ESTIMATING MINIMUM EFFECT WITH OUTLIER SELECTION

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We introduce one-sided versions of Huber’s contamination model, in which corrupted samples tend to take larger values than uncorrupted ones. Two intertwined problems are addressed: estimation of the mean of the uncorrupted samples (minimum effect) and selection of the corrupted samples (outliers). Regarding estimation of the minimum effect, we derive the minimax risks and introduce estimators that are adaptive with respect to the unknown number of contaminations. The optimal convergence rates differ from the ones in the classical Huber contamination model. This fact uncovers the effect of the one-sided structural assumption of the contaminations. As for the problem of selecting the outliers, we formulate the problem in a multiple testing framework for which the location and scaling of the null hypotheses are unknown. We rigorously prove that estimating the null hypothesis while maintaining a theoretical guarantee on the amount of the falsely selected outliers is possible, both through false discovery rate (FDR) and through post hoc bounds. As a by-product, we address a long-standing open issue on FDR control under equi-correlation, which reinforces the interest of removing dependency in such a setting.

1. Introduction. We are considering statistical problems for which some of the available data have been corrupted. Such issues have been addressed by different fields in statistics, depending on how one defines and considers the corruption. Two examples are robust estimation and sparse modeling. In the former, Huber’s contamination model [42, 43] is the prototypical setting for handling this problem. It assumes that among $n$ observations $Y_1, \ldots, Y_n$, most of them follow some normal distribution $N(\theta, \sigma^2)$ whereas the corrupted data are arbitrarily distributed. In sparse modeling, one typically assumes that the data $Y_1, \ldots, Y_n$ are normally distributed with mean $\gamma_i$, where $\gamma_i = \theta$ for uncorrupted samples and $\gamma_i \neq \theta$ is arbitrary for
corrupted samples (see [8] for a related model).

However, in some practical problems, corrupted samples do not take arbitrary values and instead satisfy a structural assumption. Consider for instance the situation where the $Y_i$’s are concentration measurements of a pollutant, coming from $n$ sensors spread out at $n$ locations of a city. This pollutant has a background concentration value $\theta$ in the city, but, due to local pollution effects, the sensors may record larger values at some locations. Health authorities are then interested in evaluating the degree of background pollution and in finding the most affected regions in the city.

In this work, we introduce one-sided contamination models to account for the structural assumption that corrupted samples tend to take larger values than uncorrupted ones. Then, we consider the twin problems of estimating the distribution of the uncorrupted samples and of identifying the corrupted samples.

1.1. Models and objectives.

1.1.1. One-sided Contamination Model (OSC). We first introduce a one-sided counterpart of Huber’s contamination model for which some of the $Y_i$’s follow a $\mathcal{N}(\theta, \sigma^2)$ distribution, whereas the remaining samples are positively contaminated, that is, they have a distribution that stochastically dominates a $\mathcal{N}(\theta, \sigma^2)$ distribution, but is otherwise arbitrary.

More formally, we assume that

$$Y_i = \theta + \sigma \varepsilon_i, \quad 1 \leq i \leq n,$$

where $\sigma > 0$ is a standard deviation parameter (either equal to 1 or unknown), $\theta \in \mathbb{R}$ is a fixed minimum effect and the $\varepsilon_i$’s are independent noise random variables. Denoting by $\pi_i$ the unknown distribution of the noise, we assume that, for some $k$, the distribution $\pi = \otimes_{i=1}^{n} \pi_i$ of $\varepsilon$ belongs to the set

$$\mathcal{M}_k = \left\{ \pi = \otimes_{i=1}^{n} \pi_i : \pi_i \succeq \mathcal{N}(0, 1), 1 \leq i \leq n, \sum_{i=1}^{n} \mathbb{1}_{\{\pi_i \succ \mathcal{N}(0, 1)\}} \leq k \right\},$$

where $\succeq$ (resp. $\succ$) denotes the stochastic domination (resp. strict stochastic domination), see Section 1.4 below for a definition. In $\mathcal{M}_k$, at most $k$ of the distributions $\pi_i$ are allowed to strictly dominate the Gaussian measure. Model (1) satisfies the heuristic explanation described above: if $\pi \in \mathcal{M}_k$, then at least $n - k$ samples are non-contaminated and follow a $\mathcal{N}(\theta, \sigma^2)$ distribution, whereas the remaining contaminated samples stochastically dominate this distribution.
In this model, henceforth referred to as the One-Sided Contamination (OSC) model, the parameter $\theta$ corresponds to the expectation of the non-contaminated samples. If $k \leq n - 1$, it also satisfies

$$
\theta = \min_{1 \leq i \leq n} E(Y_i),
$$

and can therefore be interpreted as a minimum theoretical effect. In particular, $\theta$ is identifiable for $k \in [n/2, n - 1]$, whereas this is not the case in the classical Huber’s model.

Throughout the paper, probabilities (resp. expectations) in model (1) are denoted by $P_{\theta,\pi,\sigma}$ (resp. $E_{\theta,\pi,\sigma}$). The parameter $\sigma$ is dropped in the notation whenever $\sigma = 1$.

1.1.2. One-sided Gaussian Contamination Model ($g$OSC). As in the sparse Gaussian vector model, we also consider a specific case of the OSC model for which the contaminated samples are still assumed to be normally distributed, that is, the $\pi_i$’s are Gaussian distributions with unit variance and positive mean $\mu_i/\sigma$ where $\mu \in \mathbb{R}^n_+$ is a contamination effect. In that case, the model can be rewritten as

$$
Y_i = \theta + \mu_i + \sigma \xi_i, \quad 1 \leq i \leq n,
$$

where the $\xi_i$’s are i.i.d. $\mathcal{N}(0, 1)$ distributed and $\mu \in \mathbb{R}^n_+$ is unknown. Defining the mean vector

$$
\gamma = (\theta + \mu_i)_{1 \leq i \leq n},
$$

we deduce that $Y$ follows a normal distribution with unknown mean $\gamma$ and variance $\sigma^2 I_n$, while $\theta$ corresponds to $\min_{1 \leq i \leq n} \{\gamma_i\}$, that is, the minimum component of the mean vector.

To formalize the connection with the OSC model, let $\varepsilon_i = \mu_i/\sigma + \xi_i$ and $\pi_i = \mathcal{N}(\mu_i/\sigma, 1)$ for all $i \in \{1, \ldots, n\}$. Then, (4) is a particular case of (1) because $\mathcal{N}(\mu_i/\sigma, 1) \succeq \mathcal{N}(0, 1)$. As in the OSC model, we prescribe the number of contaminated samples to be less or equal to $k$, by defining

$$
\mathcal{M}_k = \left\{ \mu \in \mathbb{R}^n_+ : \sum_{i=1}^{n} 1_{\{\mu_i \neq 0\}} \leq k \right\}.
$$

In what follows, we refer to model (4) as the One-Sided Gaussian Contamination ($g$OSC) model. The probabilities (resp. expectations) in this model are denoted by $P_{\theta,\mu,\sigma}$ (resp. $E_{\theta,\mu,\sigma}$). Whenever we assume that the variance parameter $\sigma$ is known and is equal to 1, the subscript $\sigma$ is dropped in the above notation.
1.1.3. Objectives. We are interested in the two following intertwined problems:

- **Objective 1: optimal estimation of the minimum effect.** We aim at establishing the minimax estimation rates of $\theta$, both in the OSC (1) and in the gOSC (4) models. In particular, we explore the role of the one-sided contamination assumption for the computation of such estimation rates. As explained below, this problem is at the crossroads of several lines of research, including robust estimation and non-smooth linear functional estimation.

- **Objective 2: controlled selection of the outliers.** Here, we are interested in finding the contaminated samples. In the Gaussian case (gOSC), this is equivalent to selecting the positive entries of $\mu$ in (5). Adopting a multiple testing framework, we aim at building a selection procedure with suitable false discovery rate (FDR) control [4] and providing a valid post hoc bound [37, 38]. The difficulty stems from the fact the minimum effect $\theta$ is unknown. In contrast to objective 1 where the contaminated samples were considered as nuisance quantities, for this second objective the contaminated samples are now interpreted as the signal whereas $\theta$ is a nuisance parameter.

Objective 2 is intrinsically connected to the problem of removing the correlation when making (one-sided) multiple tests from Gaussian equi-correlated test statistics: when the equi-correlation is carried by the latent factor $\theta$, we can remove this correlation by subtracting an estimator of $\theta$ from the test statistics. Although this simple strategy is quite common (see, e.g., [34] and references therein), assessing whether the theoretical performance of the resulting procedure remains suitable is a longstanding issue in multiple testing literature. In this work, we provide positive results for this problem, by showing that it is possible to (asymptotically) control the FDR while having (at least) the same power as if the test statistics had been independent.

In the remainder of the introduction, we first describe our contribution for minimum effect estimation and then turn to outlier selection.

1.2. Optimal estimation of the minimum effect.** Given $\sigma^2 = 1$ and a sparsity parameter $k \in \{1, \ldots, n-1\}$, we define the $L_1$ minimax estimation risk of $\theta$ in both gOSC (4) and OSC (1) models:

\[
\mathcal{R}[k,n] = \inf_{\hat{\theta}} \sup_{(\theta,\mu) \in \mathbb{R} \times M_k} \mathbb{E}_{\theta,\mu}[|\hat{\theta} - \theta|]; \quad \mathcal{R}[k,n] = \inf_{\hat{\theta}} \sup_{\theta \in \mathbb{R}, \pi \in M_k} \mathbb{E}_{\theta,\pi}[|\hat{\theta} - \theta|].
\]
We characterize these minimax risks by deriving matching (up to numerical constants) lower and upper bounds, uniformly over the parameter \( k \), see Sections 2 and 3. The results are summarized in Table 1 below. It is interesting to compare these orders of magnitude with those derived for the Huber contamination model with at most \( k \) contaminated samples. From, e.g., Section 2 of [14], the minimax risk is of order \( \max(n^{-1/2}, k/n) \) in Huber’s model (the results in [14] are proved for a model where the number of contaminated samples follows a Binomial distribution with parameters \((n, k/n)\), but the proofs straightforwardly extend to the case of a fixed parameter \( k \)). For \( k \leq \sqrt{n} \), the rate is parametric in all three models. For \( k \in (\sqrt{n}, n/2) \), the assumption of one-sided contamination leads to a \( \log^{1/2}(k^2/n) \) gain over the Huber model, while assuming that the contaminations are Gaussian lead to an additional logarithmic gain. For \( k \in [n/2, n-1] \), recall that Huber’s model is not identifiable while the one-sided contamination model is and we identify various minimax rates. For a fixed proportion \((k/n)\) of contaminated samples, the optimal rate converges to 0 at a polylogarithmic rate. For a proportion \((n-k)/n\) of non-contaminated samples slowly decaying with \( n \), the estimation rate still goes to 0.

<table>
<thead>
<tr>
<th>( \mathcal{R}[k, n] )</th>
<th>General bound</th>
<th>( 1 \leq k \leq \sqrt{n} )</th>
<th>( \sqrt{n} \leq k \leq n/2 )</th>
<th>( n/2 \leq k \leq n-1 )</th>
</tr>
</thead>
<tbody>
<tr>
<td>OSC</td>
<td>( \frac{\log(n^{1/2})}{\log^{1/2}(1+\frac{k}{n})} )</td>
<td>( n^{-1/2} )</td>
<td>( \frac{k/n}{\log^{1/2}(k^2/n)} )</td>
<td>( \frac{\log(n^{1/2})}{\log^{1/2}(k^2/n)} )</td>
</tr>
<tr>
<td>OSC</td>
<td>( \frac{\log^{2}(1+\sqrt{n/k})}{\log^{3/2}(1+\sqrt{n/k^2})} )</td>
<td>( n^{-1/2} )</td>
<td>( \frac{k/n}{\log^{3/2}(k^2/n)} )</td>
<td>( \frac{\log^{2}(1+\sqrt{n/k})}{\log^{3/2}(k^2/n)} )</td>
</tr>
<tr>
<td>Huber</td>
<td>( \max(n^{-1/2}, k/n) )</td>
<td>( n^{-1/2} )</td>
<td>( k/n )</td>
<td>( \infty )</td>
</tr>
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</table>

Table 1

Minimax estimation risks of \( \theta \) (up to numerical constants).

For both models (OSC and gOSC), we also provide estimation procedures that are adaptive to the unknown number \( k \) of contaminated samples. Finally, in Section 4, we consider the case where the noise level \( \sigma \) in (4) is unknown. We prove, in the OSC model, that adaptation to unknown \( \sigma \) is possible, and we characterize the optimal estimation risk for \( \sigma \).

**OSC: Technical aspects and connection to robust estimation.** As explained earlier, the OSC model (1) is a one-sided counterpart of Huber’s contamination.
tion model [42, 43] - see also [59] for the historical reference on the concept of contamination and [54, 51] for more recent reviews. From a technical perspective, minimax bounds for the OSC model proceed from the same general ideas as for Huber’s contamination model, but with a twist. In Huber’s model, the empirical median turns out to be optimal [43], while in the OSC model, there is a benefit to using other empirical quantiles. Since the contaminations are one-sided, the left tail is less perturbed than the right tail. Correcting for the bias and choosing a suitable quantile, we prove that the resulting estimator achieves (up to a constant) the optimal rate $\mathcal{R}[k,n]$. Adaptation to unknown $k$ is performed via Lepski’s method, while adaptation to unknown $\sigma$ is based on a difference of empirical quantiles.

$g$OSC: Technical aspects and connection to non-smooth functional estimation. Pinpointing the minimax risk in the Gaussian contamination model (gOSC) is much more technical. Indeed, standard estimators, such as those based on quantiles, are not optimal in this setting. The key idea of our upper bound is to invert a collection of local tests of the form “$\theta \geq u$” vs “$\theta < u$” for $u \in \mathbb{R}$, following an approach from [13] developed for sparsity testing. Recall that $\gamma_i$ in (5) stands for the expectation of $Y_i$; then under the null “$\theta \geq u$”, we have $\sum_{i=1}^{n} \mathbb{I}_{\gamma_i < u} = 0$, whereas under the alternative “$\theta < u$” we have $\sum_{i=1}^{n} \mathbb{I}_{\gamma_i < u} \geq n - k$. Thus, this approach boils down to estimating the non smooth-functional $u \mapsto \sum_{i=1}^{n} \mathbb{I}_{\gamma_i < u}$.

Starting from the seminal works [44, 24] (for respectively linear and quadratic functionals), an extensive literature on estimating smooth functionals of the mean of a Gaussian vector has been developed. Under a sparsity assumption, this problem has been investigated in [9, 67, 16, 17], and shares some strong connections with the problem of signal detection, as studied in, e.g., [46, 2].

However, estimation of non-smooth functionals (for example $\sum_{i=1}^{n} |\gamma_i|^q$ with $q \in (0,1]$) is significantly more involved, even without sparsity assumptions. See e.g. [49, 8, 10, 48, 56, 41, 69, 47, 13, 50, 15]. For this class of problems, one powerful approach, known as the polynomial approximation method [56, 41], amounts to building a suitable polynomial approximation of the non-smooth function and plugging in unbiased estimators of the moments $\sum_{i=1}^{n} \gamma_i^s$ for some integers $s \in \{1, \ldots, s_{\text{max}}\}$. Unfortunately, we cannot rely on this strategy for estimating $\sum_{i=1}^{n} \mathbb{I}_{\gamma_i < u}$, mainly because the contaminated $\gamma_i$’s may be arbitrarily large. In a related setting, where the contaminated means $\gamma_i \neq \theta$ are distributed according to some smooth prior distributions supported on $\mathbb{R}$, Cai and Jin [8] have pinpointed the optimal rate by using the empirical Fourier transform (see also [22]). However, this approach also falls down in our framework, again because the contaminated
\(\gamma_i\)'s are arbitrary. In this work, we introduce a new strategy that combines polynomial approximation methods with the empirical Laplace transform. Indeed, the empirical Laplace transform has the virtue of being almost insensitive to arbitrarily large values of \(\gamma_i\).

As for the minimax lower bound, we rely on moment matching techniques, following the approach of [56] that has been recently applied to other non-smooth functional models [10, 41, 69, 13].

1.3. Controlled selection of the outliers. Turning to the second objective, we now detail our contributions and discuss the relevant literature. Our approach falls with the multiple testing paradigm and builds upon some of our estimators for \(\theta\) (and for \(\sigma\)).

1.3.1. Multiple testing formulation. Recall that our second objective was to identify the active set of outliers in the general model (1). Again, we emphasize that what we described before as outliers is in this part the quantities of interest (e.g., the city locations with abnormal pollutant concentration in our motivating example). In the OSC model, we formulate this selection problem as \(n\) simultaneous tests of

\[
(8) \quad H_{0,i} : \pi_i = \mathcal{N}(0,1) \text{ against } H_{1,i} : \pi_i \succ \mathcal{N}(0,1), \quad 1 \leq i \leq n.
\]

(Recall that \(\succ\) stands for strict stochastic domination).

In the specific case of the gOSC model (4), this problem amounts to simultaneously testing

\[
H_{0,i} : \mu_i = 0 \text{ against } H_{1,i} : \mu_i > 0, \quad 1 \leq i \leq n.
\]

We denote the set of non-outlier coordinates by \(\mathcal{H}_0(\pi) = \{1 \leq i \leq n : \pi_i = \mathcal{N}(0,1)\}\), and the set of outlier coordinates by \(\mathcal{H}_1(\pi) = \{1 \leq i \leq n : \pi_i \succ \mathcal{N}(0,1)\}\).

The cardinality of \(\mathcal{H}_0(\pi)\) (resp. \(\mathcal{H}_1(\pi)\)) is denoted by \(n_0(\pi)\) (resp. \(n_1(\pi)\)). Hence, \(\pi \in \mathcal{M}_k\), means that the number of outliers is \(n_1(\pi) \leq k\). Our selection problem thus amounts to estimating \(\mathcal{H}_1(\pi)\) (or equivalently \(\mathcal{H}_0(\pi)\)). The dependence on \(\pi\) of \(\mathcal{H}_0(\pi), \mathcal{H}_1(\pi), n_0(\pi), n_1(\pi)\) is sometimes suppressed for convenience.

A multiple-testing procedure is a data-driven set \(R \subset \{1, \ldots, n\}\) of proposed outliers. For a given \(R\), we classically quantify the amount of false positives of \(R\) by its false discovery proportion [4]

\[
(9) \quad \text{FDP}(\pi, R) = \frac{|R \cap \mathcal{H}_0(\pi)|}{|R|} \lor 1,
\]
which records the proportion of errors among the set $R$ of selected outliers. The expectation of this quantity $\mathbb{E}_{\theta,\pi,\sigma}[\text{FDP}(\pi, R)]$ is the false discovery rate, which can be considered as the standard generalization of the single testing type I error rate to large scale multiple testing. The true discovery proportion is then defined by

$$\text{TDP}(\pi, R) = \frac{|R \cap \mathcal{H}_1(\pi)|}{n_1(\pi) \lor 1},$$

and corresponds to the proportion of (correctly) selected outliers among the set of true outliers. The expectation of this quantity $\mathbb{E}_{\theta,\pi,\sigma}[\text{TDP}(\pi, R)]$ is a widely used analogue of the power in single testing, see, e.g., [64, 1, 60]. Our contribution falls into two frameworks:

- **Multiple testing**: find a procedure selecting a subset $R \subset \{1, \ldots, n\}$ as close as possible to $\mathcal{H}_1(\pi)$, i.e. that has a TDP as high as possible while maintaining a controlled FDR;
- **Post hoc bound**: provide a confidence bound on $\text{FDP}(\pi, S)$, uniformly valid over all possible selection $S \subset \{1, \ldots, n\}$.

While the first objective is classical in the multiple testing field, see, e.g., [4, 5, 33, 35], the second objective was proposed more recently in [36, 37, 38]. It is connected to the burgeoning research field of selective inference, see, e.g., [6] and references therein. The rationale behind developing such a bound is that, since the control is uniform, the probability coverage is guaranteed even if $S$ is itself data-dependent. The obtained bound is therefore also valid when practitioners re-use the same dataset, possibly several times, in order to design $S$. We denote the outlier selected set either by $R$ or $S$ depending on the considered setting; $R$ is typically a procedure designed by the statistician, whereas $S$ is chosen by the user.

1.3.2. **Relation to the first objective and to previous literature.** In the OSC model (1), solving the above multiple testing issues is challenging primarily because the parameters $\theta$ and $\sigma$ are unknown. Indeed, this entails that the scaling of the null distribution (i.e. the distribution under the null hypothesis) is unknown. A natural idea is to design a two-stage procedure: first, we estimate $\theta$ and $\sigma$ by some estimators $\hat{\theta}$ and $\hat{\sigma}$ (as we do in the first part of this paper). Then, in a testing stage, we apply a standard multiple testing procedure to the rescaled observation $Y_i' = (Y_i - \hat{\theta})/\hat{\sigma}$.

Estimating the null distribution in a multiple testing context has been popularized in a series of works by Efron, see [26, 28, 29]. Through careful data analyses, Efron noticed that the theoretical null distribution often turns
out to be wrong in practical situations, which can lead to an uncontrolled increase in false positives. To address this issue, Efron recommends estimating the scaling parameters of the null distribution \((\theta, \sigma)\) here) by “central matching”, that is, by fitting a parametric curve to the trimmed data. In his work, Efron provides compelling empirical evidence for his approach. However, to our knowledge, the FDP and TDP of such two-stage testing procedures have never been theoretically controlled. Note that estimating the null in a multiple testing context was also the motivation of the minimax results of \([49, 8]\), although the corresponding multiple testing procedure was not studied. We recall that these previous studies are all developed in the two-sided context, whereas our focus is on a one-sided shape constraint.

Finally, let us mention that a procedure learning part of the null distribution while maintaining FDR control and a form of power optimality has been proposed in \([3, 1]\), often referred to as the Barber-Candès procedure. In the one-sided case, it is able to learn the unknown null density when this is assumed to be symmetric around 0. However, this assumption is not satisfied in our context because the symmetry point of the null distribution \((\theta)\) here) is unknown.

1.3.3. **Summary of our results.** In Sections 5 and S-1, we show that a fair modification of the quantile-based estimators \(\hat{\theta}, \hat{\sigma}\), introduced for the OSC model, can be used to estimate the null distribution to rescale the \(p\)-value process, and can then be suitably combined with classical multiple testing procedures:

1. A new \((\hat{\theta}, \hat{\sigma})\)-rescaled Benjamini-Hochberg procedure \(R\) is defined that satisfies the following FDR controlling property: in the general model (1), for any \(\pi \in \mathcal{M}_k\), with \(k = \lceil 0.9n \rceil\),

\[
\left( \mathbb{E}_{\theta, \pi, \sigma} (\text{FDP}(\pi, R)) - \frac{n_0}{n} \alpha \right)_+ \lesssim \log(n)/n^{1/16}.
\]

In addition, we derive a power result showing that the power (the expectation of the TDP) of this procedure is close to that of the \((\theta, \sigma)\)-rescaled Benjamini-Hochberg procedure, under mild conditions. The latter is an oracle benchmark that would require the exact knowledge of \(\theta\) and \(\sigma\).

2. A new \((\hat{\theta}, \hat{\sigma})\)-rescaled post hoc bound \(\text{FDP}(\cdot)\) is proposed, satisfying, for any \(\pi \in \mathcal{M}_k\), with \(k = \lceil 0.9n \rceil\),

\[
(1 - \alpha - \mathbb{P}_{\theta, \pi, \sigma} (\forall S \subset \{1, \ldots, n\}, \text{FDP}(\pi, S) \leq \text{FDP}(S)))_+ \lesssim \log(n)/n^{1/16}.
\]


To the best of our knowledge, these are the first results that theoretically validate Efron’s principle of empirical null correction in a specific multiple testing context.

For bounding the type I error rates, the technical argument used in our proof is close in spirit to recent studies [58, 45] (among others): the idea is to divide the data into two “orthogonal” parts (small or large $Y_i$’s), the first part being used for the rescaling and the second one for testing. For the power result, our formal argument is entirely new to our knowledge, see also Remark 5.3.

Finally, let us mention that the above rate $\log(n)/n^{1/16}$ is certainly not optimal and could be improved by using more involved estimators $\hat{\theta}, \hat{\sigma}$. Indeed, our concern here regarding the multiple testing procedure is to show that it is consistent with respect to the FDP/TDP metrics, this for a wide range of values of $k$, including the dense case where $k$ is of the order of $n$. Furthermore, despite the fact that this rate is relatively slow, the new rescaled BH procedure improves on the vanilla BH procedure, as illustrated in the numerical experiments of Section S-1.3 (see also Figure 1 below).

1.3.4. Application to decorrelation in multiple testing. It is well known that Efron’s methodology on empirical null correction can be applied to reduce the effect of correlations between the tests, as noted by Efron himself [27, 30] where he mentioned that “there is a lot at stake here”. Several following works supported this assertion, especially by decomposing the covariance matrix of the data into factors, see [34, 55, 32, 31]. However, strong theoretical results on such corrected multiple testing procedures are still not available.

Meanwhile, another branch of the literature aims at incorporating known and unknown dependence into multiple testing procedures, for instance, by resampling-based approaches [68, 62, 63, 61, 25, 3] or by directly plugging the known dependence structure [40, 19, 7]. However, as noted for instance in the discussion of [65], even for very simple correlation structures, no multiple testing procedure has yet been proved to control the FDR while having an optimal expected TDP.

In Section S-1.2, we apply our two-step procedure to address the multiple testing problem in the contaminated one-factor model. For instance, this model encompasses the one-sided Gaussian equi-correlation case (with nonnegative equi-correlation $\rho$) which is often used as a concrete test bed in multiple testing literature, see, e.g., [52, 18, 20] among others. It turns out that the contaminated one-factor model can be written in the form of the OSC model (1) with a random value of $\theta$ and an unknown variance.
$\sigma^2 = 1 - \rho$. Hence, we can directly apply our $(\hat{\theta}, \hat{\sigma})$-rescaled Benjamini-Hochberg procedure introduced above to solve the problem: we show that the new procedure has performances close to the BH procedure under independence (and even with a slight increase of the signal to noise ratio). Even if the model is somewhat specific, this shows that correcting the dependence can be fully theoretically justified. To illustrate numerically the benefit of such an approach, Figure 1 displays a ROC-type curve for four different versions of corrected BH procedure (in the Gaussian equi-correlated case). A full description of the simulation setting and additional experiments are provided in Section S-1.3.

Figure 1. X-axis: targeted FDR level \(\alpha \in \{0.005, 0.05, 0.1, 0.15, 0.2, 0.25, 0.3, 0.35, 0.4, 0.45, 0.5\}\), Y-axis: TDP (power) averaged over 100 replications for four different procedures (see text in Section S-1.3). The model is the one-sided Gaussian one with equi-correlation \(\rho = 0.3\). The parameters used are \(n = 10^6, \Delta = 2.5, k/n = 0.1\).

1.4. Notation. For \(x > 0\), we write \([x]^{\log_2}\) (resp. \([x]^{\log_2}\)) for \(2^{\log_2(x)}\) (resp. \(2^{\log_2(x)^1}\)), that is, the largest (resp. smallest) dyadic number no larger (resp. no smaller) than \(x\). Similarly, \([x]_{e\text{ven}}\) is the largest even integer which is not larger than \(x\).

For \(x \in \mathbb{R}^n\), we denote by \(x_{(k)}\) the \(k\)-th smallest element of \(\{x_i, 1 \leq i \leq n\}\). We also write \(x_{(\ell:m)}\) for the \(\ell\)-th smallest element among \(\{x_i, 1 \leq i \leq m\}\), for some integer \(1 \leq m \leq n\).

In what follows, \(c, c'\) denote numerical positive constants whose values may vary from line to line. For two sequences \((u_t)_{t \in T}\) and \((v_t)_{t \in T}\), we write that for all \(t \in T\), \(u_t \lesssim v_t\) (resp. for all \(t \in T\), \(u_t \gtrsim v_t\)), if there exists a universal constant \(c > 0\) such that for all \(t \in T\), \(u_t \leq cv_t\) (resp. for all \(t \in T\), \(u_t \geq c v_t\)). We write \(u_t \asymp v_t\) if \(u_t \lesssim v_t\) and \(v_t \lesssim u_t\).

For \(X, Y\) two real random variables with respective cumulative distribu-
tion functions $F_X, F_Y$, we write $X \succeq Y$ if for all $x \in \mathbb{R}$, we have $F_X(x) \leq F_Y(x)$. We write $X \succ Y$ if $X \succeq Y$ and if there exists $x \in \mathbb{R}$, such that $F_X(x) < F_Y(x)$. We also denote $P \succeq Q$ (resp. $P \succ Q$) whenever $X \succeq Y$ (resp. $X \succ Y$) for $X \sim P$ and $Y \sim Q$.

We write $\Phi$ for the cumulative distribution function of the standard normal distribution, we write $\bar{\Phi} = 1 - \Phi$ and $\phi$ for its usual density.

For space reasons, part of the outlier selection results, the numerical experiments and all the proofs of the paper are deferred to a supplemental article [11]. For clarity, the sections, equations and results of this supplement are referred to with an additional symbol “S-” in the numbering.

2. Estimation of $\theta$ in the gOSC model with known variance.

In this section, we consider the problem of estimating $\theta$ in the Gaussian contamination model (4) and investigate the $L_1$ minimax risk defined in (7). We assume throughout this section that $\sigma^2 = 1$. The case of unknown variance is discussed in Section 6.

2.1. Lower bound for the gOSC model with known variance.

**Theorem 2.1.** There exists a universal constant $c > 0$ such that for any positive integer $n$ and for any integer $k \in [1, n - 1]$,

$$
R[k, n] \geq c \frac{\log^2 (1 + \left(\frac{k}{n-k}\right)^{1/2})}{\log^{3/2} (1 + \left(\frac{k^2}{n}\right)^{1/3})}.
$$

The proof of this theorem is given in Section S-2.1. The main tool for proving this lower bound is moment matching: we build two priors on the parameter $\gamma$ that correspond to distant values of $\theta$ while they have the same $\log n$ (or so) first moments. This is done in an implicit way using the Hahn-Banach theorem together with properties of Chebychev polynomials, by using techniques close to [50, 13].

Let us distinguish between the following three regimes (see also Table 1):

- for $k \leq \sqrt{n}$, the lower bound (11) is of order $n^{-1/2}$, which is the parametric rate that would hold in the case of no contamination (i.e., $k = 0$);
- for $k \in (\sqrt{n}, \zeta n)$ with $\zeta \in (0, 1)$, the lower bound is of the order $(k/n)\log^{-3/2}(k^2/n)$. In particular, in the non-sparse case $k = \lceil n/2 \rceil$, we obtain a rate $\log^{-3/2} n$;
- for $k \in [\zeta n, n - 1]$ with $\zeta \in (0, 1)$, the lower bound on the minimax risk is of order $\log^2(\frac{n}{n-k}) \log^{-3/2}(n)$. In particular, for $k = n - 1$, the lower bound is of order $\log^{1/2} n$. 

In the remainder of this section, we match these lower bounds by considering three different estimators of \( \theta \), corresponding to the three regimes discussed above. They are then combined to derive an adaptive estimator.

### 2.2. Upper bound for small and large \( k \) in the gOSC model with known variance.

For small and for large values of \( k \) the optimal risk is achieved by simple quantile estimators. For \( k \leq n^{1/2} \), we consider the empirical median defined by

\[
\hat{\theta}_{\text{med}} = Y_{(\lceil n/2 \rceil)} .
\]

The following result holds for \( \hat{\theta}_{\text{med}} \) (note that it is stated in the more general OSC model (1)).

**Proposition 2.2.** Consider the OSC model (1) with \( \sigma = 1 \). Then there exist universal positive constants \( c_1, c_2 \) and a universal positive integer \( n_0 \) such that the following holds. For any \( n \geq n_0 \), any \( k \leq n/10 \), any \( \pi \in \mathcal{M}_k \) and any \( \theta \in \mathbb{R} \), we have

\[
\mathbb{P}_{\theta, \pi} \left( |\hat{\theta}_{\text{med}} - \theta| \geq \frac{3(k+1)}{2(n-k)} + \frac{3\sqrt{(n+1)x}}{n-k} \right) \leq e^{-x} , \quad \text{for all } x \leq c_1 n ,
\]

\[
\mathbb{E}_{\theta, \pi} [ |\hat{\theta}_{\text{med}} - \theta| ] \leq \frac{3(k+1)}{2(n-k)} + c_2 \sqrt{n} .
\]

A proof is provided in Section S-3.1. A consequence is that, for \( k \leq \sqrt{n} \), the empirical median \( \hat{\theta}_{\text{med}} \) achieves the parametric rate \( n^{-1/2} \), which turns out to be optimal in this regime, see Theorem 2.1. Note that in the regime \( k \leq \sqrt{n} \), the empirical median was already known to achieve this parametric rate in the more general Huber’s contamination model (which allows for two-sided contaminations).

When \( k \) is really close to \( n \), there are very few non contaminated data. Since \( \theta = \min_{1 \leq i \leq n} \{ \gamma_i \} \) in the model (4), we consider a debiased empirical minimum estimator

\[
\hat{\theta}_{\text{min}} = Y_{(1)} + \Phi^{-1}(1/n) ,
\]

where we recall that \( \Phi^{-1}(1/n) = \sqrt{2\log(n)} + O(1) \), see Section S-5.2. The following result holds for \( \hat{\theta}_{\text{min}} \) (note that it is again stated in the more general OSC model (1)).
Proposition 2.3. Consider the OSC model (1) with $\sigma = 1$. Then there exists some universal positive integer $n_0$ such that for any $n \geq n_0$, any $\pi \in \mathcal{M}_{n-1}$ and any $\theta \in \mathbb{R}$, the estimator $\hat{\theta}_{\min}$ satisfies

$$
P_{\theta,\pi} [ |\hat{\theta}_{\min} - \theta| \geq 2\sqrt{2 \log n} ] \leq \frac{2}{n} ; \quad \mathbb{E}_{\theta,\pi} [ |\hat{\theta}_{\min} - \theta| ] \leq 2\sqrt{2 \log n} + 1 .
$$

A proof is provided in Section S-3.1. From Theorem 2.1, the estimator $\hat{\theta}_{\min}$ turns out to be optimal when $k$ is very close to $n$, that is, when $k$ is larger than $n - n^\epsilon$ for a fixed $\epsilon \in (0, 1)$ (very few samples are non-contaminated).

2.3. Upper bound in the intermediate regime in the gOSC model with known variance. In the previous section, we have introduced estimators that are optimal in the regimes where $k \leq \sqrt{n}$ and where $k$ is very close to $n$, respectively. The intermediate case turns out to be much more involved.

Let $q \geq 2$ be an even integer whose value will be fixed below. Let $a = 3[1 + \log(3 + 2\sqrt{2})] \approx 8.29$ and $q_{\text{max}} = \lfloor \frac{1}{2a} \log n \rfloor_{\text{even}} - 2$, where $[\cdot]_{\text{even}}$ is defined in Section 1.4. Let us also introduce two rough estimators $\hat{\theta}_{\text{up}}$ and $\hat{\theta}_{\text{low}, q}$ such that $\theta$ is proved to belong to $[\hat{\theta}_{\text{low}, q}, \hat{\theta}_{\text{up}}]$ with high probability. Let $\hat{\theta}_{\text{up}} = Y(1) + 2\sqrt{\log n}$. For any positive and even integer $q$, define $\hat{\theta}_{\text{low}, q} = \hat{\theta}_{\text{med}} - \tau$ with $\tau = \pi^2/(144 q_{\text{max}}^{3/2})$ if $q \leq \frac{3}{10a} \log n$ and $\hat{\theta}_{\text{low}, q} = -\infty$ for larger $q$.

To explain the intuition behind our procedure, assume for the purpose of discussion that we have access to the mean $\gamma_i = \theta + \mu_i$, and that instead of estimating $\theta$ we simply want to test whether $\theta$ is greater than $u$ or not. Thus, our aim is to define a suitable function $\psi_{q, \lambda}(u)$ of $u$ and the $\gamma_i$’s which is close to zero when all $\gamma_i$’s are higher than $u$ and is as large as possible when many $\gamma_i$’s are smaller than $u$. Since at least $n - k$ of the $\gamma_i$’s are equal to $\theta$, having large values of $\psi_{q, \lambda}(u)$ means that $\theta < u$. Assuming without loss of generality that $u = 0$, this can be achieved by constructing an auxiliary function $g_q : \mathbb{R} \mapsto \mathbb{R}$ such that $|g_q(x)| \leq 1$ for $x \in (-\infty, 0]$ and $g_q(x)$ large for $x > 0$. If the interval $(-\infty, 0]$ were replaced by $[-1, 1]$ and the function $g_q$ was restricted to be a polynomial, constructing such a $g_q$ would look like a polynomial extremum problem, which is solved by a Chebychev polynomial (see Section S-5.1 for some definitions and properties). To handle the non-bounded interval $(-\infty, 0]$, we map $(-\infty, 0]$ to $(-1, 1]$ using the function $x \mapsto 2e^x - 1$. Denoting by $T_q$ the Chebychev polynomial of degree $q$, this leads us to considering the function

$$
g_q(x) = T_q(2e^x - 1) = \sum_{j=0}^q a_{j,q} e^{xj}, \quad x \in \mathbb{R} ,
$$

(14)
where the coefficients $a_{j,q}$ are defined in (S-65). It follows from the definition of Chebyshev polynomials that $g_q(x)$ belongs to $[-1, 1]$ for $x \leq 0$ and $g_q(x) = \cosh[q \arccosh(2e^x - 1)]$ for $x > 0$.

Now, for $\lambda > 0$ and $u \in \mathbb{R}$, consider the function $\psi_{q, \lambda}(u)$ defined by

$$
\psi_{q, \lambda}(u) = \frac{1}{n} \sum_{i=1}^{n} g_q(\lambda(u - \gamma_i)) = \frac{1}{n} \sum_{i=1}^{n} g_q(\lambda(u - \theta - \mu_i)).
$$

This function depends on the $\gamma_i$’s. Since all $\mu_i$’s are non negative, it follows from the above observation that $|\psi_{q, \lambda}(u)| \leq 1$ for all $u \leq \theta$. Conversely, for $u \geq \theta$, $\psi_{q, \lambda}(u)$ is lower bounded as follows:

$$
\psi_{q, \lambda}(u) \geq -\frac{k}{n} + \frac{n - k}{n} g_q(\lambda(u - \theta)),
$$

hence it is bounded away from 1 as long as $u - \theta$ is large enough. As a consequence, the smallest number $u_*$ that satisfies $\psi_{q, \lambda}(u_*) > 1$ should be close (in some sense) to $\theta$.

Obviously, we do not have access to the function $\psi_{q, \lambda}$ as it requires the knowledge of the $\gamma_i$’s (or, more precisely, of quantities of the form $e^{-j\lambda \gamma_i}$). Nevertheless, we can still build an unbiased estimator of such quantities using the empirical Laplace transform of $Y$. Given $\lambda > 0$ and $u \in \mathbb{R}$, define

$$
\hat{\eta}_{\lambda}(u) = n^{-1} \sum_{i=1}^{n} e^{\lambda(u-Y_i)-\lambda^2/2}, \quad \eta_{\lambda}(u) = n^{-1} \sum_{i=1}^{n} e^{\lambda(u-\theta-\mu_i)}.
$$

Since all $Y_i$’s are independent and normally distributed with unit variance, we have $E[\hat{\eta}_{\lambda}(u)] = \eta_{\lambda}(u)$. This leads us to considering the statistic

$$
\hat{\psi}_{q, \lambda}(u) = \sum_{j=0}^{q} a_{j,q} \hat{\eta}_{j, \lambda}(u),
$$

which is an unbiased estimator of $\psi_{q, \lambda}(u)$ for any fixed $\lambda > 0$ and $u \in \mathbb{R}$. Since $\hat{\psi}_{q, \lambda}(u)$ approximates $\psi_{q, \lambda}(u)$, we roughly want to take $\hat{\theta}_q$ to be the smallest value such that $\hat{\psi}_{q, \lambda}(u)$ is bounded away from 1. This is why we define $\hat{\theta}_q$ by inverting the function $\hat{\psi}_{q, \lambda}(\cdot)$. More precisely, for an even integer $q \leq q_{\text{max}}$, taking $\lambda_q = \sqrt{2/qa}$, we define the estimator $\hat{\theta}_q$ by

$$
\hat{\theta}_q = \inf \left\{ u \in \left[ \hat{\theta}_{\text{low}, q}, \hat{\theta}_{\text{up}} \right] : \hat{\psi}_{q, \lambda_q}(u) > 1 + \frac{e^{aq}}{\sqrt{n}} \right\},
$$

with the convention $\inf \emptyset = \hat{\theta}_{\text{up}}$. 

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**Theorem 2.4.** Consider the gOSC model (4) with known variance $\sigma^2 = 1$. There exist universal positive constants $c_1, c_2, c_3,$ and $n_0$ such that the following holds for any $n \geq n_0$, any integer $k \in [e^{2a \sqrt{n}} n - 64 n^{1-1/(4a)}]$, any $\mu \in M_k$ and any $\theta \in \mathbb{R}$. The estimator $\hat{\theta}_{q_k}$ defined by (19) with $q_k = \lfloor \frac{1}{a} \log \left( \frac{k}{\sqrt{n}} \right) \rfloor_{\text{even}} \wedge q_{\text{max}}$ satisfies

\begin{equation}
\mathbb{P}_{\theta, \mu} \left( \hat{\theta}_{q_k} \notin \left[ \theta, \theta + c_1 \frac{\log^2 \left( 1 + \sqrt{\frac{k}{n-k}} \right)}{\log^{3/2} \left( \frac{k^2}{n} \right)} \right] \right) \leq c_3 \left( \frac{\sqrt{n}}{k} \right)^{4/3} \log^3 \left( \frac{k^2}{n} \right),
\end{equation}

and

\begin{equation}
\mathbb{E}_{\theta, \mu} \left[ \left| \hat{\theta}_{q_k} - \theta \right| \right] \leq c_2 \frac{\log^2 \left( 1 + \sqrt{\frac{k}{n-k}} \right)}{\log^{3/2} \left( \frac{k^2}{n} \right)}.
\end{equation}

A proof is provided in Section S-3.3. This result shows that $\hat{\theta}_{q_k}$ has a maximum risk of order $\frac{k}{n} \log^{-3/2} \left( \frac{k^2}{n} \right)$ in the regime $k \in [e^{2a \sqrt{n}} n - 64 n^{1-1/(4a)}]$. Combined with the lower bound of Theorem 2.1, we have shown that $\hat{\theta}_{q_k}$ is minimax in the intermediate regime.

**Remark 2.5.** Let us emphasize that in the regime $e^{2a \sqrt{n}} \leq k \leq \lfloor n/2 \rfloor$, the minimax risk is of order $(k/n) \log^{-3/2} (n)$, which is faster than the minimax rate $(k/n) \log^{-1/2} (n)$ that we would obtain in a two-sided deconvolution problem, as in [8] where $k/n \propto n^{-\beta}$ (and by considering the extreme case where there is no regularity assumption, that is, $\alpha = 0$ with their notation).

**Remark 2.6.** If we are only interested in a probability bound (20) and not in the moment bound (21), the preliminary estimators $\hat{\theta}_{\text{low}, q}$ and $\hat{\theta}_{\text{up}}$ are not needed: the estimator could be computed by taking the minimum over $\mathbb{R}$ in (19).

**Remark 2.7.** The behavior of the estimator family $\{ \hat{\theta}_q \}_q$ given by (19) is illustrated in Section S-6.3 by using numerical experiments. While a fine tuning of the constants is required, this reinforces the theoretical finding that $q$ should be chosen as a function of $k$ in order to provide a small risk.

### 2.4. Adaptative estimation in the gOSC model with known variance.

In this section, we combine the three estimators studied in the above section to obtain an estimator that is adaptive with respect to the parameter $k$. The method stems from the Goldenshluger-Lepski principle, see, e.g., [57, 39, 53].
To unify notation, we henceforth write \( \hat{\theta}_0 \) for the median estimator \( \hat{\theta}_{\text{med}} \) and \( \hat{\theta}_{q_{\text{max}}+2} \) for the minimum estimator \( \hat{\theta}_{\text{min}} \). In order to obtain an adaptive procedure, we select one of the estimators \( \{ \hat{\theta}_q, q \in \{0, 2, \ldots, q_{\text{max}}, q_{\text{max}} + 2\} \} \) as follows:

\[
q = \min \{ q \in \{0, \ldots, q_{\text{max}} + 2\} \text{ s.t. } |\hat{\theta}_q - \hat{\theta}_{q'}| \leq \delta_{q'} \text{ for all } q' > q \},
\]

where the thresholds are chosen such that \( \delta_q = 10^{a(q+2)} \sqrt{nq^{3/2}} \) for \( q \in \{2, \ldots, q_{\text{max}} - 2\} \), \( \delta_{q_{\text{max}}} = \frac{25}{q_{\text{max}}^{3/2}} \) and \( \delta_{q_{\text{max}}+2} = 4 \sqrt{2 \log n} \) (the value of \( a \) being the same as in Section 2.3).

**Theorem 2.8.** Consider the gOSC model (4) with known variance \( \sigma^2 = 1 \). There exist universal positive constants \( c_1, c_2, c_3 \), and \( n_0 \) such that the following holds. For any \( n \geq n_0 \), for any integer \( k \in [1, n-1] \), any \( \theta \in \mathbb{R} \), and any \( \mu \in \mathcal{M}_k \), the adaptive estimator \( \hat{\theta}_{\text{ad}} = \hat{\theta}_q \) satisfies

\[
\mathbb{P}_{\theta, \mu} \left[ |\hat{\theta}_{\text{ad}} - \theta| > c_1 \frac{\log^2 (1 + \sqrt{\frac{k}{n-k}})}{\log^{3/2} (1 + (k^2/n)^{1/3})} \right] \leq c_2 \left( \frac{\sqrt{n}}{k} \vee \sqrt{n} \right)^{4/3} \log^3 \left( \frac{k \vee (2 \sqrt{n})}{\sqrt{n}} \right),
\]

and

\[
\mathbb{E}_{\theta, \mu} \left[ |\hat{\theta}_{\text{ad}} - \theta| \right] \leq c_3 \frac{\log^2 (1 + \sqrt{\frac{k}{n-k}})}{\log^{3/2} (1 + (k^2/n)^{1/3})}.
\]

A proof is given in Section S-3.4. The risk bound in (24) matches the minimax lower bound of Theorem 2.1 for all \( k \in [1, n-1] \). The estimator \( \hat{\theta}_{\text{ad}} = \hat{\theta}_q \) is therefore minimax adaptive with respect to \( k \). The qualitative behavior of \( \hat{\theta}_{\text{ad}} = \hat{\theta}_q \) is illustrated in Section S-6.3 via numerical experiments.

**Remark 2.9.** Theorem 2.8 implies that the estimation rate is not affected by knowledge of \( k \). This contrasts with other statistical setting where costless adaptation is impossible. For instance, the optimal adaptive rate is slower than the minimax rate in signal detection for the smooth Gaussian white noise model, see [66]. A related phenomenon appears also in extreme value theory, for the problem of adaptively estimating the tail coefficient [12]. In both cases, in the non-adaptive setting, the bias and deviations are of the same order. Intuitively, the deviations have to be slightly inflated to enable adaptivity, which leads to slower rates. This does not happen in our problem because the bias of the statistic turns out to be significantly larger than the deviations (see the proof in the supplementary material for more details).
and therefore a slight inflation of the deviations to account for adaptivity has no impact on the rate.

3. Estimation of $\theta$ in the general OSC model with known variance. In this section, we study the estimation problem in the general OSC model (1). Hence, the contaminations are no longer assumed to be Gaussian. Throughout this section, $\sigma$ is still assumed to be known and equal to 1 (for unknown variance, see Section 4). Recall that the $L_1$ minimax risk is defined in (7).

3.1. Lower bound in the OSC model with known variance. We first show that estimating $\theta$ is more difficult in this model than for the gOSC case.

**Theorem 3.1.** There exists a universal positive constant $c$ such that for any positive integer $n$ and for any integer $k \in [1, n - 1]$,

$$R_k[n] \geq c \frac{\log \left( \frac{n}{n-k} \right)}{\log^{1/2}(1 + \frac{k^2}{n^2})}.$$  

A proof is provided in Section S-2.2. Let us comment briefly the order of this lower bound, by considering again the three aforementioned regimes (see also Table 1):

- For $k \leq \sqrt{n}$, the lower bound (25) is of order $n^{-1/2}$, which is the parametric rate, hence is the same as for the Gaussian case;
- For $k \in (\sqrt{n}, \zeta n)$ with $\zeta \in (0, 1)$, the lower bound is of the order $(k/n)\log^{-1/2}(k^2/n)$, and so is strictly slower than with the Gaussian assumption (by an additional factor of order $\log(k^2/n)$). In particular, in the non-sparse case $k = \lceil n/2 \rceil$, this gives a lower bound of order $\log^{-3/2}(n)$ (by contrast to the $\log^{-3/2}(n)$ bound in the Gaussian model);
- For $k \in [\zeta n, n-1]$ with $\zeta \in (0, 1)$, the lower bound is of order $\log(n/(n-k))\log^{-1/2}(n)$. Compared to the gOSC model, there is an additional factor of order $\log(n)/\log(n/(n-k))$. Nevertheless, in the extreme case $k = n - 1$, the two lower bounds are both of order $\log^{1/2}(n)$.

3.2. Upper bound in the OSC model with known variance. In this subsection, we introduce a bias-corrected quantile estimator that matches the minimax lower bound of Theorem 3.1. Consider some $\pi \in \mathcal{M}_k$. Let $\xi = (\xi_1, \ldots, \xi_n)$ denote a standard Gaussian vector. The starting point is the following: on the one hand, all random variables $Y_i - \theta$ stochastically dominate $\xi_i$ so that $Y_{(q)} - \theta \succeq \xi_{(q)}$. On the other hand, $Y_{(q)}$ is stochastically...
dominated by the $q$-th smallest observation among the non-contaminated data $Y_j$. As a consequence, we have

\begin{equation}
(26) \quad \xi(q) \leq Y(q) - \theta \leq \xi_{(q,(n-k))},
\end{equation}

where we recall that $\xi_{(q,(n-k))}$ is the $q$-th largest observation among the $n-k$ first observations of $\xi$. Since $\xi(q)$ is concentrated around $\Phi^{-1}(q/n)$, this leads to introducing the debiased estimator

\begin{equation}
(27) \quad \tilde{\theta}_q = Y(q) + \Phi^{-1}(q/n) \quad 1 \leq q \leq \lceil n/2 \rceil.
\end{equation}

In view of (12), we have that $\tilde{\theta}_1 = \hat{\theta}_{\min}$ while $\tilde{\theta}_{\lceil n/2 \rceil}$ is almost equal to the empirical median $\hat{\theta}_{\text{med}}$ (up to the additive $\Phi^{-1}([n/2]/n)$ term which is of order $1/n$ and so is negligible). The following theorem bounds the error of $\tilde{\theta}_q$ for a wide range of $q$.

**Theorem 3.2.** Consider the OSC model (1) with known variance $\sigma^2 = 1$. There exist universal positive constants $c_1, c_2, c'_2, c_3, c_4$ such that the following holds. For any integer $k \in [1,n-1]$, any integer $q$ such that $c_4 \log n \leq q \leq (0.7(n-k)) \wedge \lceil n/2 \rceil$, any $\theta \in \mathbb{R}$ and any $\pi \in \mathcal{M}_k$, the estimator $\tilde{\theta}_q$ satisfies

\begin{equation}
(28) \quad P_{\theta,\pi} \left[ -c_2 \sqrt{\frac{x}{q \log(n-k) \vee 1}} \leq \tilde{\theta}_q - \theta \leq c_1 \frac{\log \left( \frac{n}{n-k} \right)}{\sqrt{\log \left( \frac{n}{n-k} \right) \vee 1}} + c_2 \sqrt{\frac{x}{q \log(n-k) \vee 1}} \right] \geq 1 - 2e^{-x},
\end{equation}

for all $0 < x < c_3 q$, and

\begin{equation}
(29) \quad \mathbb{E}_{\theta,\pi} \left[ |\tilde{\theta}_q - \theta| \right] \leq c_1 \frac{\log \left( \frac{n}{n-k} \right)}{\sqrt{\log \left( \frac{n}{n-k} \right) \vee 1}} + c'_2 \frac{1}{\sqrt{q \log(n-k) \vee 1}}.
\end{equation}

A proof is given in Section S-3.5. The risk bound in (29) exhibits a bias/variance trade-off as a function of $q$ via the quantities

\begin{align*}
b(q) &= \frac{\log \left( \frac{n}{n-k} \right)}{\sqrt{\log \left( \frac{n}{n-k} \right) \vee 1}} ; \\
s(q) &= \frac{1}{\sqrt{q \log(n-k) \vee 1}}.
\end{align*}
The quantity $s(q)$ is a deviation term that decreases with $q$ and whose minimum is of the order of $n^{-1/2}$. This minimum is achieved for $q = [n/2]$ and the corresponding estimator is close to the empirical median. The quantity $b(q)$ is a bias term which increases slowly with $q$. Its minimum is of the order of $\log \left( \frac{n}{n-k} \right) \log^{-1/2} (n-k)$ and is achieved for $q$ constant (or of the order of $\log n$). The corresponding estimators are extreme quantiles such as $\tilde{\theta}_1 = \hat{\theta}_{\min}$.

Note also that the condition $c_4 \log(n) \leq q \leq 0.7(n-k)$ cannot be met when $k$ is too close to $n$ (i.e. $n-k < (c_4/0.7) \log(n)$). Hence, Theorem 3.2 is silent in that regime. Nevertheless, this case is addressed by the minimum estimator $\tilde{\theta}_1 = \theta_{\min}$ already studied in Proposition 2.3.

To achieve the minimax risk, it remains to suitably choose $q$ as a function of $k$. Examining $b(.)$ and $s(.)$, we see that when $k$ increases we should decrease $q$ (and therefore use a more extreme quantile) in order to decrease the bias. More precisely, we define

$$q_k = \begin{cases} 
  \left\lceil \frac{n}{2} \right\rceil & \text{if } k \in [1, 4\sqrt{n}) ; \\
  \left\lceil \frac{n^{5/4}}{k^{1/2}} \right\rceil \log_2(k) & \text{if } k \in [4\sqrt{n}, n - n^{4/5}] ; \\
  1 & \text{if } k \in (n - n^{4/5}, n - 1]. 
\end{cases}$$

In the very sparse situation ($k \leq 4\sqrt{n}$), $\tilde{\theta}_{q_k}$ corresponds to the empirical median. As $k$ increases towards $n$, $q_k$ goes smoothly to $n^{1/4}$. Finally, when $k$ is very close to $n$, we consider the minimum estimator $\tilde{\theta}_1$. Other choices of $q_k$ may also lead to optimal risk bounds and the choice (30) is made to simplify the proofs.

**Corollary 3.3.** Consider the OSC model (1) with known variance $\sigma^2 = 1$. There exist universal positive constants $c$ and $n_0$ such that the following holds. For any integer $n \geq n_0$, any integer $k \in [1, n-1]$, any $\theta \in \mathbb{R}$ and any $\pi \in \mathcal{M}_k$, the estimator $\tilde{\theta}_{q_k}$ satisfies

$$\mathbb{E}_{\theta, \pi} \left[ |\tilde{\theta}_{q_k} - \theta| \right] \leq c \frac{\log \left( \frac{n}{n-k} \right)}{\log^{1/2} \left( 1 + \frac{k^2}{n} \right)}.$$

A proof is given in Section S-3.5. The estimation rate of the estimator $\tilde{\theta}_{q_k}$ matches the minimax lower bound given in Theorem 3.1. However, it is not adaptive because it uses the value of $k$.

**3.3. Adaptive estimation in the OSC model with known variance.** We now provide a procedure that adapts to $k$, by following a Goldenshluger-Lepski approach. Let $Q$ denote the collection of values of $q_k$ when $k$ goes
from 1 to $n - 1$. This collection contains 1, $[n/2]$ and a dyadic sequence from $n^{1/4}$ to $n/2$ (roughly). To build an adaptive procedure, we select among the estimators $\{\tilde{\theta}_q, q \in Q\}$ in the following way:

$$(32) \quad \hat{q} = \max \{q \in Q \text{ s.t. } |\hat{\theta}_q - \hat{\theta}_{q'}| \leq \delta_{q'} \text{ for all } q' < q\},$$

where

$$(33) \quad \delta_q = c_0 \begin{cases} \frac{\sqrt{\log(n)}}{n^{1/6}} & \text{if } q < \sqrt{2n^{1/4}}; \\ \frac{q^{2/3}}{\sqrt{\log(n^q)}} & \text{otherwise}, \end{cases}$$

where the constant $c_0$ is large enough (depending on the constants $c_1$ and $c'_2$ of Theorem 3.2). Note that at most three elements in $Q$ are less than $\sqrt{2n^{1/4}}$.

**Proposition 3.4.** Consider the OSC model (1) with known variance $\sigma^2 = 1$. There exist universal positive constants $c$ and $n_0$ such that the following holds. For any integer $n \geq n_0$, any integer $k \in [1, n - 1]$, any $\theta \in \mathbb{R}$, and any $\pi \in \mathcal{M}_k$, the estimator $\tilde{\theta}_{ad} = \tilde{\theta}_{\hat{q}}$ (see (30) and (32)) satisfies

$$\mathbb{E}_{\theta, \pi, \sigma} \left[ |\tilde{\theta}_{ad} - \theta| \right] \leq c \frac{\log \left( \frac{n}{n - k} \right)}{\log^{1/2} \left( 1 + \frac{k^2}{n} \right)}.$$ 

A proof is given in Section S-3.5. The above result shows that, as in the Gaussian case, adaptation with respect to $k$ can be achieved without any loss.

4. Estimation in the OSC model with unknown variance. In this section, we consider the OSC model (1) for which the noise variance $\sigma^2$ is unknown. We derive the minimax risks and estimators for $\theta$ and $\sigma$ in that setting.

4.1. Lower bound in the OSC model with unknown variance. First note that, obviously, the lower bound (25) for estimating $\theta$ is also a valid lower bound for the minimax risk

$$\inf_{\theta} \sup_{\theta \in \mathbb{R}, \sigma > 0, \pi \in \mathcal{M}_k} \mathbb{E}_{\theta, \pi, \sigma} \left[ \left| \frac{\hat{\theta} - \theta}{\sigma} \right| \right],$$

corresponding to the OSC model (1) where $\sigma$ is unknown.

Now, let us provide a lower bound for the estimation risk of $\sigma$. Since the problem of estimating $\sigma$ when $\theta$ is unknown is at least as difficult as
the problem when it is known that $\theta = 0$, we provide a lower bound for
the latter. Interestingly, the procedure that will be presented in Section 4.2
achieves the same risk even in the more general setting where $\theta \in \mathbb{R}$ is
unknown. We introduce the minimax risk
\begin{equation}
\overline{R}_v[k,n] = \inf_{\tilde{\sigma}} \sup_{\sigma > 0, \pi \in \mathcal{M}_k} \mathbb{E}_{0,\pi,\sigma} \left[ \left| \tilde{\sigma} - \sigma \right| / \sigma \right].
\end{equation}

The following theorem provides a lower bound for $\overline{R}_v[k,n]$ (and therefore
also a lower bound on the minimax risk with arbitrary unknown $\theta$):

**Theorem 4.1.** There exists a universal positive constant $c$ such that for
any integer $n \geq 2$ and any integer $k \in [1, n-2]$, we have
\begin{equation}
\overline{R}_v[k,n] \geq c \frac{\log \left( \frac{n}{n-k} \right)}{\log \left( 1 + \frac{k}{n^2} \right)}.
\end{equation}

A proof is given in Section S-2.3. For $k \leq \sqrt{n}$, the lower bound (35)
is of order $n^{-1/2}$. For $k \in [\sqrt{n}, \zeta n]$ (with $\zeta \in (0,1)$ fixed), the risk is of
order $k/[n \log(k^2/n)]$ which is faster by a $\log^{1/2}(k^2/n)$ term than for mean
estimation. When $n - k = n^\gamma$ with $\gamma \in (0,1)$ (almost no uncontaminated
data), the relative risk of convergence is at least of constant order.

In the next section, we prove that these lower bounds on $\theta$ and $\sigma$ are all
sharp (up to numerical constants).

4.2. Upper bound for the OSC model with unknown variance. Since the
model is translation invariant, the variance can be estimated without knowing $\theta$. This is done by considering a rescaled difference of empirical quantiles.
More precisely, for two positive integers $1 \leq q' \leq q \leq n$, let
\begin{equation}
\tilde{\sigma}_{q,q'} = \frac{Y(q) - Y(q')}{\Phi^{-1}(q'/n) - \Phi^{-1}(q/n)},
\end{equation}
with the convention $0/0 = 0$. When $k = 0$ (no contamination), $Y(q)$ (resp.
$Y(q')$) should be close to $\theta - \sigma \Phi^{-1}(q/n)$ (resp. $\theta - \sigma \Phi^{-1}(q'/n)$) so that,
intuitively, $\tilde{\sigma}_{q,q'}$ should be close to $\sigma$. Then, to estimate $\theta$, we simply plug
$\tilde{\sigma}_{q,q'}$ into the quantile estimator considered in Section 3.2. More precisely,
we consider
\begin{equation}
\tilde{\theta}_{q,q'} = Y(q) + \tilde{\sigma}_{q,q'} \Phi^{-1} \left( \frac{q}{n} \right).
\end{equation}
Given $k \in \{1, \ldots, n-1\}$, $q_k$ is taken as in (30) and

\begin{equation}
q'_k = \begin{cases} 
\lceil n/3 \rceil & \text{if } k \in [1, 4\sqrt{n}) ; \\
\left\lfloor \frac{n^{7/4}}{k^{3/2}} \right\rfloor (\log_2 n) & \text{if } k \in [4\sqrt{n}, n-4^{1/5}) ; \\
1 & \text{if } k \in (n - n^{4/5}, n - 2] .
\end{cases}
\end{equation}

For sparse contaminations ($k < 4\sqrt{n}$), $\tilde{\sigma}_{q_k,q'_k}$ is a rescaled difference of the empirical median and the empirical quantile of order $1/3$. For a larger number of contaminations, more extreme quantiles are considered. For $k \geq n - n^{4/5}$, we simply take $\tilde{\sigma}_{q_k,q'_k} = 0$.

**Proposition 4.2.** Consider the OSC model (1) with unknown variance $\sigma^2$ and the quantities $q_k$ and $q'_k$ defined in (30) and (38). There exist universal positive constants $c$, $c'$, and $n_0$ such that the following holds. For any $n \geq n_0$, for any integer $k \in [1, n-2]$, any $\theta \in \mathbb{R}$, any $\sigma > 0$ and any $\pi \in \mathcal{M}_k$, we have

\begin{align}
\mathbb{E}_{\theta, \pi, \sigma} \left[ |\tilde{\sigma}_{q_k,q'_k} - \sigma|/\sigma \right] &\leq c \frac{\log \left( \frac{n}{n-k} \right)}{\log(1 + \frac{k}{n^{1/2}})} ; \\
\mathbb{E}_{\theta, \pi, \sigma} \left[ |\tilde{\theta}_{q_k,q'_k} - \theta|/\sigma \right] &\leq c' \frac{\log \left( \frac{n}{n-k} \right)}{\log^{1/2}(1 + \frac{k}{n})} .
\end{align}

A proof is given in Section S-3.5. The above proposition together with the lower bounds of Section 4.1 implies that $\tilde{\sigma}_{q_k,q'_k}$ and $\tilde{\theta}_{q_k,q'_k}$ are minimax estimator of $\sigma$ and $\theta$, respectively.

Interestingly, the estimators $\tilde{\sigma}_{q_k,q'_k}$ and $\tilde{\theta}_{q_k,q'_k}$ do not require the knowledge of either $\theta$ or $\sigma$ whereas our lower bounds were respectively restricted to known $\theta$ and known $\sigma$ settings. This entails that the knowledge of one parameter ($\theta$ or $\sigma$) does not significantly ease the estimation of the other.

5. Controlled selection of outliers. In this section, we consider the OSC model (1) (with unknown variance) and now turn to the identification of the outliers. As described in Section 1.3, this can be reformulated as a multiple testing problem (see also the notation therein).

5.1. Rescaled $p$-values. As already discussed in Section 1.3, ensuring good multiple testing properties in the OSC model is challenging because the scaling parameters $\theta$ and $\sigma$ are unknown. A natural approach is then to use the rescaled observations $Y'_i = (Y_i - \bar{\theta})/\hat{\sigma}$, $1 \leq i \leq n$, where $\bar{\theta}$, $\hat{\sigma}$ are
some suitable estimators of \( \theta \) and \( \sigma \). To formalize further this idea, let us consider the corrected \( p \)-values
\[
(41) \quad p_i(u, s) = \frac{\Phi \left( \frac{Y_i - u}{s} \right)}{s}, \quad u \in \mathbb{R}, \ s > 0, \ 1 \leq i \leq n.
\]
The perfectly corrected \( p \)-values thus correspond to
\[
(42) \quad p^*_i = p_i(\theta, \sigma), \ 1 \leq i \leq n.
\]
These oracle \( p \)-values cannot be used in practice, because they depend on the unknown parameters \( \theta \) and \( \sigma \). Our general aim is to build estimators \( \hat{\theta}, \hat{\sigma} \) such that the theoretical performance of the corrected \( p \)-values \( p_i(\hat{\theta}, \hat{\sigma}) \) mimic those of the oracle \( p \)-values \( p^*_i \), when plugged into standard multiple testing or post hoc procedures. Although the use of modified \( p \)-values and plug-in estimators has often been advocated since the seminal work of Efron [26], proving that the behavior of the plug-in procedure is asymptotically similar to that of the oracle procedure is, to our knowledge, new. The challenge is to precisely quantify how the estimation error affects the FDP/TDP metrics. For this, a key point is the following relation between \( p_i(u, s) \) and \( p^*_i \):

\[
(43) \quad \{ p_i(u, s) \leq t \} = \{ p^*_i \leq U_{u,s}(t) \}, \ i \in \{1,\ldots,n\}, \ t \in [0,1],
\]
where
\[
(44) \quad U_{u,s}(t) = \frac{s \Phi^{-1}(t) + u - \theta}{\sigma}; \quad U_{u,s}^{-1}(v) = \frac{\sigma \Phi^{-1}(v) + \theta - u}{s}.
\]
Furthermore, a useful property is that the order of the \( p \)-values does not change after rescaling. We will denote
\[
(45) \quad 0 = p(0)(u, s) \leq p(1)(u, s) \leq \cdots \leq p(n)(u, s),
\]
the ordered elements of \( \{ p_i(u, s), 1 \leq i \leq n \} \). We also denote by \( 0 = p(0;H_0)(u, s) \leq p(1;H_0)(u, s) \leq \cdots \leq p(n;H_0)(u, s) \) the ordered elements of the subset \( \{ p_i(u, s), i \in H_0 \} \), that is, of the \( p \)-value set corresponding to false outliers (or, equivalently, true null hypotheses).

5.2. Upper-biased estimators. This section provides estimators \( \tilde{\theta}_+, \tilde{\sigma}_+ \) that will be suitable to make the \( p \)-value rescaling. They are similar to the estimators introduced in Sections 3.2 and 4.2. However, since minimax estimation and false outliers control do not use the same risk metrics, we need to slightly modify these estimators, especially by making them upper-biased (which roughly means that the null hypotheses are favored).
For $q_n = \lfloor n^{3/4} \rfloor$ and $q'_n = \lfloor n^{1/4} \rfloor$, let us consider

$$
\begin{cases}
\tilde{\theta}_+ = Y_{(q_n)} + \tilde{\sigma}_+ \Phi^{-1} \left( \frac{q_n}{n} \right); \\
\tilde{\sigma}_+ = \Phi^{-1} \left( \frac{q'_n}{n-k_0} \right) - \Phi^{-1} \left( \frac{q_n}{n} \right),
\end{cases}
$$

for some parameter $k_0 \leq \lfloor 0.9 n \rfloor$. The key difference with the estimators $\tilde{\theta}_{q,q'}, \tilde{\sigma}_{q,q'}$ of Section 4 is the quantity $k_0$ appearing in the denominator of $\tilde{\sigma}_+$. The following result holds.

**Proposition 5.1.** Consider the OSC model (1) with unknown variance $\sigma^2$. Then there exist two universal positive constants $c, c'$ such that the following holds for any positive integer $n$, for any $\theta \in \mathbb{R}$, $\sigma > 0$, for any $\pi \in \mathcal{M}_k$ with $k = \lfloor 0.9 n \rfloor$. Choosing $k_0$ such that $n_1(\pi) \leq k_0 \leq \lfloor 0.9 n \rfloor$ within the estimators $\tilde{\theta}_+, \tilde{\sigma}_+$ (46), we have

$$
\begin{align*}
\mathbb{P}_{\theta,\pi,\sigma}(\tilde{\theta}_+ - \theta \leq -\sigma n^{-1/16}) &\leq c/n; \\
\mathbb{P}_{\theta,\pi,\sigma}(\tilde{\sigma}_+ - \sigma \leq -n^{-1/16}) &\leq c/n;
\end{align*}
$$

$$
\begin{align*}
\mathbb{P}_{\theta,\pi,\sigma}(\left| \tilde{\theta}_+ - \theta \right| \geq \sigma \left( c'(k_0/n) \log^{-1/2}(n) + n^{-1/16} \right)) &\leq c/n; \\
\mathbb{P}_{\theta,\pi,\sigma}(\left| \tilde{\sigma}_+ - \sigma \right| \geq \sigma \left( c'(k_0/n) \log^{-1}(n) + n^{-1/16} \right)) &\leq c/n.
\end{align*}
$$

Proposition 5.1 is proved in Section S-3.5. It is closely related to Theorem 3.2 above, although the statement is slightly different because of the introduced bias in the estimators. Inequalities (47–48) entail that the estimators are (with high probability) above the targeted quantity minus a polynomial term, which will be particularly suitable for obtaining a control on the false positives (FDR control and post hoc bounds). Inequalities (49–50) are two-sided, which is useful for studying the power (TDP) of the rescaled procedures: there is an additional error term of order $(k_0/n) \log^{-a}(n)$, $a \in \{1/2, 1\}$, where $k_0$ corresponds to a known upper bound of the number of contaminated coordinates of $\pi$.

The assumption $n_1(\pi) \leq k_0 \leq \lfloor 0.9 n \rfloor$ in Proposition 5.1 is not very restrictive: it means that the number of outliers is bounded by above by some quantity $k_0$, which is used in the definition of the estimators (46). For instance, taking $k_0 = \lfloor 0.7 n \rfloor$ means that we assume that there is no more than 70% of outliers in the data.

Finally note that our multiple testing analysis will only rely on the deviation bounds (47–50). As a consequence, other estimators satisfying these properties can be used for scaling the $p$-values.
5.3. **FDR control and power optimality for selected outliers.** The Benjamini-Hochberg (BH) procedure, introduced in [4], has probably been the most widely used multiple testing procedure. Here, the rescaled BH procedure (of nominal level $\alpha$), denoted $\text{BH}_\alpha(u,s)$ is defined from the $p$-value family $p_i(u,s), 1 \leq i \leq n$, as follows:

- Order the $p$-values as in (45);
- Consider $\hat{\ell}_\alpha(u,s) = \max\{\ell \in \{0,1,\ldots,n\} : p_\ell(u,s) \leq \alpha \ell / n\}$;
- Reject $H_{0,i}$ for any $i$ such that $p_i(u,s) \leq \hat{\ell}_\alpha (u,s)$, for $\hat{\ell}_\alpha (u,s) = \alpha \hat{\ell}_\alpha (u,s) / n$.

The procedure, identified to the set of selected outliers, is then given by

$$\text{BH}_\alpha(u,s) = \{1 \leq i \leq n : p_i(u,s) \leq \hat{\ell}_\alpha (u,s)\}.$$  

The classical FDR-controlling result of [4, 5] can be re-interpreted as follows: the BH procedure using the perfectly corrected $p$-values (42), that is, $\text{BH}^*_\alpha = \text{BH}_\alpha(\theta,\sigma)$, satisfies

$$E_{\theta,\pi,\sigma}(\text{FDP}(\pi,\text{BH}^*_\alpha)) = \frac{n_0}{n} \leq \alpha, \text{ for all } \theta, \pi, \sigma.$$  

This comes from the fact that the perfectly corrected $p$-values (42) are independent, with uniform marginal distributions under the null hypothesis.

Recall the estimators $\tilde{\theta} +$ and $\tilde{\sigma} +$ defined in (46) with the tuning parameter $k_0$. The next result gives the behavior of the rescaled procedure $\text{BH}_\alpha(\tilde{\theta} +, \tilde{\sigma} +)$ both in terms of FDP and TDP.

**Theorem 5.2.** Consider the OSC model (1) with unknown variance $\sigma^2$. There exists a universal positive constant $c$ such that the following holds. Considering the estimators $\tilde{\theta} +, \tilde{\sigma} +$ given by (46), for any $\alpha \in (0,0.4), \theta \in \mathbb{R}$, $\sigma > 0$, and any $\pi \in \mathcal{M}_{[0.9n]}$, if $k_0$ in $\tilde{\sigma} +$ satisfies $n_1(\pi) \leq k_0 \leq [0.9n]$, we have

$$E_{\theta,\pi,\sigma}(\text{FDP}(\pi,\text{BH}_\alpha(\tilde{\theta} +, \tilde{\sigma} +))) - \frac{n_0}{n} \alpha \leq c \log(n)/n^{1/16}.$$  

Additionally, for any sequence $\epsilon_n \in (0,1)$ tending to zero with $\epsilon_n \gg \log^{-1/2}(n)$, for any sequence $\pi = \pi_n$, $k_0 = k_{0,n}$ with $n_1(\pi_n) \leq k_{0,n} \leq [0.9n]$ and $n_1(\pi_n)/n \asymp k_{0,n}/n$, we have for all $\theta \in \mathbb{R}, \sigma > 0$,

$$\limsup_n \left\{ -E_{\theta,\pi_n,\sigma}(\text{TDP}(\pi_n,\text{BH}^*_\alpha)) - E_{\theta,\pi_n,\sigma}(\text{TDP}(\pi_n,\text{BH}_\alpha(1+\epsilon_n)(\tilde{\theta} +, \tilde{\sigma} +))) \right\} \leq 0.$$
In a nutshell, inequalities (52) and (53) show that the selection procedure $BH_{\alpha}(\hat{\theta}_+, \hat{\sigma}_+)$ behaves similarly to the oracle procedure $BH^*_\alpha$, both in terms of false discovery rate control and power. Note that the power result (53) is shown to be valid only when $k_0/n$ is taken of the same order as $n_1/\pi n$, that is, for a procedure using information on $n_1/\pi n$. On the positive side, it entails an optimal procedure even without sparsity: if we know that $n_1/\pi n$ is of the order of a constant, then using the estimators $\hat{\theta}_+, \hat{\sigma}_+$ for say $k_0 = \lfloor 0.5n \rfloor$, we will always achieve both (52) and (53). In particular, Theorem 5.2 can be seen as a first step towards the validation of Efron’s empirical null principle, in a specific one-sided situation.

The proof of this theorem is given in Section S-4.1. Compared to the usual FDR proofs of the existing literature, there are two additional difficulties: first the independence assumption between the corrected $p$-values is not satisfied anymore, because the correction terms are random; second, the quantity $FDP(\pi, BH_{\alpha}(\hat{\theta}_+, \hat{\sigma}_+))$ is not monotone in the estimators $\hat{\theta}_+, \hat{\sigma}_+$, because of the denominator of the FDP. However, the specific properties of $\hat{\theta}_+, \hat{\sigma}_+$ given in Proposition 5.1 will be enough to get our result: first, these estimators are biased upwards with an error term vanishing at a polynomial rate $n^{-1/16}$, which is enough for false positive control. As for the power, we consider the bias downwards, which is of order $(k_0/n) \log^{-a}(n)$, $a \in \{1/2, 1\}$. It turns out that the error term induced in the power is of order $(k_0/n) \log^{-a}(n) \log(n/n_1)$, which tends to 0 when $k_0/n \approx n_1/n$ both in the sparse and non-sparse cases.

The other multiple testing inferences described in the introduction (Section 1.3): post-hoc bounds, decorrelation and the corresponding numerical experiments are postponed to the supplementary material (see Section S-1).

Remark 5.3. Let us define the risk of a procedure as $FDR + FNR$, where $FNR$ denotes the average of $1 - TDP$. Theorem 5.2 shows that the risk of the selection procedure $BH_{\alpha(1+\epsilon_n)}(\hat{\theta}_+, \hat{\sigma}_+)$ is smaller than the risk $R^*(n, \pi_n, \alpha)$ of the oracle procedure $BH^*_\alpha$ plus some vanishing terms. Interestingly, the risk $R^*(n, \pi_n, \alpha)$ has been studied in [1] for a particular choice of $\pi_n$ coming from the classical Gaussian sequence model in the specific Donoho-Jin sparse regime [23]. It is shown that $R^*(n, \pi_n, \alpha)$ tends to zero when the signal strength is above the estimation boundary (as defined, e.g., in [23]), for some proper calibration of $\alpha_n$, typically $\alpha_n = 1/(\log n)$. Inspecting our proofs (see Section S-4.1), we can derive that the procedure $BH_{\alpha_n(1+\epsilon_n)}(\hat{\theta}_+, \hat{\sigma}_+)$ also has a risk tending to zero in that regime. In addition, the work [60] provides a polynomial decaying rate for $R^*(n, \pi_n, \alpha_n)$ when $\alpha_n = n^{-\kappa}$ for some $\kappa > 0$. Again, inspecting our proofs (see Section S-4.1), this also holds for the risk
of $BH_{\alpha_n(1+\epsilon_n)}(\tilde{\theta}_+, \tilde{\sigma}_+)$, as our remainder term also vanishes at a polynomial rate in that regime, although our exponent may be smaller (e.g., $n^{-c}$ with $c < 1/16$). Finally, we emphasize that Theorem 5.2 does not rely on the Donoho-Jin sparse regime. Hence, this type of power result is new in multiple testing literature to the best of our knowledge.

6. Discussion. In this section, we describe some possible extensions and future research directions.

Knowledge of the variance in the gOSC model. In the gOSC model (Section 2), we assumed exact knowledge of the variance. Unfortunately, plugging an estimator of the variance into the empirical Laplace transform from (17) seems to deteriorate the deviations of the corresponding estimator $\hat{\psi}_{q, \lambda}(u)$. For this reason, we were not able to derive an estimator of $\theta$ with $\sigma$ unknown achieving the optimal rate $R[k, n]$ in the intermediate regime where $k \in [\sqrt{n}; cn]$, $c \in (0, 1)$. Since the gOSC model is a specific case of the OSC model, it is possible to achieve the rate $\tilde{R}[k, n]$ using $\tilde{\theta}_{q, q_k}$ (37). Since there is $\log(k^2/n)$ gap between the gOSC and OSC models, it still remains unclear if perfect adaptation to unknown $\sigma$ is possible in the gOSC model or if a logarithmic price must be paid.

Simultaneous adaptation to $k$ and $\sigma$ in the OSC model. In Sections 3 and 4, we assumed that either $\sigma$ or $k$ is known. If both quantities are unknown, one needs to select a suitable estimator $\tilde{\theta}_{q, q_k}$ in the collection $\{\tilde{\theta}_{q, q_k}\}$ introduced in Section 4. This turns out to be possible using a Goldenshluger-Lepski scheme in the spirit of the one studied in Section 3.3 provided that we have at our disposal an estimator $\tilde{\sigma}$ of $\sigma$ satisfying $|\tilde{\sigma}^2 - \sigma^2|/\sigma^2 \in [1/2, 2]$ with high probability. In view of Section 4, this is the case when $k$ is not close to $n$ ($k \leq n - n/e^{c\log^{3/2}(n)})$. As a result, simultaneous adaptation to $k$ and $\sigma$ is possible provided that the true number of contaminations is not too large. If $k$ is allowed to be as large as $n - 2$, it remains an open question whether there is price to pay.

Multivariate one-Sided contaminations. Throughout this manuscript, we assumed that the observations are univariate. One could think of a multivariate extension: $Y_i = \theta + \sigma \varepsilon_i$ where $Y_i \in \mathbb{R}^d$ and $\varepsilon_i$ is either distributed as a standard normal distribution or, for a contaminated observation, satisfies $\varepsilon_{i,j} \sim \mathcal{N}(0, 1)$ for each coordinate $j = 1, \ldots, d$. In Huber’s original multivariate contamination model, it is known that estimating each coordinate $\theta$ independently is suboptimal. This led Huber to introduce Tukey depth estimators [42]. More recently, polynomial time estimators were proved to achieve similar performances [21] (up to some logarithmic loss). It would be
interesting to investigate whether it is possible to build upon the one-sided assumption to improve their convergence rates.

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**SUPPLEMENTARY MATERIAL**

**Supplement to “Estimating minimum effect with outlier selection”**

(doi: COMPLETED BY THE TYPESETTER; .pdf). This supplement contains additional materials for the outlier selection problem, the proofs of the paper, auxiliary results and numerical experiments.

**References.**


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