1\}^{|I^c|}} |\eta^\top X_{I^c} h_{I^c}| = \|X_{I^c}^\top \eta\|_1.$

We shall now establish the equivalence (i) \Leftrightarrow (ii). Let $\theta \in \mathbb{R}^n$ and $\boldsymbol{y} \in \mathbb{R}^N$ be any vectors such that $I = |\mathbb{I}^0(\theta)| = d$. The so-defined I can be any subset of \mathbb{I} provided |I| = d. Hence any θ satisfying this cardinality constraint solves problem (8) if and only if (20) holds for any partition (I, I^c) of \mathbb{I} with |I| = d and for any $h_{I^c} \in \{-1, +1\}^{|I^c|}$. This is equivalent to Eq. (18).

Finally, let us observe that

$$\max_{(I,I^c):|I|=d} \max_{\eta \in \mathbb{R}^n} \left\{ \left\| X_{I^c}^\top \eta \right\|_1 \text{ s. t. } \left\| X^\top \eta \right\|_1 = 1 \right\}$$

is a decreasing function of d so that if (18) holds for some d_0 , it holds also for any $d \ge d_0$. It follows that (i) \Leftrightarrow (ii), hence completing the proof. \Box

It should be mentioned that the equivalence (i) \Leftrightarrow (ii) was also obtained in earlier papers, see e.g., [42,43]. Uniqueness of the solution follow in a similar way as in the proof of Corollary 7 by invoking conditions S1' and S2' of Theorem 4.

Corollary 8 (Uniqueness) Let d be an integer. Then the following statements are equivalent. (i')

$$\forall \, \theta \in \mathbb{R}^n, \forall \boldsymbol{y} \in \mathbb{R}^N, \, |\mathbb{I}^0(\theta)| \ge d \\ \Rightarrow \operatorname*{arg\,min}_{w \in \mathbb{R}^n} \|\phi(w)\|_1 = \left\{\theta\right\}$$
(22)

(ii') Eq. (18) holds with strict inequality.(iii')

$$\max_{\substack{(I,I^c):\\|I|=d}} \max_{h\in\{\pm 1\}^{|I^c|}} \min_{\alpha} \left\{ \|\alpha\|_{\infty} \text{ s. t.} \right.$$

$$X_{I^c}h = X_I\alpha,$$

$$\exists S \subset I, \operatorname{rank}(X_S) = n, \|\alpha_S\|_{\infty} < 1 \right\} \le 1$$
(23)

Remark 9 Corollaries 7 and 8 imply the following. If there exists an integer d such that (18) or (19) holds and $\{\theta \in \mathbb{R}^n : \|\phi(\theta)\|_0 \leq N-d\} \neq \emptyset$, then $\arg\min_{\theta} \|\phi(\theta)\|_0 \subset \arg\min_{\theta} \|\phi(\theta)\|_1$. It follows under these conditions that whenever θ^o solves the ℓ_0 problem (4), it also solves the ℓ_1 problem (8). In particular $\arg\min_{\theta} \|\phi(\theta)\|_1 = \{\theta^o\} \Rightarrow \arg\min_{\theta} \|\phi(\theta)\|_0 = \{\theta^o\}.$

It should be noted that when the data are noise-free, there always exists a d such that (17)-(19) hold. For example d = N is the maximum possible value that satisfies these conditions. Let us denote by $\pi^o(X)$ the minimum integer d such that the conditions (17)-(19) hold, that is,

$$\pi^{o}(X) = \min \left\{ d \in \mathbb{I} \text{ s.t. Eq. (18) is true} \right\}.$$
(24)

Such a number $\pi^{o}(X)$ depends only on the matrix X. It can be viewed as a measure of the richness properties of the regressor matrix X. Recoverability of the true parameter vector θ^{o} by the least ℓ_{1} -norm estimator (8) in the face of gross errors is enhanced when $\pi^{o}(X)$ is small. We may hence say that the smaller $\pi^{o}(X)$, the richer (or more generic) the regressors in X are.

Computing $\pi^{o}(X)$ directly from the definition (24) is a hard combinatorial problem with a complexity comparable to that of the ℓ_0 problem (4). An algorithm of slightly reduced complexity but still combinatorial has been derived in [37] for this purpose. Here, we ask the question of whether $\pi^{o}(X)$ can be more cheaply estimated in a somewhat efficient way. Such estimates are most likely over-estimates and lead to sufficient conditions for exact recoverability of the parameter vector θ^{o} in the presence of gross errors sequence $\{f_t\}$.

2.4 Sufficient conditions of recoverability by convex optimization

We start by introducing the following notations :

$$v_1(k) = \max_{(I,I^c):|I|=k \ge \nu_n(X)} \left\| X_I^\top (X_I X_I^\top)^{-1} X_{I^c} \right\|_{\infty}$$
(25)

$$v_2(k) = \max_{(I,I^c):|I|=k} \left\| X_{I^c}^{\top} (XX^{\top})^{-1} X \right\|_1,$$
(26)

where the maximum is taken over the set of those partitions (I, I^c) of \mathbb{I} that satisfy |I| = k. In addition, let

$$k_1(X) = \min_{k \in \mathbb{I}, k \ge \nu_n(X)} \left\{ k : v_1(k) \le 1 \right\}$$
(27)

and

$$k_2(X) = \min_{k \in \mathbb{I}} \{k : v_2(k) \le 1/2\}.$$
 (28)

Assuming that rank(X) = n, it can be seen that the numbers $k_i(X)$, i = 1, 2, are well-defined. First, note that $\nu_n(X) \leq N$ so that the condition $k_i(X) \geq \nu_n(X)$ is achievable. Moreover, by considering the trivial partition (I, I^c) with $I = \mathbb{I}$ and $I^c = \emptyset$, we see that a possible (the largest indeed) value for $k_i(X)$ is N.

Theorem 10 (Sufficient condition for exact recovery) Assume $\operatorname{rank}(X) = n$. Then the numbers $k_1(X)$ and $k_2(X)$ satisfy

$$\forall \theta \in \mathbb{R}^{n}, \forall \boldsymbol{y} \in \mathbb{R}^{N} \left| \mathbb{I}^{0}(\theta) \right| \geq \min(k_{1}(X), k_{2}(X)) \\ \Rightarrow \theta \in \operatorname*{arg\,min}_{w \in \mathbb{R}^{n}} \left\| \phi(w) \right\|_{1}$$
(29)

where $\phi(w) = \boldsymbol{y} - X^{\top} w$. If in addition all the inequalities involved in the definition of $k_1(X)$ and $k_2(X)$ are strict, then the second part of (29) becomes $\arg\min_{w\in\mathbb{R}^n} \|\phi(w)\|_1 = \{\theta\}$, that is, θ is the unique minimizer of (8).

PROOF. To prove the first statement, we just need to show that

$$\min(k_1(X), k_2(X)) \ge \pi^o(X).$$
 (30)

Part 1: $k_1(X) \ge \pi^o(X)$. Define

$$v_{0}(k) = \max_{\substack{(I,I^{c}): \\ |I|=k}} \max_{\substack{\alpha \in \mathbb{R}^{|I|} \\ \alpha \in \mathbb{R}^{|I|}}} \min_{\alpha \in \mathbb{R}^{|I|}} \left\{ \|\alpha\|_{\infty} \text{ s. t. } X_{I^{c}}h = X_{I}\alpha \right\}$$

that is, $v_0(k)$ corresponds to the left hand side of (19) (with *d* replaced by *k*). By making use of Corollary 7 and the definitions (25) and (27), it is enough to show that $v_0(k) \leq v_1(k)$. For this purpose, let (I, I^c) be an arbitrary partition of \mathbb{I} such that $|I| \geq \nu_n(X)$. Consider the problem

$$\min_{\alpha \in \mathbb{R}^{|I|}} \left\{ \left\| \alpha \right\|_{\infty} : X_{I^c} h = X_I \alpha \right\},\tag{31}$$

where $h \in \{-1, +1\}^{|I^c|}$ but otherwise arbitrary. Let $p^*(h)$ be the optimal value of problem (31) and pose

$$\alpha^*(h) = \operatorname*{arg\,min}_{\alpha \in \mathbb{R}^{|I|}} \big\{ \|\alpha\|_2 \text{ s.t. } X_{I^c}h = X_I\alpha \big\}.$$

Since $\alpha^*(h)$ is a feasible point for problem (31), it must hold that $p^*(h) \leq ||\alpha^*(h)||_{\infty}$. The so-defined $\alpha^*(h)$ is the well-known least Euclidean-norm solution to an underdetermined system of linear equations [4]; $\alpha^*(h)$ can be analytically expressed as $\alpha^*(h) = X_I^{\top}(X_I X_I^{\top})^{-1} X_{I^c} h$ for all $h \in \{-1, +1\}^{|I^c|}$. As a consequence,

$$\max_{h \in \{\pm 1\}^{|I^c|}} p^*(h) \leq \max_{h \in \{\pm 1\}^{|I^c|}} \left\| X_I^\top (X_I X_I^\top)^{-1} X_{I^c} h \right\|_{\infty} \\
\leq \max_{h \in \{\pm 1\}^{|I^c|}} \left\| X_I^\top (X_I X_I^\top)^{-1} X_{I^c} \right\|_{\infty} \left\| h \right\|_{\infty} \\
= \left\| X_I^\top (X_I X_I^\top)^{-1} X_{I^c} \right\|_{\infty}.$$

The last equality uses $||h||_{\infty} = 1$. It follows that if $v_0(k) \leq v_1(k)$ hence proving that $k_1(X) \geq \pi^o(X)$. Part 2: $k_2(X) \geq \pi^o(X)$

Proceeding from Corollary 7 and the definitions (26) and (28), we just need to show that

$$\max_{(I,I^c):|I|=k} \max_{\eta \in \mathbb{R}^n} \left\{ \left\| X_{I^c}^\top \eta \right\|_1 \text{ s. t. } \left\| X^\top \eta \right\|_1 = 1 \right\} \le v_2(k).$$

To this end, set $b = X^{\top} \eta$. Then $b \in im(X^{\top})$ and $\eta = (XX^{\top})^{-1}Xb$. It follows that

$$\begin{split} \max_{\eta \in \mathbb{R}^n} & \left\{ \left\| X_{I^c}^\top \eta \right\|_1 \text{s.t.} \left\| X^\top \eta \right\|_1 = 1 \right\} \\ &= \max_{b \in \text{im}(X^\top)} \left\{ \left\| X_{I^c}^\top (XX^\top)^{-1} Xb \right\|_1 \text{s.t.} \left\| b \right\|_1 = 1 \right\} \\ &\leq \left\| X_{I^c}^\top (XX^\top)^{-1} X \right\|_1. \end{split}$$

Taking now the maximum over all partitions (I, I^c) of \mathbb{I} , |I| = d, the result follows.

Uniqueness. This is a straightforward consequence of Corollary 8. $\ \square$

Evaluating numerically $k_1(X)$ and $k_2(X)$ is still a combinatorial problem. Next we investigate some overestimates of $\pi^o(X)$ which are free from the maximization over sets (I, I^c) . The new thresholds have the important advantage of being more easily computable. **Theorem 11 (Another sufficient condition)** Assume that $\nu_n(X) \leq N - 1$ and define the following numbers

$$r(X) = \max_{k \in \mathbb{I}} \left| x_k (XX^{\top})^{-1} x_k \right|$$
(32)

$$\xi(X) = \max_{k \in \mathbb{I}} \min_{\gamma_k \in \mathbb{R}^{N-1}} \left\{ \left\| \gamma_k \right\|_{\infty} \text{s.t.} x_k = X_{\neq k} \gamma_k \right\}$$
(33)

where $X_{\neq k} \triangleq X_{\mathbb{I} \setminus \{k\}}$ is the matrix obtained from X by removing its k-th column. Then the following statement is true: $\forall p \in \left\{\frac{1}{r(X)}, 1 + \frac{1}{\xi(X)}\right\}$,

$$\forall \boldsymbol{\theta} \in \mathbb{R}^{n}, \forall \boldsymbol{y} \in \mathbb{R}^{N}, \ \left| \mathbb{I}^{0}(\boldsymbol{\theta}) \right| > N - \frac{p}{2} \\ \Rightarrow \operatorname*{arg\,min}_{w \in \mathbb{R}^{n}} \left\| \boldsymbol{\phi}(w) \right\|_{1} = \left\{ \boldsymbol{\theta} \right\}.$$
 (34)

PROOF. The proof is decomposed into two cases.

Case 1: p = 1/r(X). From Theorem 10, it is known that $\|X_{I^c}^{\top}(XX^{\top})^{-1}X\|_1 < 1/2$, $I^c = \mathbb{I} \setminus \mathbb{I}^0(\theta)$, is a sufficient condition for θ to be the unique minimizer of (8). Now we use the fact that the 1-norm of a matrix is the maximum of the 1-norms of its columns:

$$\begin{split} \|X_{I^c}^{\top}(XX^{\top})^{-1}X\|_1 &= \max_{t=1,\dots,N} \|X_{I^c}^{\top}(XX^{\top})^{-1}x_t\|_1 \\ &= \max_{t=1,\dots,N} \sum_{k \in I^c} |x_k^{\top}(XX^{\top})^{-1}x_t| \\ &\leq |I^c| \, r(X). \end{split}$$

Therefore a sufficient condition for θ to be the unique solution of (8) is that $|I^c|r(X) < 1/2$. The conclusion follows immediately.

Case 2: $p = 1 + 1/\xi(X)$.

Since $\nu_n(X) \leq N-1$, each $x_k, k \in \mathbb{I}$, can be written as a linear combination of the columns of $X_{\neq k}$. Let $\gamma_k \in \mathbb{R}^{N-1}$ be any vector satisfying $x_k = X_{\neq k}\gamma_k$. It follows that for any $\eta \in \mathbb{R}^n$,

$$\left|x_{k}^{\top}\eta\right| \leq \sum_{t \neq k} \left|\gamma_{k,t}\right| \left|x_{t}^{\top}\eta\right| \leq \left\|\gamma_{k}\right\|_{\infty} \left(\left\|X^{\top}\eta\right\|_{1} - \left|x_{k}^{\top}\eta\right|\right)$$

with $\gamma_{k,t}$ denoting the entry of $\gamma_k \in \mathbb{R}^{N-1}$ indexed by t. Since this holds for any γ_k such that $x_k = X_{\neq k} \gamma_k$, it holds also for

$$\gamma_k^{\star} = \operatorname*{arg\,min}_{\gamma \in \mathbb{R}^{N-1}} \left\{ \left\| \gamma \right\|_{\infty} \mathrm{s.\,t.\,} x_k = X_{\neq k} \gamma \right\}$$

Hence,

$$\left|x_{k}^{\top}\eta\right| \leq \xi(X)\left(\left\|X^{\top}\eta\right\|_{1} - \left|x_{k}^{\top}\eta\right|\right) \quad \forall k \in \mathbb{I}, \forall \eta \in \mathbb{R}^{n}$$

or, equivalently,

$$x_k^{\top} \eta \Big| \le \frac{\xi(X)}{1+\xi(X)} \| X^{\top} \eta \|_1 \quad \forall k \in \mathbb{I}, \, \forall \eta \in \mathbb{R}^n.$$

Summing over the set I^c yields

$$\max_{\eta \neq 0} \frac{\left\| X_{I^c}^{\top} \eta \right\|_1}{\left\| X^{\top} \eta \right\|_1} \le \frac{\xi(X)}{1 + \xi(X)} \left| I^c \right|$$
(35)

In virtue of (18), it appears that for θ to be the unique minimizer of (8), it is sufficient that $\frac{\xi(X)}{1+\xi(X)} |I^c| < 1/2$ from which we see that $|I^c| < 1/2(1+1/\xi(X))$ is a sufficient condition. \Box

It should be noted that the numbers r(X) and $\xi(X)$ defined in (32) and (33) are both computable from matrix X. r(X) is less expensive to evaluate numerically than $\xi(X)$ but leads in general to a more pessimistic bound than $\xi(X)$ on the number of tolerable outliers. Computing $\xi(X)$ literally from the definition (33), for example by interior point methods, requires solving about N linear programs having each a worst-case complexity bounded by $O(\sqrt{N}\ln(1/\epsilon))$ where ϵ refers to the precision demanded [17]. Empirical evidence tend to suggest that the bound obtained from $\xi(X)$ on the number of correctable outliers is very close to $N - \pi^{o}(X)$ (see Section 5.4). As it turns out, while the computational complexity (polynomial) of $\xi(X)$ is lower than that of the algorithm developed in [37] for estimating directly $\pi^{o}(X)$, it still provides a competitive bound.

Remark 12 $\xi(X)$ can be approximated at a cheaper computational cost by replacing the infinity norm with the 2-norm. This provides an over-estimate $\widehat{\xi}(X) \ge \xi(X)$ defined by

$$\widehat{\xi}(X) = \max_{k \in \mathbb{I}} \min_{\gamma \in \mathbb{R}^{N-1}} \left\{ \|\gamma\|_2 \text{ s.t. } x_k = X_{\neq k} \gamma \right\}$$
$$= \max_{k \in \mathbb{I}} \left\| X_{\neq k}^\top (X_{\neq k} X_{\neq k}^\top)^{-1} x_k \right\|_2.$$

We conclude this section with a few technical remarks concerning some interesting properties of the numbers r(X) and $\xi(X)$.

Lemma 13 (Some properties of r(X)) Under the assumption that $\nu_n(X) \leq N-1$, r(X) and $\xi(X)$ satisfy:

$$\max\left\{\frac{1}{r(X)}, 1 + \frac{1}{\xi(X)}\right\} \le N - \nu_n(X) + 1, \quad (36)$$

$$N - \frac{1}{2r(X)} \ge k_2(X) \ge \pi^o(X).$$
(37)

PROOF. Proof of (36):

First case: $1/r(X) \leq N - \nu_n(X) + 1$. We know from the proof of Theorem 11 (see also Part 2 in the proof of Theorem 10) that

$$r(X) \geq \frac{1}{|I|} \left\| X_{I}^{\top} (XX^{\top})^{-1} X \right\|_{1} \geq \frac{1}{|I|} \max_{\eta \neq 0} \frac{\left\| X_{I}^{\top} \eta \right\|_{1}}{\left\| X^{\top} \eta \right\|_{1}}$$

for any $I \subset \mathbb{I}$. A special case is when the subset I is a singleton of the form $I = \{q\}$. For any $\eta \in \mathbb{R}^n$, let $T(\eta) = \sup (X^\top \eta) = \{t \in \mathbb{I} : x_t^\top \eta \neq 0\}$. When $\eta \neq 0$, consider an index $q(\eta)$ such that $q(\eta) \in \arg \max_{k \in T(\eta)} |x_k^\top \eta|$. Then by applying the above inequality with $I = \{q(\eta)\}$, we get

$$r(X) \geq \max_{\eta \neq 0} \frac{|x_{q(\eta)}^{\top}\eta|}{\left\|X_{T(\eta)}^{\top}\eta\right\|_{1}} \geq \max_{\eta \neq 0} \frac{1}{\left|T(\eta)\right|}$$

with $|T(\eta)|$ standing for the cardinality of $T(\eta)$. When $\eta \neq 0$, the smallest value $|T(\eta)|$ can take is $N - \nu_n(X) + 1$ where $\nu_n(X)$ is the number defined by Eq. (6). It can therefore be concluded that $r(X) \geq 1/(N - \nu_n(X) + 1)$. Second case: $1 + 1/\xi(X) \leq N - \nu_n(X) + 1$. The second case follows by a similar reasoning as in the first one. In effect, according to [6], the following equality holds,

$$\min_{\gamma_k \in \mathbb{R}^{N-1}} \left\{ \|\gamma_k\|_{\infty} \text{ s. t. } x_k = X_{\neq k} \gamma_k \right\}$$

$$= \max_{\eta \in \mathbb{R}^n} \left\{ x_k^\top \eta \text{ s. t. } \|X_{\neq k}^\top \eta\|_1 = 1 \right\}$$

$$= \max_{\eta \neq 0} \left\{ \frac{|x_k^\top \eta|}{\|X_{\neq k}^\top \eta\|_1} \text{ s. t. } x_k^\top \eta \ge 0 \right\}.$$

For a given $\eta \neq 0$, pick $q_1(\eta)$ such that

$$q_1(\eta) \in \underset{k \in \mathbb{I}}{\operatorname{arg\,max}} \left\{ x_k^\top \eta : x_k^\top \eta \ge 0 \right\}.$$

By exploiting the equalities above and using the notation $T(\eta)$ defined earlier we get that

$$\xi(X) \ge \max_{\eta \neq 0} \frac{|x_{q_1(\eta)}^\top \eta|}{\|X_{\neq k}^\top \eta\|_1}.$$

Now the conclusion can be reached by arguing similarly as in the *first case*.

Proof of (37): Let (I, I^c) be a partition of \mathbb{I} and set k = N - 1/(2r(X)). First note that

$$|I| \ge k \implies |I^c| r(X) \le 1/2.$$

On the other hand, we know (from the proof of Theorem 11) that

$$\|X_{I^c}^{\top}(XX^{\top})^{-1}X\|_1 \le |I^c|r(X).$$

It follows that $|I| \geq k \Rightarrow ||X_{I^c}^{\top} (XX^{\top})^{-1} X||_1 \leq 1/2$ and hence $v_2(k) \leq 1/2$. By invoking the definition of the number $k_2(X)$ in (28), it can be concluded that $k \geq k_2(X)$. \Box

Remark 14 For any nonsingular matrix $T \in \mathbb{R}^{n \times n}$, $r(TX) = r(X), \xi(TX) = \xi(X), k_i(TX) = k_i(X), i = 1, 2$. It follows that the numbers $r(X), \xi(X)$ and $k_i(X)$, i = 1, 2, depend only on the subspace spanned by the rows of the regressor matrix X.

3 Some implementation aspects

3.1 Enforcing recoverability by iterative re-weighting

The parameter vector θ^o from the model (1) can be uniquely recovered by solving the convex problem (8) if θ^o satisfies, for example, condition (34) of Theorem 11. In case this condition is not naturally satisfied, an interesting question is how we can process the data in order to promote it. In this section we discuss an algorithmic strategy for enhancing the recoverability of θ^o by ℓ_1 minimization. Our discussion is inspired by [9]. The idea is to solve a sequence of problems of the type (8) with different weights computed iteratively [9,1]. The iterative scheme can be defined for a fixed number r_{\max} of iterations as follows. At iteration $r = 0, \ldots, r_{\max}$, compute

$$\theta^{(r)} = \operatorname*{arg\,min}_{\theta \in \mathbb{R}^n} \sum_{t=1}^N w_t^{(r)} \left| y_t - \theta^\top x_t \right|, \qquad (38)$$

with weights defined, for all t, by $w_t^{(0)} = 1/N$, and

$$w_t^{(r)} = \frac{\xi_t^{(r)}}{\sum_{t=1}^N \xi_t^{(r)}}, \quad \text{if } r \ge 1,$$

where

$$\xi_t^{(r)} = \frac{1}{\left|y_t - x_t^\top \theta^{(r-1)}\right| + \delta},$$

with $\delta > 0$ a small number whose role is to prevent division by zero and r is the iteration number. Note that there are many other reweighting strategies which can be used in (38), see *e.g.*, [12,44,22]. Since we are dealing here with a sequence of convex optimization problems, they can be numerically implemented using any convex solver. In particular the CVX software package [18] solves efficiently this category of problems in a Matlab environment.

3.2 On the treatment of the noise $\{e_t\}$

The formulations (4) and (8) are convenient when the noise $\{e_t\}$ is equal to zero. Nevertheless, they are expected to work in the presence of a moderate amount of

noise. To take explicitly the noise $\{e_t\}$ into account, we propose to compute estimates $\hat{e} \in \mathbb{R}^N$ and $\varphi \in \mathbb{R}^N$ (of e and f respectively) by minimizing a cost of the form $\|\hat{e}\|_2^2 + \lambda \|\varphi\|_0$ under an equality constraint of the form (1). In other words, we consider the problem

$$\underset{(\theta,\varphi)\in\mathbb{R}^{n}\times\mathbb{R}^{N}}{\text{minimize}} \left[\frac{1}{2}\left\|\boldsymbol{y}-\boldsymbol{X}^{\top}\boldsymbol{\theta}-\boldsymbol{\varphi}\right\|_{2}^{2}+\lambda\left\|\boldsymbol{\varphi}\right\|_{0}\right]$$
(39)

and its convex relaxation,

$$\underset{(\theta,\varphi)\in\mathbb{R}^n\times\mathbb{R}^N}{\text{minimize}} \left[\frac{1}{2} \|\boldsymbol{y} - \boldsymbol{X}^{\top}\boldsymbol{\theta} - \varphi\|_2^2 + \lambda \|\varphi\|_1\right].$$
(40)

where $\lambda \geq 0$ is a regularization parameter.

Lemma 15 A pair $(\theta^*, \varphi^*) \in \mathbb{R}^n \times \mathbb{R}^N$ solves (40) if and only if it satisfies

$$XX^{\top}\theta^{\star} - X(\boldsymbol{y} - \varphi^{\star}) = 0$$
(41)

$$X^{\top}\theta^{\star} - (\boldsymbol{y} - \varphi^{\star}) = -\lambda s(\varphi^{\star}), \qquad (42)$$

where $s(\varphi^{\star})$ is a vector in \mathbb{R}^N whose entries $s_t(\varphi^{\star})$, $t = 1, \ldots, N$, are defined by: $s_t(\varphi^{\star}) = \operatorname{sign}(\varphi_t^{\star})$ if $\varphi_t^{\star} \neq 0$ and $s_t(\varphi^{\star}) \in [-1, 1]$ if $\varphi_t^{\star} = 0$.

PROOF. Let $l(\theta, \varphi) = \frac{1}{2} \| \boldsymbol{y} - \boldsymbol{X}^{\top} \theta - \varphi \|_{2}^{2} + \lambda \| \varphi \|_{1}$ be the objective function of the problem (40). Then lis a proper convex function which is differentiable with respect to variable θ on \mathbb{R}^{n} and admits a subdifferential at any variable $\varphi \in \mathbb{R}^{N}$. $(\theta^{*}, \varphi^{*})$ minimizes $l(\theta, \varphi)$ iff $0 = \nabla_{\theta} l(\theta^{*}, \varphi^{*})$ and $0 \in \partial_{\varphi} l(\theta^{*}, \varphi^{*})$. These conditions translate immediately into $XX^{\top}\theta^{*} - X(\boldsymbol{y} - \varphi^{*}) = 0$ and $-(\boldsymbol{y} - \varphi^{*} - X^{\top}\theta^{*}) + \lambda s(\varphi^{*}) = 0$, where $s(\varphi^{*}) \in$ $\partial \| \varphi^{*} \|_{1}$ is a subgradient of $\| \varphi \|_{1}$ at φ^{*} . \Box

It is interesting to note that (41)-(42) imply $Xs(\varphi^*) = 0$, which is very similar to (9). The following lemma characterizes the uniqueness of the solution of (40).

Lemma 16 (Uniqueness of solution to (40)) A

pair (θ^*, φ^*) is the unique solution to problem (40) if and only if both of the following statements are true

(i) $(\theta^{\star}, \varphi^{\star})$ satisfies conditions (41)-(42) for some $s(\varphi^{\star}) \in \partial \|\varphi^{\star}\|_{1}$

(*ii*) rank
$$(X) = n$$
 and rank $(\Psi_{\mathcal{S}^c}) = |\mathcal{S}^c|$.

Here, $\Psi = I_N - X^{\top} (X X^{\top})^{-1} X$, with I_N being the identity matrix of order N, Ψ_{S^c} is a matrix formed with the columns of Ψ indexed by S^c defined by $S^c = \mathbb{I} \setminus S$, with $S = \{t \in \mathbb{I} : |s_t(\varphi^*)| < 1\}.$

The expression of $(\theta^{\star}, \varphi^{\star})$ is then given by:

$$\theta^{\star} = (XX^{\top})^{-1}X(\boldsymbol{y} - \boldsymbol{\varphi}^{\star}), \qquad (43)$$

If
$$|\mathcal{S}^{c}| = 0$$
, then $\varphi^{\star} = 0$, otherwise
 $\varphi^{\star}_{\mathcal{S}^{c}} = \left(\Psi^{\top}_{\mathcal{S}^{c}}\Psi_{\mathcal{S}^{c}}\right)^{-1}\Psi^{\top}_{\mathcal{S}^{c}}\left(\Psi\boldsymbol{y} - \lambda s(\varphi^{\star})\right), \quad \varphi^{\star}_{\mathcal{S}} = 0.$
(44)

PROOF. $l(\theta, \varphi)$ is a quadratic function of θ . For a fixed φ^* , the minimizer θ^* of $l(\theta, \varphi^*)$ is unique if and only if X has full row rank, i.e., rank(X) = n. The unique value of θ^* is expressed in function of φ^* by (43). Plugging the expression (43) of θ^* in the objective gives

$$\tilde{l}(\varphi) \triangleq l(\theta^{\star}, \varphi) = \frac{1}{2} \left\| \Psi \boldsymbol{y} - \Psi \varphi \right\|_{2}^{2} + \lambda \left\| \varphi \right\|_{1}.$$

The rest of the proof then boils down to showing that the minimizer φ^* of $\tilde{l}(\varphi)$ is unique if and only if rank $(\Psi_{\mathcal{S}^c}) = |\mathcal{S}^c|$. To begin with, let us point out the following (see also ⁴ [39]). If φ^* and ξ^* are two minimizers of $\tilde{l}(\varphi)$, then we have necessarily

$$\Psi\varphi^{\star} = \Psi\xi^{\star} \tag{45}$$

$$s(\varphi^{\star}) = s(\xi^{\star}). \tag{46}$$

The relation (45) follows from the strict convexity of $l(\varphi)$ as a function of $\Psi\varphi$. In effect, by changing the optimization variable into $\delta = \Psi\varphi$, $\tilde{l}(\varphi)$ becomes $\frac{1}{2} ||\Psi \boldsymbol{y} - \delta||_2^2 + \lambda ||\Psi^{\dagger}\delta + v||_1$, with v a vector in ker(Ψ) and \dagger referring to generalized inverse. This last function is strictly convex with respect to δ . As a consequence, its minimizer is unique and equal to $\delta^* = \Psi\varphi^*$. To see why the relation (46) holds, plug the expression (43) of θ^* into (42). We get $\lambda s(\varphi^*) = \Psi \boldsymbol{y} - \Psi \varphi^*$. Combining this with (45) (i.e., the uniqueness of $\Psi\varphi^*$) yields immediately (46).

Let us examine first the case where $|S^c| = 0$. This is indeed equivalent to $S = \mathbb{I}$ and so, $\varphi^* = 0$. Would there exist another minimizer ξ^* of $\tilde{l}(\varphi)$, it should obey (46), which implies that ξ^* is necessarily equal to zero. Now consider the case $|S^c| > 0$.

Sufficiency. Assume that rank(Ψ_{S^c}) = $|S^c|$. As argued above, any two minimizers φ^* and ξ^* of $\tilde{l}(\varphi)$ obey (45)-(46). From (46) we get that $S \subset \{t \in \mathbb{I} : \xi_t^* = 0\}$, which implies that $S^c \supset \operatorname{supp}(\xi^*)$. As a consequence, we can write (45) in the following reduced form $\Psi_{S^c}(\varphi_{S^c}^* - \xi_{S^c}^*) = 0$. With rank(Ψ_{S^c}) = $|S^c|$, this implies that $\varphi^* = \xi^*$ and that the minimizer of $\tilde{l}(\varphi)$ is unique.

Necessity. Assume that rank($\Psi_{\mathcal{S}^c}$) $< |\mathcal{S}^c|$. Consider a nonzero vector $\eta \in \mathbb{R}^N$ such that $\eta_{\mathcal{S}} = 0$ and $\eta_{\mathcal{S}^c} \in \ker(\Psi_{\mathcal{S}^c})$. Let $\eta_1 = \nu \eta$, with $\nu \neq 0$. It is straightforward to verify that $\Psi \varphi^* = \Psi(\varphi^* + \eta_1)$. Note that ν can be chosen sufficiently small such that φ^*_t and $\varphi^*_t + \eta_{1,t}$ have the same sign whenever $\varphi^*_t \neq 0$. Following a similar path as in the proof of Theorem 4, we can establish that $s(\varphi^*) = s(\varphi^* + \eta_1)$. Finally, with $\Psi \varphi^* = \Psi(\varphi^* + \eta_1)$, $s(\varphi^*) = s(\varphi^* + \eta_1)$ and the fact that φ^* is an optimal solution (hence satisfying (42)), it is easy to check that $\varphi^* + \eta_1$ also satisfies (42). By Lemma 15, $\varphi^* + \eta_1 (\neq \varphi^*)$ solves (40). Hence, the solution is not unique. Derivation of Eqs (43)-(44). These relations result from simple rearrangements of (41)-(42). \Box

From Lemma 16, it appears that the true vector \boldsymbol{f} can be found by problem (40) if and only if there is a vector $\hat{\theta} \in \mathbb{R}^n$ such that $(\hat{\theta}, \boldsymbol{f})$ satisfies the conditions (i)-(ii) of Lemma 16. In particular, $(\hat{\theta}, \boldsymbol{f})$ must satisfy (42). A necessary condition for this is that $\Psi \boldsymbol{e} = \lambda \boldsymbol{s}(\boldsymbol{f})$. And this implies that the regularization parameter must verify $\lambda \geq ||\Psi \boldsymbol{e}||_{\infty}$ when $\boldsymbol{f} = 0$, and $\lambda = ||\Psi \boldsymbol{e}||_{\infty}$ when $\boldsymbol{f} \neq 0$. Note further that if $\boldsymbol{e} = 0$ and $\boldsymbol{f} \neq 0$, then λ must be equal to zero! However, if λ is set to zero in (40), then the solution set is

$$\{(\theta, \varphi) : \theta = (XX^{\top})^{-1}X(\boldsymbol{y} - \varphi), \varphi \in \boldsymbol{y} + \operatorname{im}(X^{\top})\}.$$

Since this set contains infinitely many elements, we conclude that it is unlikely to get exactly the true f by solving (40) irrespective of the value of the regularization parameter λ .

In any case, the estimation error can be bounded as follows.

Proposition 17 Assume that the conditions of Lemma 16 are satisfied and denote with (θ^*, φ^*) the solution to problem (40). Then

$$\|\theta^{\star} - \theta^{o}\|_{2} \leq (K_{1}\varepsilon + \lambda K_{2}) + K_{1}M\sqrt{\frac{|\mathcal{J} \cap \mathbb{I}^{c}(\theta^{o})|}{|\mathcal{J}|}} \quad (47)$$

where $\varepsilon = \max_{t \in \mathbb{I}} |e_t|, M = \max_{t \in \mathbb{I}} |f_t|,$

$$K_{1} = \max_{|J| \ge \nu_{n}(X)} \sqrt{|J|} \left\| (XX^{\top})^{-1} (I_{n} + E_{J} + 2E_{J}^{2} + E_{J}^{3}) X_{J} \right\|_{2}$$

$$K_{2} = \max_{|J| \ge \nu_{n}(X)} \sqrt{|J|} \left\| (X_{J}X_{J}^{\top})^{-1} X_{J} \right\|_{2}$$
(48)
(49)

with $E_J = (X_{J^c} X_{J^c}^{\top}) (X_J X_J^{\top})^{-1}$, $J \subset \mathbb{I}$. In (47) I_n is the identity matrix of order n; the set \mathcal{J} denotes the maximizing argument of (48) and $\mathbb{I}^c(\theta^o) = \mathbb{I} \setminus \mathbb{I}^0(\theta^o)$.

PROOF. The idea of the proof consists in deriving first an expression of $\theta^* - \theta^o$ and then working out a bound on its norm. From (43) and the data model (1), we have

$$\theta^{\star} = (XX^{\top})^{-1}X \left(X^{\top} \theta^{o} + \boldsymbol{e} + \boldsymbol{f} - \varphi^{\star} \right)$$

This, by noting that $\varphi_{\mathcal{S}}^{\star} = 0$, can be written as

$$(XX^{\top})(\theta^{\star} - \theta^{o}) = X_{\mathcal{S}}(\boldsymbol{e}_{\mathcal{S}} + \boldsymbol{f}_{\mathcal{S}}) + X_{\mathcal{S}^{c}}(\boldsymbol{e}_{\mathcal{S}^{c}} + \boldsymbol{f}_{\mathcal{S}^{c}} - \varphi_{\mathcal{S}^{c}}^{\star})$$

Using formula (44) and manipulating a little, we arrive at

$$(XX^{\top})(\theta^{\star} - \theta^{o}) = \begin{bmatrix} X_{\mathcal{S}} - X_{\mathcal{S}^{c}}(\Psi_{\mathcal{S}^{c}}^{\top}\Psi_{\mathcal{S}^{c}})^{-1}\Psi_{\mathcal{S}^{c}}^{\top}\Psi_{\mathcal{S}} \end{bmatrix} (\boldsymbol{e}_{\mathcal{S}} + \boldsymbol{f}_{\mathcal{S}}) + \lambda X_{\mathcal{S}^{c}}(\Psi_{\mathcal{S}^{c}}^{\top}\Psi_{\mathcal{S}^{c}})^{-1}\Psi_{\mathcal{S}^{c}}^{\top}s(\varphi^{\star}).$$

 $^{^4\,}$ It is to be noted that the analysis in [39] provides only a sufficient condition.

Further calculations using the Woodbury's matrix identity and exploiting the relation $Xs(\varphi^*) = 0$, yield

$$(XX^{\top}) (\theta^{\star} - \theta^{o}) = (I_n + E_{\mathcal{S}} + 2E_{\mathcal{S}}^2 + E_{\mathcal{S}}^3) X_{\mathcal{S}} (\boldsymbol{e}_{\mathcal{S}} + \boldsymbol{f}_{\mathcal{S}}) - \lambda (I + E_{\mathcal{S}}) X_{\mathcal{S}} s(\varphi_{\mathcal{S}}^*)$$

with $E_{\mathcal{S}} = (X_{\mathcal{S}^c} X_{\mathcal{S}^c}^{\top}) (X_{\mathcal{S}} X_{\mathcal{S}}^{\top})^{-1}$. The result follows by multiplying with $(XX^{\top})^{-1}$, remarking that $(XX^{\top})^{-1} = (X_{\mathcal{S}} X_{\mathcal{S}}^{\top})^{-1} (I_n + E_{\mathcal{S}})^{-1}$ and taking the euclidean norm.

It is interesting to notice that the numbers K_1 , K_2 and $|\mathcal{J}|$ depend solely on the data matrix X. Moreover, when the sequence $\{f_t\}$ contains only a few nonzero elements (but otherwise arbitrarily large), the last term in (47) is likely to vanish. As a consequence, even though the bound M can be large in principle, the bound on the estimation error can be kept at a reasonable level.

4 Extension to multivariable systems

We consider now the multivariable analogue of model (1) written in the form

$$y_t = A^o x_t + f_t + e_t, (50)$$

where $y_t \in \mathbb{R}^m$ is the output vector at time $t, \{f_t\} \subset \mathbb{R}^m$ is the sequence of errors, $\{e_t\} \subset \mathbb{R}^m$ is the noise sequence, $A^o \in \mathbb{R}^{m \times n}$ is the parameter matrix.

The question of interest is how to recover the matrix A^o from measurements corrupted by a vector sequence of sparse errors $\{f_t\}$. The sparse optimization approach is still applicable to this case, that is, we can formulate the estimation problem as

$$\min_{A \in \mathbb{R}^{m \times n}} \left| \left\{ t : y_t - Ax_t \neq 0 \right\} \right| \tag{51}$$

with $|\cdot|$ standing for cardinality. It can be easily verified that Theorem 3 applies to (51) as well.

The convex relaxation takes the form of a nonsmooth optimization with a cost functional consisting of a sumof-norms of errors [29,11],

$$\underset{A \in \mathbb{R}^{m \times n}}{\operatorname{minimize}} \sum_{t=1}^{N} \|y_t - Ax_t\|_2$$
(52)

with $\|\cdot\|_2$ referring to the Euclidean norm.

Theorem 18 A matrix $A^* \in \mathbb{R}^{m \times n}$ solves the sum-ofnorms problem (52), if and only if any of the following equivalent statements holds:

T1. There exists a sequence of vectors $\{\beta_t\}_{t\in\mathbb{I}^0(A^\star)} \subset \mathcal{B}_2(0,1)$ such that

$$\sum_{t \notin \mathbb{I}^0(A^\star)} v_t^\star x_t^\top + \sum_{t \in \mathbb{I}^0(A^\star)} \beta_t x_t^\top = 0, \qquad (53)$$

where $v_t^* = (y_t - A^* x_t) / \|y_t - A^* x_t\|_2$. Here, $\mathcal{B}_2(0,1) \subset \mathbb{R}^m$ is the Euclidean unit ball of \mathbb{R}^m . T2. For any matrix $\Lambda \in \mathbb{R}^{m \times n}$,

$$\Big|\sum_{t\notin\mathbb{I}^{0}(A^{\star})}v_{t}^{\star\top}\Lambda x_{t}\Big| \leq \sum_{t\in\mathbb{I}^{0}(A^{\star})}\left\|\Lambda x_{t}\right\|_{2}.$$
 (54)

T3. The optimal value of the problem

$$\min_{Z \in \mathbb{R}^{m \times p}} \|Z\|_{2,\infty} \text{ subject to } V^* X_{\mathbb{I}^c(A^*)}^\top = Z X_{\mathbb{I}^0(A^*)}^\top$$
(55)

 $p = |\mathbb{I}^{0}(A^{\star})|$ and V^{\star} being a matrix formed with the unit 2-norm vectors v_{t}^{\star} , for $t \in \mathbb{I} \setminus \mathbb{I}^{0}(A^{\star})$, is smaller than 1.

Moreover, the solution A^* is unique if and only if any of the following assertions is true:

- T1'. (53) holds and rank $(X_{\mathcal{T}}) = n$ where $\mathcal{T} = \{t \in \mathbb{I}^0(A^\star) : \|\beta_t\|_2 < 1\}.$
- T2'. (54) holds with strict inequality symbol for all $\Lambda \in \mathbb{R}^{m \times n}$, $\Lambda \neq 0$.

PROOF. The proof is similar to that of Theorem 4. It is therefore omitted here.

It is interesting to note that based on Theorem 18, the analysis carried out in the previous sections can be easily generalized to the multivariable case. In particular, Proposition 6 and Theorems 10-11 can be restated for the multivariable model (50) with only some slight modifications. For illustration purpose, we just state below the multivariable counterpart of Corollary 7.

Corollary 19 Let d be an integer. Then the following three statements are equivalent.

(j)

$$\forall A \in \mathbb{R}^{m \times n}, \forall Y \in \mathbb{R}^{m \times N}, \left| \mathbb{I}^{0} \left(A \right) \right| \geq d \\ \Rightarrow A \in \operatorname*{arg\,min}_{W \in \mathbb{R}^{m \times n}} \left\| Y - WX \right\|_{2, \operatorname{col}}$$
(56)

(jj)

$$\max_{\substack{(I,I^c):\\|I|=d}} \max_{\Lambda \in \mathbb{R}^{m \times n}} \left\{ \|\Lambda X_{I^c}\|_{2,\operatorname{col}} \operatorname{s.t.} \|\Lambda X\|_{2,\operatorname{col}} = 1 \right\} \le 1/2$$
(57)

(jjj)

$$\max_{\substack{(I,I^c): V \in \mathbb{B}^{m \times |I^c|} \ Z \in \mathbb{R}^{m \times |I|}}} \min_{\substack{\{\|Z\|_{2,\infty} \text{ s. t.} \\ |I| = d \\ X_{I^c} V^\top = X_I Z^\top \}} \leq 1$$
(58)

with
$$\mathbb{B}^{m \times q} = \left\{ \begin{bmatrix} b_1 & \cdots & b_q \end{bmatrix} \in \mathbb{R}^{m \times q}, \ b_i \in \mathcal{B}_2(0,1) \right\}$$

PROOF. The proof is similar to that of Corollary 7.

(jj) \Leftrightarrow (jjj) : We exploit the equivalence between (54) and (55). First by letting $I = \mathbb{I}^0(A^*)$, $I^c = \mathbb{I}^c(A^*)$, V_{I^c} be a matrix collecting all the vectors $v_t^* \in \mathbb{B}^m$, $t \in I^c$, (54) can equivalently be written as

$$\max_{\Lambda \in \mathbb{R}^{m \times n}} \left[\left| \operatorname{tr} \left(V_{I^c}^\top \Lambda X_{I^c} \right) \right| + \left\| \Lambda X_{I^c} \right\|_{2, \operatorname{col}} \operatorname{s.t.} \left\| \Lambda X \right\|_{2, \operatorname{col}} = 1 \right] \\ \leq 1$$

Maximizing over all the sets (I, I^c) satisfying |I| = d and over all $V \in \mathbb{B}^{m \times |I^c|}$ yields (57) after remarking that $\max_{V_{I^c} \in \mathbb{B}^{m \times |I^c|}} |\operatorname{tr} (V_{I^c}^\top \Lambda X_{I^c})| = ||\Lambda X_{I^c}||_{2,\operatorname{col}}$. Proceeding similarly from (55), yields (58). Hence (jj) \Leftrightarrow (jj). (j) \Leftrightarrow (jj) : By Theorem 18 and the first part of the proof, any matrix A with $|\mathbb{I}^0(A)| = d$ minimizes the objective $||Y - WX||_{2,\operatorname{col}}$ (with variable W) if and only if (57) holds. The conclusion is obtained by observing that

$$\max_{\substack{(I,I^c): \Lambda \in \mathbb{R}^{m \times n} \\ |I| = d}} \max_{\substack{\{ \|\Lambda X_{I^c}\|_{2, \text{col}} \text{ s. t. } \|\Lambda X\|_{2, \text{col}} = 1 \}}$$

is decreasing as a function of d. \Box

An analogue of Corollary 8 can be obtained similarly. It is interesting to note that the statement (34) of Theorem 11 holds unchanged in the multivariable case with $p = \xi(X)$.

Remark 20 (Geometric median) In the special case where n = 1, the matrix A^o in (50) reduces to a vector $a^o \in \mathbb{R}^m$. Assuming $x_t = 1$ for all t, the problem (52) then becomes

$$\underset{a \in \mathbb{R}^m}{\operatorname{minimize}} \sum_{t=1}^N \|y(t) - a\|_2.$$
(59)

This is the so-called geometric median problem. By applying (54), we can see that a° solves (59) if $|\mathbb{I}^{0}(a^{\circ})|/N \geq 1/2$.

5 Numerical illustration

5.1 Static models subject to intermittent gross errors

In our first experiment we consider static linear and affine models of the form (1) with n = 4 and N = 500. The affine model refers to the case where the regressor x_t has the form $x_t = [\tilde{x}_t^\top \ 1]^\top$. The goal is to estimate the probability of exact recovery of the true parameter vector by problem (8) in function of the number of nonzero elements in the sequence $\{f_t\}$. For this purpose, the noise $\{e_t\}$ is set to zero. The nonzero elements of $\{f_t\}$ are drawn from a Gaussian distribution with mean 100 and variance 1000^2 . For each level of sparsity (*i.e.*, proportion of nonzeros), a Monte Carlo simulation of size 100 is carried out with randomly generated static/affine models and 500 data samples at each run. Repeating this for four situations (linear/affine and linear/affine with positive f_t 's), we obtain the results depicted in Figure 1.

We observe that in the linear case, problem (8) provides the true parameter vector when the output is affected by up to 80% of nonzero gross errors. This is because the data $\{x_t\}$ which were sampled from a Gaussian distribution are very generic. In the case of affine models, the performance is a little less good. If we set all f_t 's to have the same sign, then as suggested by condition (15), the percentage of outliers that can be corrected by the optimization problem (8) cannot exceed 50%.

5.2 Static models with both noise and gross errors

Consider now the case of static models of the form (1)in the presence of both $\{e_t\}$ and $\{f_t\}$ sampled from Gaussian distributions $\mathcal{N}(0, \sigma_e^2)$ and $\mathcal{N}(100, 1000^2)$ respectively. The variance σ_e^2 is selected so as to achieve a certain signal to noise ratio before the gross error sequence is added to the output. Again, by carrying out a Monte-Carlo simulation of size 100 with different sparsity levels and randomly generated models at each run, we obtain the average errors plotted in Figure 2. It turns out that the results returned by problems (8) and (40) with $\lambda = 0.10$ are almost the same for an SNR in {10 dB, 20 dB}. The performance can be assessed by comparing with an "oracle" estimate *i.e.*, the least squares estimate one would obtain if the locations of zeros in the sequence $\{f_t\}$ were known. The results in Figure 2 tend to suggest that the proposed approach performs very well. For the current numerical experiment, our results are very close to the ideal estimate when the proportion of nonzeros is less than 70%.

5.3 Dynamic linear models subject to sensor intermittent faults

In the case when (1) represents a dynamic ARX model subject to gross errors, it can be observed (see Fig. 3) that the probabilities of exact recovery are much smaller than in the static case studied in Section 5.1. This difference is related to the richness (or genericity) of the regression vectors (columns of X) involved in each case. In the static example above, the vectors $\{x_t\}$ are freely sampled in any direction of \mathbb{R}^n by following a Gaussian distribution. In the dynamic system case however, the data vectors $\{x_t\}$ constructed as in (3) are constrained to lie on a manifold. As a result, the data matrix X generated by the dynamic system is less generic. According to conditions of the paper, and (34) in particular, there is a threshold depending on the richness of the data such that exact recovery is guaranteed whenever the number of zero entries in \boldsymbol{f} is larger than this threshold. So, the more generic the data contained in X are, the more outliers can be removed by problem (8). Note that the lack of sufficient genericity can be compensated (to some extent) by implementing the iterative sparsity enhancing technique (ℓ_1 reweighted algorithm) described in Section 3.1. This leads, for only two iterations, to significantly improved results as represented in Figure 4.

5.4 Numerical evaluation of sufficient bounds

This subsection presents a numerical evaluation of the estimates of number of outliers that can be corrected



Figure 1. Estimates of probabilities of exact recovery when noise $\{e_t\}$ is equal to zero. From a numerical point of view, the recovery is said to be exact if $\|\hat{\theta} - \theta^o\|_2 \leq 10^{-5}$ and inexact otherwise, with $\hat{\theta}$ being the estimated parameter vector.



Figure 2. Average relative estimation error versus sparsity level.



Figure 3. Estimates of probabilities of exact recovery when noise $\{e_t\}$ is equal to zero. Results of a Monte-Carlo simulation of size 100 with randomly generated linear ARX systems of order $n_a = n_b = 2$.



Figure 4. Estimates of probabilities of exact recovery by reweighed ℓ_1 minimization when noise $\{e_t\}$ is equal to zero. Results of a Monte-Carlo simulation of size 100 with randomly generated linear ARX systems with orders $n_a = n_b = 2$.

by the nonsmooth optimization-based estimator. Note that the numbers $k_i(X)$, i = 1, 2 from Theorem 10 are hard to compute numerically because this would require a combinatorial optimization. To be more specific, the complexity of evaluating literally $k_i(X)$, i = 1, 2, is about $\sum_{k=1}^{k_i(X)} {N \choose k} C_i(N, n, k)$ where ${N \choose k}$ refers to the binomial coefficient, $C_1(N, n, k) = O(n^3 + kn(N + 2n - k))$ and $C_2(N, n, k) = O((n^2 + nN)(N + n - k))$ denote the complexity induced by the computation of $||X_I^{\top}(X_I X_I^{\top})^{-1} X_{I^c}||_{\infty}$ and $||X_{I^c}^{\top}(X X^{\top})^{-1} X||_1$ respectively with I and I^c being some sets such that $|I| = k \leq N$ and $|I^c| = N - k$.

Therefore we just compare those bounds which are easier to compute. More specifically, four thresholds are compared: PSfrag replacements

- The bounds 1/(2r(X)) and $1/2 + 1/(2\xi(X))$ obtained in Theorem 11.
- The mutual coherence-based bound $1/2(1 + 1/\mu(P_X))$ with $P_X = I_N - X^{\top} (XX^{\top})^{-1} X$ obtained in [5,1]. Here μ represents the so-called mutual coherence.
- The bound T(X) [37] which is used as a reference since it corresponds indeed to a direct computation of $N - \pi^o(X) + 1$ (assuming the inequality in (18) is replaced with a strict one), see Eq. (24). Recall that computing such a bound has a combinatorial complexity in the dimensions (n, N) of the matrix X. Therefore, to make it feasible at a reasonable time on a standard computer, we have to set n = 2 and $N \leq 200$.

Figure 5 compares the sufficient thresholds in the case of static data drawn from a Gaussian distribution $\mathcal{N}(0, I_3)$. Figure 6 compares the same thresholds for dynamic data in the form (3). The generating system in this case is an ARX model defined by $y_t = -0.40y_{t-1} - 0.15u_{t-1}$ and driven by a normally distributed input sequence. In all cases, the data matrix X is normalized so as to have unit $\|\cdot\|_{\Sigma_X}$ -norm columns before being processed. Here, the norm $\|\cdot\|_{\Sigma_X}$ is defined by $\|x\|_{\Sigma_X} = (x^{\top}\Sigma_X^{-1}x)^{1/2}$ with $\Sigma_X = XX^{\top}$. The plots in Figure 5 and Figure 6 draw the average values obtained over 100 independent runs in term of percentage with respect to the total number of data. The results suggest three interesting facts :

- All the bounds are very loose that is, they largely underestimate the number of admissible gross errors. For example Figure 1 shows that exact recovery can be achieved in the face a relatively large proportion (more than 70%) of corrupted data while the sufficient bounds in Figure 5 indicate a value around 20. This is normal since the bounds reflect worst-case distributions of the outliers and their signs (see Theorem 4 and Corollary 7).
- The bound based on $\xi(X)$ approaches the bound T(X)[37] while still enjoying less numerical complexity. The other bounds based respectively on mutual-coherence and r(X) are overall very close. These two last bounds seem to be more sensitive to the richness of the data and probably to their magnitudes also. This fact is more apparent when the data are not normalized.
- As could be intuitively expected, the dynamic data generated by a linear system are less generic. The bounds obtained in this case are smaller. The question as to which type of dynamic system can generate more generic data is open.



Figure 5. Comparison of sufficient bounds on the number of gross errors for exact recovery : static data drawn from a Gaussian distribution.

Comparison of execution times. Evaluating the *r*-based and μ -based bounds defined above is clearly very cheap as compared to the two other bounds. Therefore we shall only compare the execution times for the bounds T(X) [37] and $1/2(1+1/\xi(X))$ (see Theorem 11) for $X \in \mathbb{R}^{n \times N}$. This is done by measuring the average time over 10 runs.⁵ The results reported in Table 1 show that for a

⁵ The computation is performed in a Matlab environment (version 2013a,64-bit), on a computer equipped with a processor Intel(R) Core(TM) i7-3630QM CPU@2.4Ghz, RAM 16Go. Only 10 runs have been considered because the computation time for T(X) grows very quickly beyond the capacity of the computer. Indeed T(X) is computed only once when n = 5 because the algorithm takes too long to complete (about 2 hours for each run in this case). Note that



Figure 6. Comparison of sufficient bounds on the number of gross errors for exact recovery : dynamic data generated by an ARX model.

given number N of data points, the computation time for the algorithm in [37] is small for $n \leq 3$ but grows very fast (at a combinatorial rate) when n increases. In contrast, the cost associated with the evaluation of the ξ -based bound grows only at a logarithmic rate. This numerical experiment confirms that the proposed ξ -based bound is algorithmically cheaper to compute than T(X). For example, for $X \in \mathbb{R}^{5 \times 200}$ (see last column of Table 1), computing T(X) takes nearly 2 hours while the ξ -based bound derived in the current paper is obtained in less than 27 seconds.

n	2	3	4	5
T(X) [sec.]	0.02	1.40	101.82	5.62×10^3
ξ -based [sec.]	22.70	24.56	25.44	26.80

Table 1

Empirical comparison of the numerical complexities associated with evaluating T(X) [37] and the ξ -based bound (see Theorem 11) in term of execution times. Here the number N of data points is fixed and equal to 200.

6 Conclusion

In this paper we have discussed the potential of nonsmooth convex optimization for addressing the problem of robust estimation. Considering in particular the problem of inferring an unknown parameter vector from measurements which are subject to possibly large gross errors, we have shown that an exact recovery is possible regardless of the number of gross errors provided certain conditions of genericity hold. Then we investigated worst-case conditions which depend solely on the number of gross errors affecting the data. Necessary and sufficient conditions have been derived in this case. Since such conditions are numerically expensive to test directly, we have relaxed them into some sufficient but relatively tight conditions for exact recovery. Simulations results reveal that the proposed worst-case conditions for exact recovery are somewhat pessimistic when compared to the potential of the nonsmooth estimator in practice. Concerning the identification problem, future work will consider the problem of designing the excitation of a dynamic system so as to achieve such strong genericity properties on the regressor matrix.

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what matters in this experiment is not the numerical values of the execution times but the trend they exhibit.

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