Sparse Analysis Model Based Dictionary Learning for Signal Declipping

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Abstract—Clipping is a common type of distortion in which the amplitude of a signal is truncated if it exceeds a certain threshold. Sparse representation has underpinned several algorithms developed recently for reconstruction of the original signal from clipped observations. However, these declipping algorithms are often built on a synthesis model, where the signal is represented by a dictionary weighted by sparse coding coefficients. In contrast to these works, we propose a sparse analysis-model-based declipping (SAD) method, where the declipping model is formulated on an analysis (i.e. transform) dictionary, and additional constraints characterizing the clipping process. The analysis dictionary is updated using the Analysis SimCO algorithm, and the signal is recovered by using a least-squares based method or a projected gradient descent method, incorporating the observable signal set. Numerical experiments on speech and music are used to demonstrate improved performance in signal to distortion ratio (SDR) compared to recent state-of-the-art methods including A-SPADE and ConsDL.

Index Terms—ASimCO, clipping signal, nonlinear measurement, sparse analysis.

I. INTRODUCTION

S PARSE representation plays an important role in inverse problems and signal recovery tasks, such as denoising, inpainting, declipping and super-resolution. By exploiting the sparsity of signals in some domain, such as a transform domain, the original signals can be estimated from the observed signals

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using a sparse representation model. Clipping is a common type of distortion whereby the amplitude of a signal is truncated if it exceeds a certain threshold. Clipping may occur when a signal is recorded by a sensor limited by the range of data that can be physically measured, when the signal is digitized, or when an analog or digital signal is transformed through a nonlinear system. The clipped signal $y \in \mathbb{R}^m$ is a distorted observation of the ground truth signal $x \in \mathbb{R}^m$,

$$y = f(x) \tag{1}$$

where f is a nonlinear clipping function. The observed signal y may be divided into clipped regions and reliable (i.e. unclipped) regions. Declipping aims to recover the original signal x from the clipped signal y. This can be achieved with a variety of methods, as summarized in a recent survey paper [1]. Typically, this problem can be formulated using a cost function based on reconstruction errors, together with constraints or regularizers that are consistent with the clipping process [2], [3], [4], [5], [6].

A popular model used for the declipping problem is the sparse synthesis model [7], [8]. In this model, the original signal x is represented using an overcomplete dictionary $\Omega^T \in \mathbb{R}^{m \times p}$, i.e. $x = \Omega^T \alpha$, where $\alpha \in \mathbb{R}^p$ is a sparse coefficient vector, p > m, and the superscript T denotes matrix transpose. The original signal x can be reconstructed from the observations y based on the dictionary and sparse coefficients [9], [10]. Often, l_0 or l_1 regularizations are used to enforce the sparsity constraint on the representation coefficients of the signal [11], [12], [13], [14], [15]. The dictionary Ω can be fixed [2], or learned [16], [17], [18], [19]. The declipping problem has been addressed by imposing constraints on the samples in the clipped region. More specifically, the magnitude of the original signal should be greater than or equal to that of the clipped signal in the clipped region. An alternative to the synthesis model is the sparse analysis model, where an analysis dictionary $\Omega \in \mathbb{R}^{p \times m}$ with p > m is used to sparsify the signal $x \in \mathbb{R}^m$ [20], [21], [22], i.e. making Ωx sparse. Only few algorithms, such as the A-SPADE algorithm [23], [24], [25], have considered the use of an analysis model for signal declipping, with a fixed dictionary.

A. Related Work

In [2], the original signal is estimated using the reliable (i.e. unclipped) samples from the observed signal, with the dictionary formed using positions of the unclipped samples. This algorithm

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does not involve the update of dictionary atoms as they are preselected from the unclipped samples of the observed signal. To reconstruct the samples that are clipped, additional constraints are imposed on their amplitudes. This algorithm works better for low clipping levels than for high [2]. When the clipping region is enlarged, the number of reliable (i.e. unclipped) samples that can be used for constructing the dictionary decreases, and as a result, the quality of the recovered signal deteriorates [2].

In [19], [26], the concept of clipping consistency is exploited to improve the reconstruction of the original signal, where additional constraints are enforced for signal recovery and dictionary learning so that they are "consistent" with the measurement process. For example, the magnitude of the original signal in the clipped region of the observed signal should be greater than the clipping level. Therefore, such information can be used to design consistency constraints to improve signal recovery. However, the algorithm is based on extended linear least-squares, which may be limited for the signals with rapidly changing dynamics (e.g. highly non-stationary signals). It is worth noting that the idea of measurement consistency could also be used to address the problem of signal recovery from other type of nonlinear observations such as quantized or 1-bit samples, as shown in a recent study [27]. Another algorithm called S-SPADE [23], [24] is also based on the synthesis model, which uses l_0 pseudo-norm to measure the sparsity, and l_2 norm to bound the distance between the signal and its sparse approximation.

With a similar heuristic as in the S-SPADE algorithm, the A-SPADE algorithm [23], [24], [25] uses a fixed analysis dictionary such as a discrete Fourier transform (DFT) or a discrete cosine transform (DCT) for iterative signal reconstruction, where the previous estimates are used to obtain the new estimate. There are two main steps in the A-SPADE algorithm. The first step is to obtain the representation matrix using the fixed analysis dictionary and enforce sparsity of the representation matrix. The second step is to apply a constraint on the recovered signal to meet the clipping conditions. The empirical results in [25] show that the A-SPADE algorithm.

Recent research, e.g., [27], has shown that learned dictionaries have the potential to outperform pre-defined dictionaries in representing a signal and providing recovery of the original signal from the clipped signal. Nevertheless, this research has considered only the synthesis model. To the best of our knowledge, dictionary learning with a sparse analysis model has not been applied in the context of signal declipping. In this paper, we will bridge this gap and examine the use of a dictionary learned with a sparse analysis model for signal declipping. This idea is partially inspired by the work [20], where an analysis dictionary was shown to outperform a synthesis dictionary for signal recovery from corrupted observations.

B. Contributions

Our contributions can be summarized as follows. (1) We build a new model (cost function) for signal declipping with the consistency constraint defined on an analysis dictionary. (2) We then develop a two-stage method for optimizing the cost function. This involves dictionary learning using the Analysis SimCO (ASimCO) algorithm [20] in the first stage, and signal recovery in the second stage. To recover the original signal from the clipped signal, we developed two methods based on least squares and projected gradient descent, respectively. (3) We evaluate the performance of the proposed algorithms with different parameters, in terms of signal to distortion ratio (SDR). We also compare our algorithms with two baseline methods, i.e. A-SPADE [23] and ConsDL [27], and show that our algorithms perform significantly better than the ConsDL algorithm and slightly better than the A-SPADE algorithm.

C. Notations

In this paper, uppercase letters denote matrices, e.g. $Y \in$ $\mathbb{R}^{m \times n}$, and lowercase letters denote vectors, e.g. $y \in \mathbb{R}^m$. The *i*th row of X is denoted as $X_{i,:}$, and the *j*th column of X is denoted as $X_{i,j}$. For a matrix X, $||X||_0$ denotes the l_0 pseudo-norm, measuring the sparsity, i.e. counting the number of nonzero elements of X. $||X||_F$ denotes Frobenius-norm of the matrix X. $\|y\|_1$ denotes the l_1 -norm of the vector y. The symbols \leq and \succeq denote the element-wise smaller (or equal) and larger (or equal), respectively. I represents an identity matrix, $\underline{1}$ is an all-one vector, and $\underline{1}$ an all-one matrix. The symbol $(x)_+$ denotes $\max(0, x), (X)_+$ denotes $\max(0, X), (x)_-$ denotes $-(-x)_+,$ and $(X)_{-}$ denotes $-(-X)_{+}$. The operators $\max(X, Y)$ and $\min(X, Y)$ denote taking the element-wise maximum and minimum of X and Y, respectively. We use C to denote a set. For convenience, we use Ω and Ω^T to represent the analysis and synthesis dictionary, respectively.

D. Organization of the Paper

In Section II, we provide some background materials about the Analysis SimCO algorithm. In Section III, we formulate the declipping problem. In Section IV, we present our method for the declipping problem, including the cost function and the optimization algorithms. In Section V, we present simulation studies of the proposed algorithm, with a focus on the choices of the parameters and comparisons between the least-squares based method and projected gradient descent based method for the recovery of the original signal. In Section VI, we compare our algorithm with the two baseline algorithms. Finally, Section VII draws conclusions and discusses future works.

II. ANALYSIS SIMCO ALGORITHM

In this section, we provide an overview of the ASimCO algorithm that has been already published in [20]. Such materials are not novel, but included to make the algorithms presented later self-contained. In an analysis model, the analysis dictionary $\Omega \in \mathbb{R}^{p \times m}$ is sought to sparsify x [28], [29], [30] as follows

$$a = \Omega x \quad \text{s.t.} \quad \|a\|_0 = p - l \tag{2}$$

where $x \in \mathbb{R}^m$, $a \in \mathbb{R}^p$, $0 \le l \le p$, and l is the co-sparsity [31] indicating the number of zero elements in a, and a is the analysis representation vector of the signal x with respect to Ω . In this model, if x is a clean signal, we can obtain the representation

vector a directly via multiplying x by the analysis dictionary Ω . Note that the sparsity constraint is an upper bound. In other words, if the value of Ωx is more than l co-sparse, no extra nonzero values will be added to increase its l_0 pseudo-norm. In practice, Ω can be either a pre-defined fixed analysis dictionary, or an analysis dictionary learned from observed signals. To learn the analysis dictionary, we can use an iterative algorithm alternating between two steps: analysis pursuit and analysis dictionary update. This problem is often called as analysis dictionary learning (ADL) [22], [31], [28], [32], [33]. Given X which contains the training signals, the ADL problem can be described as follows [34]

$$\left\{\hat{A},\hat{\Omega}\right\} = \arg\min_{A,\Omega} \|A - \Omega X\|_F^2 \text{ s.t. } \|A_{:,i}\|_0 = p - l, \forall i \quad (3)$$

where $X \in \mathbb{R}^{m \times n}$ is a matrix containing *n* training examples each of dimension *m*, and $A \in \mathbb{R}^{p \times n}$ is the analysis representation matrix. In order to mitigate the scale ambiguities, the ASimCO algorithm optimizes the following cost [20], [35], by imposing a constraint on the rows of Ω ,

$$\hat{A}, \hat{\Omega} \Big\} = \underset{A,\Omega}{\operatorname{arg\,min}} \|A - \Omega X\|_F^2$$
s.t.
$$\|A_{:,i}\|_0 = p - l, \forall i$$

$$\|\Omega_{j,:}\|_2 = 1, \forall j.$$

$$(4)$$

Different from other ADL algorithms, such as Analysis K-SVD [28], multiple atoms in the analysis dictionary Ω can be updated simultaneously in the ASimCO algorithm. The ASimCO algorithm alternates between two stages, i.e. analysis pursuit and analysis dictionary updating, as discussed next.

A. Analysis Pursuit

In this stage, we need to compute the analysis representation matrix A, given the analysis dictionary Ω and the signal X. Similar to (2), we can obtain A as follows

$$A = \Omega X. \tag{5}$$

Since the analysis dictionary is generated arbitrarily, it does not necessarily satisfy the co-sparsity constraints in (2), therefore a hard thresholding operation is applied to ensure the co-sparsity constraint

$$\hat{A} = HT_l(A). \tag{6}$$

This is a nonlinear operation to set the smallest l elements (in magnitude) in each column of A to zeros. Among all the matrices satisfying the co-sparsity constraints, \hat{A} is the best approximation of A in terms of the Frobenius norm error.

B. Dictionary Update

Given X and A, the analysis dictionary Ω can be found by optimizing the following cost function [35], [20]

$$\underset{\Omega}{\arg\min} g\left(\Omega\right) = \underset{\Omega}{\arg\min} \left\|A - \Omega X\right\|_{F}^{2} \quad \text{s.t.} \left\|\Omega_{j,:}\right\|_{2} = 1, \forall j.$$
(7)

The unit-norm constraint on Ω suggests that columns of Ω^T sit on the unit sphere $S = \{u \in \mathbb{R}^m : u^T u = 1\}$, i.e. $(\Omega_{j,:})^T \in S$. To optimise this cost function, we can use the gradient descent based method, as detailed in [36] and [20]. The negative gradient of the objective function in (7) with respect to Ω can be obtained as

$$H = -\nabla g(\Omega)$$

= $-\frac{\partial ||A - \Omega X||_F^2}{\partial \Omega}$ (8)
= $2AX^T - 2\Omega X X^T$.

To update Ω , a line search method accounting for the unit-norm constraint $\|\Omega_{j,:}\|_2 = 1$ was proposed in [20], where the search direction of the *j*th row of Ω is projected onto the tangent space of *S* as follows

$$\bar{h}_j = H_{j,:} \left(I - \Omega_{j,:}^T \Omega_{j,:} \right) \tag{9}$$

The *j*th row of Ω can then be updated as follows [20]

$$\hat{\Omega}_{j,:} \left(\xi \right) = \begin{cases} \Omega_{j,:} & \text{if } \|\bar{h}_{j}\|_{2} = 0, \\ \Omega_{j,:} \cos\left(\xi \|\bar{h}_{j}\|_{2} \right) + \left(\frac{\bar{h}_{j}}{\|\bar{h}_{j}\|_{2}}\right) \sin\left(\xi \|\bar{h}_{j}\|_{2} \right) \\ & \text{otherwise} \end{cases}$$
(10)

where ξ is the step size, which can be either fixed, or varied by searching for the optimal value at each iteration in terms of the golden section rule, as detailed in [36], [37]. Readers may refer to [20] for more details about the derivation of (9) and (10) by incorporating $\|\Omega_{j,:}\|_2 = 1$. For the convenience of later use, we abbreviate (10) as follows

$$\hat{\Omega} = w\left(\Omega\right). \tag{11}$$

III. MODELLING OF CLIPPED OBSERVATIONS

The clipped signal y in (1) is a distorted version of the original signal x, and can be written as follows

$$y_i = f(x_i) = \begin{cases} \theta^+ & \text{if } x_i \ge \theta^+ \\ \theta^- & \text{if } x_i \le \theta^- \\ x_i & \text{others} \end{cases}$$
(12)

where $\theta^+ > 0$ and $\theta^- < 0$ are positive and negative clipping levels, respectively, as illustrated in Fig. 1. This representation can also be written in vector form

$$y = f(x) = M^{r}x + \theta^{+}M^{c+}\underline{1} + \theta^{-}M^{c-}\underline{1}$$
(13)

where <u>1</u> is an all-one vector, M^r , M^{c+} and M^{c-} are diagonal binary sensing matrices defining the reliable, positive and negative clipped samples, respectively. In practice, we can select the maximum and minimum value of the observed signal as θ^+ and θ^- , i.e. $\theta^+ = \max(y)$, and $\theta^- = \min(y)$.

Declipping can be regarded as an inverse problem with discarded samples. Given the observed signal y with clipped samples missing, the aim is to reconstruct the original signal x from y, by solving the inverse problem [4], [16], [38], [39].

In declipping problems, the amplitudes of the original signal in the clipping region are greater than or equal to the amplitudes of the observed signal. Therefore, we can define a so-called



Fig. 1. Hard clipping: y is the observed signal, which contains the nonlinear observations of the original signal x. θ^+ and θ^- are the maximum and the minimum of the observed signal.

consistency set, as in [19], in closed form as follows

$$C(y) = \left\{ x | M^{r}y = M^{r}x, M^{c+}y \preceq M^{c+}x, M^{c-}y \succeq M^{c-}x \right\}.$$
(14)

This set contains all the possible x that could have generated the observation y that is *consistent* with the observation (i.e. measurement) process, which, in this case, is described by the clipping function.

IV. PROPOSED METHOD

A. Cost Function

Our method is built on the following cost function

$$\left\{ \hat{\Omega}, \hat{X}, \hat{A} \right\} = \underset{\Omega, X, A}{\operatorname{arg\,min}} \|A - \Omega X\|_{F}^{2}$$

s.t. $\|A_{:,i}\|_{0} = p - l, \forall i$
 $\|\Omega_{j,:}\|_{2} = 1, \forall j$
 $X \in C(Y)$ (15)

where $X \in \mathbb{R}^{m \times n}$ represents the signal to be estimated, $Y \in \mathbb{R}^{m \times n}$ represents the clipped signal, and C(Y) is a consistency set defined on matrices, i.e. replacing y by Y and x by X in (14). The cost function (15) consists of two parts. The first part, $||A - \Omega X||_F^2$, aims to learn the analysis dictionary Ω that can sparsify the signal X. The constraint on $||A_{\cdot,i}||_0$ is used to enforce the sparsity of the representation coefficients A, and the constraint on $||\Omega_{j,:}||_2$ is applied to mitigate the scale ambiguity of the analysis dictionary Ω . The second part takes into account the nonlinear observations from the clipped signal Y, and enforces the recovered signal X to be on the set C(Y). In practice, Ycould be formed by splitting the clipped signal into n segments each of dimension m. Our aim is to learn the analysis dictionary Ω and estimate the signal X, given the observed signal Y.

Algorithm 1: Proposed SAD.					
1.	Initialization: $\Omega^{[0]}, X^{[0]} = Y, t = 0, \mu$				
2.	While not converged do				
3.	Update Ω and A by solving (16) using ASimCO.				
4.	Update X by solving (17).				
5.	t = t + 1				
6	End while				

B. Optimization Process

To optimize the cost function (15), we develop a two-stage method where the unknown variables are estimated in an alternating manner, as commonly adopted in sparse representations and analysis dictionary learning.

In the first stage, given $X^{[t]}$ obtained in iteration t, $\Omega^{[t+1]}$ and $A^{[t+1]}$ are updated in the (t + 1)-th iteration, based on the following cost function

$$\left\{ \Omega^{[t+1]}, A^{[t+1]} \right\} = \underset{\Omega, A}{\operatorname{arg\,min}} \left\| A - \Omega X^{[t]} \right\|_{F}^{2}$$
s.t. $\|A_{:,i}\|_{0} = p - l, \forall i$
 $\|\Omega_{j,:}\|_{2} = 1, \forall j.$

$$(16)$$

Note that $X^{[0]}$ is initialized as Y. This stage can be achieved by the ASimCO algorithm [20] as reviewed in Section II.

The second stage is to find X, given Ω and A, under the constraint on X to enforce consistency, i.e. $X \in C(Y)$, as follows

$$\hat{X} = \underset{X}{\operatorname{arg\,min}} \|A - \Omega X\|_{F}^{2}$$
s.t. $X \in C(Y)$.
(17)

The proposed sparse analysis model based declipping (SAD) algorithm can be outlined in Algorithm 1.

To optimise the cost function (17), we present two methods, i.e. the least squares based method and projected gradient based method, respectively.

The least squares based method can be achieved by approximating the constrained optimization problem (17) as an unconstrained one. We note that the term $||M^r(Y - X)||_F^2$ promotes the recovered signal to be close to the original signal for the unclipped part of the signal, and the terms $||M^{c+}(\theta^+\underline{1} - X)_+||_F^2$ and $||M^{c-}(\theta^-\underline{1} - X)_-||_F^2$ promote the amplitudes of the recovered samples in the clipped regions to be consistent with the clipping function, where $\underline{1} \in \mathbb{R}^{m \times n}$ is an all-one matrix. As a result, instead of optimizing (17), we can optimize the following cost function

$$\hat{X} = \arg\min_{X} \left\{ \|A - \Omega X\|_{F}^{2} + \mu \left[\|M^{r}(Y - X)\|_{F}^{2} + \|M^{c+}(\theta^{+}\underline{1} - X)_{+}\|_{F}^{2} + \|M^{c-}(\theta^{-}\underline{1} - X)_{-}\|_{F}^{2} \right] \right\}$$
(18)

where a regularization term controlled by a parameter μ is used to approximate the constraint $X \in C(Y)$. A relaxed problem similar to the one proposed in (18) has been used in [4], [27], [40], [41] in the context of the synthesis model with a synthesis dictionary [42], [43], [44].

We then introduce an auxiliary function $Z(X) \in \mathbb{R}^{m \times n}$ defined as follows

$$Z(X) = M^{r}Y + M^{c+} \max\left(\theta^{+}\underline{1}, X\right) + M^{c-} \min\left(\theta^{-}\underline{1}, X\right)$$
(19)

With Z(X), we can simplify (18) to

$$\hat{X} = \underset{X}{\arg\min} \|A - \Omega X\|_{F}^{2} + \mu \|Z(X) - X\|_{F}^{2}$$
(20)

With a least squares based method [45], [46], [47], the solution of X can be obtained analytically from (20) as

$$X^{[t+1]} = \left(\Omega^T \Omega + \mu I\right)^{-1} \left(\Omega^T A + \mu Z(X^{[t]})\right)$$
(21)

where $I \in \mathbb{R}^{m \times m}$ is an identity matrix, A and Ω are obtained by following the dictionary update step in (16). In our work, Ω is real-valued, and $\Omega^T \Omega$ is a positive definite matrix. As a result $(\Omega^T \Omega + \mu I)$ is a full-rank invertible positive definite matrix. Empirically, we observed that the matrix inversion was stable in our experiments. Note that the use of the auxiliary function Z(X) is to facilitate the solution of (20). Although the three types of samples are regularized with a single μ in (18), there may be a potential scope to treat them differently with different regularization parameters.

The second method for optimising (17) is based on the projected gradient algorithms [47], [48]. The gradient of the objective function (17) with respect to X is

$$\nabla \|A - \Omega X\|_F^2 = \frac{\partial \|A - \Omega X\|_F^2}{\partial X}$$

= $-2\Omega^T A + 2\Omega^T \Omega X.$ (22)

Thus, $X^{[t+1]}$ can be obtained as

$$X^{[t+1]} = X^{[t]} - \lambda (-2\Omega^T A + 2\Omega^T \Omega X^{[t]})$$
(23)

where λ is a step size. A projection step is then used to enforce $X^{[t+1]} \in C(Y)$ as follows

$$X^{[t+1]} \leftarrow M^{r}Y + M^{c+} \max\left(\theta^{+}\underline{1}, X^{[t+1]}\right) + M^{c-} \min\left(\theta^{-}\underline{1}, X^{[t+1]}\right)$$
(24)

where the measurement matrices are the same as those in Equation (14), and \leftarrow means updating $X^{[t+1]}$ with the formula on the right hand side.

With the SAD algorithm, we can finally obtain the analysis dictionary Ω , the analysis representation matrix A and the estimated signal X. Algorithm 2 summarizes the implementation details of the SAD algorithm.

C. Algorithm Convergence and Initialization

The SAD algorithm alternates among the steps of analysis sparse coding, analysis dictionary update and signal estimation. The algorithms for estimating X, i.e. the operation 4.a in Algorithm 2 have been shown to be convergent in [20] and [49]. However, convergence of Step 3 cannot be ensured. We observed

Algorithm 2: Implementation Details of SAD.

Initialization: $\Omega^{[0]}, X^{[0]} = Y, t = 0, l, \mu, \lambda$ 1. While not converged do 2. **Input:** $X^{[t]}, \Omega^{[t]}, \overline{\Omega}_{[1]} = \Omega^{[t]}, k = 0, g(\Omega_{[0]}) = 0$ 3. a. do b. $\Omega_{[k+1]} = w(\Omega_{[k]})$ using equation (11) for function $w(\Omega)$ $A_{[k+1]} = HT_l(\Omega_{[k+1]}X^{[t]})$ c. using equation (6) for function $HT_l(\Omega X)$ d. k = k + 1until a pre-defined number of iterations is reached. e. f. **output:** $\Omega^{[t+1]} = \Omega_{[k+1]}, A^{[t+1]} = A_{[k+1]}, Z^{[t+1]}$ $\Omega = \Omega^{[t+1]}, A = A^{[t+1]}, Z^{[t+1]}$ 4. a. using equation (21) or (23) (with (24)) to get $X^{[t+1]}$ b. output: $X^{[t+1]}$ 5. t = t + 16. End while 7. **Output:** X

that $g(\Omega_{[k]})$ may increase occasionally, which could be caused by the update of X with (20) and (19) in which the comparison between X and $\theta^+ \mathbb{1}$ or $\theta^- \mathbb{1}$ is performed. However, the decreasing trend of $g(\Omega_{[k]})$ is often restored automatically with further iterations. In practice, we did not observe any other issues about the convergence of the algorithm. We have used a pre-defined number of iterations (e.g. 10) as the stopping criterion for the inner loop for updating the analysis dictionary. In Algorithm 2, we use the superscript t to denote the iteration index for the outer loop, while the subscript k for the inner loop.

In the SAD algorithm, the initial dictionaries Ω can be set as random matrices with normalized rows, or as pre-defined dictionaries such as the DCT dictionary. The initial dictionaries can affect the results, as observed in our experiments. We will examine the impact of the initial analysis dictionary on the declipping performance in Section V.

V. SIMULATIONS

In this section, we focus on demonstrating the performance of the proposed SAD algorithm using different parameters, on sound datasets. Our algorithms can be easily applied to other data, such as images. We use SDR as the performance metric, which is defined as follows

SDR
$$(\hat{x}, x) = 20 \log \frac{\|x\|_2}{\|x - \hat{x}\|_2}$$
 (25)

where \hat{x} is the estimated signal, and x is the original signal.

In the following subsections, we will show the setup of our experiments, the performance of the SAD algorithm, including the influence of the parameters m and l, and the initialization of Ω on the performance of the algorithm. The code of our proposed SAD algorithm can be downloaded from Github.¹ We

¹[Online]. Avaiable: http://github.com/BinLi504/SAD_matlabopen



Fig. 2. The waveform of the original speech signal and the clipped signal used for performance tests. The original signal was normalized to [-1,1], before being clipped at the level $\theta = 0.15$.

will present the comparisons of our algorithms with the baseline algorithms, i.e. ConsDL [19] and A-SPADE [23], in Section VI.

A. Experimental Setup

To form Y, we take the clipped signal and divide it into n segments, each of length m, with overlaps between the adjacent segments, using a sliding window. For example, if the clipped signal has 16 384 samples, we can form Y as a matrix 64×16321 , using a sliding window of m = 64 samples with a hop size equal to one sample. In the experiments in this section, we tested several different values for m, i.e. m = 16, 32, 64, and 128. The specific values used are discussed in the corresponding subsections.

In our experiments, the clipping level θ was set using a routine from [27], which takes an input SDR and outputs the clipping level. The clean signal was normalized before it was clipped at the specified level. More details about the generation of clipped signals are given in Section VI-A.

Our experiments were performed using Matlab on the Intel Core i5-7500 CPU, with memory 8 G-Bytes, on Windows 10 operating system.

B. Empirical Tests of Parameters

We first test the performance of the proposed algorithm (with the least-squares based update for X) using a speech signal "dev_male2_150ms_1_ch12.wav" from the SISEC dataset [50]. These signals contain stereo recordings, with 8 bits per channel, sampled at 16 kHz. One of the channels is used as test data. The original signal is relatively long, therefore we cut it into 16 384 samples, as shown in Fig. 2(a). We study empirically the set up of important parameters such as co-sparsity and dimension of dictionary atoms. Here we use a relatively short signal for parameter tuning for the proposed SAD method, however, we have also used longer and more test signals in Section VI for performance comparisons on speech and music datasets. The clipping level θ is set as 0.15. The waveform of the clipped signal is shown in Fig. 2(b). We set m = 64, l = 120, and the number of iterations as 20 000.

1) Selection of Co-Sparsity l and m: In the SAD algorithm, the observed signal y is divided into segments each of length m (i.e. equal to the dimension of dictionary atoms in terms of



Fig. 3. The SDR achieved using different co-sparsity l. The setting of l is related to the dimension of dictionary atoms m. For a given m, the choice l can be found empirically.

our formulations discussed in Section IV), and the estimated signal is finally obtained by concatenating the segments with an overlap-and-add technique. The dimension of dictionary atoms m and the selection of co-sparsity l affect the quality of the recovered signal. We study the impact of the co-sparsity l on the performance of the proposed algorithm. The value of l is varied, with a step increase of 8. The clipping level is fixed as $\theta = 0.15$. Fig. 3 shows the results with different *l*'s, when m = 128, 64, 32, and 16, respectively. It can be seen that the setting of l is related to m. For example, when m = 128, the best choice for l is 232, while for m = 64, the best choice for l is 120. When m = 32, the best choice for l is 56, while for m = 16, the best choice for l is 24. When we use different m's, l should be set accordingly. For each m value, we choose l empirically in terms of the reconstruction quality. The SDR results with different m's are presented in Fig. 4. It can be observed that the SDR increases with the increase of m, but such improvement is flattened when m is relatively large. Although only a single signal was used in this test, we observed empirically that the performance trend is similar when other signals were used. Thus, the parameters tuned in this way seem to be appropriate for subsequent experiments.

2) Signal Recovery: To illustrate the performance of the proposed algorithm, we show the plots of three segments of the recovered signal and the original signal in Fig. 5. Here m = 64 and l = 120. It can be seen that the recovered signal is similar to the original signal.

C. Dictionary Initialization: Random Versus DCT

With the speech signal used in the previous section, we compare the performance difference for the dictionaries initialized as a random Gaussian matrix and a uniform distribution matrix. We run the experiments ten times and observe the effect of these



Fig. 4. The SDR of the recovered signal, achieved with different dimension of dictionary atoms m. The signal reconstruction performance by the proposed algorithm increases with the increase of m, however, the performance increase tends to be saturating with larger m.

initial dictionaries on the signal recovery results. The estimate of X is based on the least-squares update. The average results of these tests are presented in Fig. 6. It can be seen that the results are similar. The standard deviation of the SDR results in the 10 tests with Gaussian matrix is 0.083 dB and the uniform matrix is 0.052 dB. It appears that the randomly initialised dictionaries with these two distributions give similar SDR results.

We also compare the randomly initialised analysis dictionary with the dictionary initialised as DCT. As the two random dictionaries tested above give similar results, we initialise the random matrix by generating the elements following a Gaussian distribution with zero mean and unit variance. The atoms in the initial dictionaries were normalized. In this experiment, we have increased the number of test signals to obtain more reliable results. We used 10 speech signals,² and the length of each signal is about 5 to 6 seconds, including about 80 000 to 96 000 samples (as the sampling rate is 16 kHz). The parameters of the SAD algorithm are set as m = 72 and l = 136, and X is solved with the least-squares update.

Fig. 7 shows the SDR of the recovered signal with the dictionary learning algorithm in which the dictionary is initialised as either a random Gaussian matrix or as a DCT dictionary. Overall, using the DCT dictionary performs similarly to using a Gaussian dictionary in the initialization. When the clipping level is at 0.1, the SDR obtained by using DCT is 0.3 dB higher than that using the Gaussian dictionary.

D. Least Squares Versus the Projected Gradient

In the SAD algorithm, two methods can be used to obtain X, i.e. the least squares based method and the projected gradient based method. In this section, we compare the performance of these two methods. We use the same experimental set up as in Section V-C, with the analysis dictionary initialised with the

²[Online]. Avaiable: http://www.repository.voxforge1.org/downloads/ SpeechCorpus/Trunk/ Audio/Main/



Fig. 5. Signal recovery examples: (a) (c) and (e) show three segments of an original clean speech signal and the clipping level. (b), (d) and (f) show the corresponding segments of the estimated (i.e. declipped) signal.



Fig. 6. The SDR of the declipped signal, achieved with the analysis dictionary initialized as a random matrix whose elements follow Gaussian or uniform distributions with zero mean and unit variance.

DCT dictionary. Fig. 8 shows that the SDRs obtained by the two methods are very similar.

We have tuned the parameters such as the step size λ , by running a set of experiments. In particular, we tested different combinations of the parameters $\theta = 0.2$, and 0.8, $\lambda = 0.1$, 0.25, and 0.4, and the number of iterations equal to 1, 10 and 20,



Fig. 7. The SDR of the declipped signal, achieved either with the dictionary initialized as a random matrix whose elements follow Gaussian distribution or with a DCT dictionary.



Fig. 8. The SDR of the declipped signal, obtained with the proposed SAD algorithm, where the least squares method and the projected gradient method are used for updating X, respectively.

respectively. The test results are given in Table I. The results show that the step size and the number of iterations can affect the SDR results of the recovered signals. Based on these tests, we chose the step size at 0.25 and the number of iterations at 10 in the subsequent experiments.

We also analyzed the impact of μ on the results by changing μ from 0.4 to 1.8, while keeping other parameters unchanged. The test results for the clipping level at 0.15 (i.e. input SDR at 5 dB) are given in Table II. From the table, we can see that the selection of μ has a very small influence on the results.

These two methods seem to offer very similar performance, and both need to tune a parameter (i.e. μ in the least squares based method, and λ in the projected gradient method). However, we found empirically that their computational complexities can be different. The projected gradient based method contains an additional loop (i.e. the inner loop for updating X), and if multiple iterations in the inner loop are performed, its complexity will

TABLE I TESTS FOR DIFFERENT COMBINATIONS OF PARAMETERS (CLIPPING LEVEL, STEP SIZES AND ITERATION NUMBERS) IN THE PROPOSED METHOD WITH PROJECTED GRADIENT

θ	λ	iterations	SDR (dB)		
0.2	0.1	1	15.36		
0.2	0.1	10	14.70		
0.2	0.1	20	14.76		
0.2	0.25	1	15.35		
0.2	0.25	10	14.54		
0.2	0.25	20	14.54		
0.2	0.4	1	13.47		
0.2	0.4	10	12.10		
0.2	0.4	20	12.01		
0.8	0.1	1	39.88		
0.8	0.1	10	39.90		
0.8	0.1	20	39.98		
0.8	0.25	1	39.94		
0.8	0.25	10	40.91		
0.8	0.25	20	41.12		
0.8	0.4	1	42.32		
0.8	0.4	10	41.20		
0.8	0.4	20	41.31		

 TABLE II

 The Influence of Different μ Values on Signal Recovery

 Performance in Terms of SDR of the Recovered Signal

μ	0.4	0.6	0.8	1.0	1.2	1.4	1.6	1.8
SDR	8.64	8.64	8.66	8.66	8.66	8.66	8.66	8.64

become higher than the least squares based method. For this reason, we will only use the least-squares based method for the comparisons in the next section.

VI. COMPARISON WITH BASELINE ALGORITHMS

In this section, we compare the performance of our proposed SAD algorithm with two recent baseline algorithms, i.e. the ConsDL algorithm [19] and the A-SPADE algorithm [23].

A. Datasets and Test Preparation

Both speech and music data were used in our evaluations. The speech signals were taken from the Trunk dataset. We used four folders from this dataset, which are respectively, AT-20 130 718-lws, Adminvox-05 232 006, 1snoke-20 120 412-hge, and 1337ad-20 170 321-ajg. In total, there are 50 speech signals from these four folders, sampled at 16 kHz. The length of each signal is about 4 to 7 seconds. The music signals were taken from the DSD100 dataset.³ The music files used are the vocals of 055 – Angels In Amplifiers - I'm Alright, vocals of 081 - Patrick Talbot - Set Me Free and vocals of 049 - Young Griffo - Facade, downloaded from DSD100 dataset. Each of these files contains sound recordings of approximately 45 to 49 seconds, sampled at 44.1 kHz. We then pre-processed the data, and generated 50 signals each of 1.9 seconds, i.e. 81 920 samples.

The clipped signals were generated in the same way for speech and music signals, by following [27]. More specifically, we set

³[Online]. Avaiable: https://www.loria.fr/ aliutkus/DSD100subset.zip

TABLE III MAPPING BETWEEN AVERAGE CLIPPING LEVEL θ and Input SDR of the Speech Signals (S/S), and Music Signals (S/M) Used in the Experiments

θ	0.1	0.2	0.3	0.4	0.5	0.6	0.7	0.8	0.9
S/S	4.2	7.5	11.3	15	19	23.5	29	36	45
S/M	3.5	7	10.8	14	18	22	27	33	42

the same input SDR for each of the 50 signals, and tune the level of clipping to match the input SDR. We also tune the input SDR so that the level of clippping is achieved approximately at the level specified. We then calculate the average level of clipping over these 50 signals. For example, when the input SDR is 4.2 dB for 50 speech signals, the average θ is approximately 0.1. We varied the clipping level θ from 0.1 to 0.9, which corresponds to the input SDR from 4.2 dB to 45 dB, as shown in the second row of Table III. Note that for the same input SDR, different signals may lead to slightly different θ values. Therefore, the θ is an average θ for all the clipped signals used in our tests, tuned to be approximately at the corresponding clipping level.

The implementation of the ConsDL algorithm [19] comes from the author's personal homepage.⁴ For the A-SPADE algorithm [23] we used its version 1.0, downloaded from the author's homepage.⁵ In our experiments, the default settings of the ConsDL (i.e. the size of overlapping time frames N = 256with rectangular windows and 75% overlap) and A-SPADE algorithms were used. The parameters used in the SAD algorithm are set as p = 144, m = 72 and l = 120, with initial analysis dictionary Ω set as a DCT dictionary matrix. The condition for stopping algorithm iterations is that the difference between the values of the cost function at two consecutive iterations is less than 0.001. The same experimental setup has been used for the experiments on speech and music data. We take the average of the SDRs calculated from the recovered signals from the 50 tests as the performance metric.

B. Results on Speech Signals

Fig. 9 shows the average SDR of all the recovered speech signals for each clipping level. From this figure, we can see that the proposed SAD algorithm performs better than other algorithms, except when $\theta = 0.1$, and 0.2, in which case, A-SPADE performs better. Fig. 10 shows the variance of the SDR results for all the tests on the 50 speech signals obtained by the SAD for each clipping level, as compared with the baseline algorithms. It can be observed that the ConsDL algorithm gives the smallest variance, while the variance of A-SPADE is higher than those of other compared methods including the proposed SAD algorithms, except when $\theta < 0.2$. It is interesting to note that the variance of all the algorithms increases with the increase in θ , which is probably not surprising considering the fact that the average SDR improvements also increase with the increase in θ .



Fig. 9. Average SDR over all the recovered speech signals obtained by the SAD algorithm as compared with A-SPADE and ConsDL, for different clipping levels.



Fig. 10. Variance of the SDRs for all the recovered speech signals obtained by the SAD as compared with the A-SPADE and ConsDL algorithms, for each clipping level.

C. Results on Music Signals

Fig. 11 shows the average SDR results of all the tests for the music signals. The proposed method performs better than ConsDL for all the clipping levels, and slightly better than A-SPADE, except for low clipping levels. Fig. 12 shows the variance of the SDR results of the recovered music signals, for each clipping level. From this figure, we can see that the variance of the SAD algorithm is smaller than that of the A-SPADE algorithm but larger than that of the ConsDL algorithm.

In the A-SPADE algorithm, we used the default parameter set up in its original codes, which was for signals sampled at 16 kHz, and may not be optimal for the experiments on the music signals sampled at 44.1 kHz. In practice, it may not be a trivial task to tune the parameters of A-SPADE to obtain the optimal performance for a different sampling rate. To address this issue, we run another set of experiments, where we first re-sample music signals from 44.1 kHz to 16 kHz. We then run the tests of all the compared algorithms on the re-sampled music signals. Fig. 13 shows the average SDR of the recovered 50 music signals obtained by the proposed SAD algorithm, as compared with A-SPADE and ConsDL. From this figure, we

⁴[Online]. Available: https://www.cvssp.org/Personal/LucasRencker/ software.html#DL_for _declipping

⁵[Online]. Available: http://spade.inria.fr



Fig. 11. Average SDR of all the recovered music signals obtained by the proposed SAD algorithm, as compared with the A-SPADE and ConsDL algorithms for different clipping levels. For comparison, the average input SDR is also shown for each clipping level.



Fig. 12. Variance of the SDR results for all the recovered music signals obtained by the SAD, as compared with the A-SPADE and ConsDL algorithms for each clipping level.



Fig. 13. Average of the SDR results for all the tests on the recovery of the down-sampled music signals, obtained by the SAD, as compared with the A-SPADE and ConsDL algorithms for different clipping levels. For comparison, the average input SDR is also shown for each clipping level.



Fig. 14. Variance of the SDR results for all the tests on the recovery of the down-sampled music signals, obtained by the SAD, as compared with the A-SPADE and ConsDL algorithms for different clipping levels.

can see that the SDR results obtained are quite similar to those in Fig. 11. This shows that the influence of sampling rate on the performance of the algorithms compared is very small. Fig. 14 shows the variance of the SDR results for all the tests on the re-sampled music signals obtained by the SAD for each clipping level, as compared with the baseline algorithms. From the figure, we can see that the general trend of performance is similar to those shown in Fig. 12, although it appears to be more stable with respect to different clipping levels. The ConsDL algorithm still gives the smallest variance. The variance of A-SPADE is higher than those of other compared methods including the proposed SAD algorithm.

D. A Note on Running Speed

We found empirically that the proposed SAD algorithm is computationally less efficient as compared with A-SPADE and ConsDL. To process a signal of 81 920 samples with a dictionary of dimension 144×64 , the proposed SAD algorithm using least-squares based update for X took about 18 minutes to run for 500 (outer) iterations, with the simulation environment we mentioned earlier. Using the projected gradient based method, the proposed SAD algorithm needs to run an additional loop (i.e. the inner loop for updating X). With the same number of iterations for the outer loop, i.e. 500, the SAD algorithm took about 17.5 and 26.5 minutes, respectively, when the inner loop was run for one iteration and 10 iterations, respectively. However, the A-SPADE and ConsDL algorithms took about 6 and 6.5 minutes, respectively.

VII. CONCLUSION

We have presented a sparse analysis-model based signal declipping algorithm. In our proposed SAD algorithm, the cost function is established on a sparse analysis model, and the clipped signal is restored by using the analysis dictionary learned directly from the observed signal, via the Analysis SimCO algorithm. The evaluation results showed that our proposed algorithm offers better performance in terms of SDR and more stable results in terms of variance, than two recent baselines, namely, the ConsDL and A-SPADE algorithms.

In our work, we considered only real-valued analysis dictionary Ω . The method could potentially be extended to complexvalued dictionaries, such as the dictionaries learned with complex non-negative matrix factorization [51], which uses another matrix to encode the phase term of the complex numbers. Including phase term could help maintain phase coherence between different frequency components within the signal [51].

In addition, we used a fixed step size in the projected gradient method, which could be replaced by a variable step size, and optimized using e.g. the golden section rule, as used in [36] and [37]. Other possible directions of future work include extension of the SAD algorithm for noisy signal declipping, the employment of the dictionary pair learning (DPL) [52] for declipping, and comparisons with other analysis dictionary learning algorithms such as [53].

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