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HIGH RESOLUTION SPARSE ESTIMATION OF EXPONENTIALLY DECAYING SIGNALS

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ABSTRACT

We consider the problem of sparse modeling of a signal consisting of an unknown number of exponentially decaying sinusoids. Since such signals are not sparse in an oversampled Fourier matrix, earlier approaches typically exploit large dictionary matrices that include not only a finely spaced frequency grid but also a grid over the considered damping factors. The resulting dictionary is often very large, resulting in a computationally cumbersome optimization problem. Here, we instead introduce a novel dictionary learning approach that iteratively refines the estimate of the candidate damping factor for each sinusoid, thus allowing for both a quite small dictionary and for arbitrary damping factors, not being restricted to a grid. The performance of the proposed method is illustrated using simulated data, clearly showing the improved performance as compared to previous techniques.

Index Terms—Sparse signal modeling, Spectral analysis, Sparse reconstruction, Parameter estimation.

1. INTRODUCTION

Exponentially decaying sinusoidal signals occur naturally in a wide range of fields, such as, for instance, radio frequency spectroscopy, wireless communications, sonar and radar (see, e.g., [1, 2] and the references therein). Commonly, the measurements suffer from various kinds of interference signals or corrupting additive colored noise. Furthermore, the number of modes present in the signal are generally unknown or may vary over time, typically necessitating some form of model order selection decision. Given such difficulties, it is often of interest to formulate non-parametric or semi-parametric modeling techniques, imposing only mild assumptions of the a priori knowledge of the signal structure. Popular solutions include the so-called dCapon, dAPES, and dIAA spectral estimators which form a generalized spectral estimate of the signal, constructing their spectral representation over both the frequency and damping dimensions [1, 2] (see also [3, 4]). Although this form of techniques are robust to the model assumptions, they suffer difficulties in separating closely spaced frequency and damping modes from each other, and typically require notable computational efforts if not implemented carefully [4]. As an alternative, one may use sparse modeling of the signal, forming a large dictionary of all potential frequencies and damping candidates, thus generally having vastly more columns than rows. For a given signal and the resulting dictionary matrix, one thus wishes to find the sparsest solution to the resulting linear set of equations mapping the signal to a linear combination of a few of the columns of the dictionary. Such techniques have successfully been applied to line spectral data, and the topic has attracted notable attention in the recent literature (see, e.g., [5–11]). Although these algorithms appear quite different from each other, they share the property that the considered dictionary grid should be selected sufficiently fine to allow for a sparse signal representation (see also [12, 13]), which, if extended to also consider damped modes, necessitates a large dictionary matrix containing elements with a sufficiently fine grid over the range of both the potential frequencies and damping candidates (see, e.g., [2, 14, 15]); this will be particularly noticeable if treating large data sets, or data sets with multiple measurement dimensions, such as in NMR measurements. In order to mitigate this problem, we here propose a novel dictionary learning approach wherein we iteratively decompose the signal with a fixed small dictionary, adaptively learning the dictionary elements best suited to enhance sparsity. To this effect, we initially form a coarsely spaced dictionary with undamped modes over the range of considered frequency candidates, iteratively adapting both the frequency and damping settings for the dictionary elements, thereby also allowing for both a reduction and an expansion of the number of dictionary elements considered in the optimization.

2. SPARSE DICTIONARY LEARNING

Let \( y(n) \) denote a signal of interest that may be well modelled as a sum of \( K \) exponentially decaying sinusoids, i.e.,

\[
y(n) = \sum_{k=1}^{K} \alpha_k e^{-2\pi j f_k n - \beta_k n} + e(n) \quad (1)
\]

for \( n = 1, \ldots, N \), where \( f_k, \beta_k, \text{ and } \alpha_k \) are the frequency, damping factor, and complex amplitude of the \( k \)-th component, respectively, whereas \( e(n) \) denotes an additive circularly

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symmetric Gaussian noise. Here, both the number of modes, $K$, and the parameters detailing each such mode, i.e., $f_k$, $\beta_k$, and $\alpha_k$, for $k = 1, \ldots, K$, are considered unknown, as is the variance of the additive noise, $\sigma_e^2$. Stacking the observations of the signal into a vector

$$y = [ y(1) \ldots y(N) ]^T$$

(2)

where $(\cdot)^T$ denotes the transpose, allows the signal to be expressed using a sparse signal representation, such that it may be formulated as a linear combination of only a few out of a range of potential candidate modes, i.e.,

$$y = \sum_{k=1}^P \alpha_k w_{f_k}(\beta_k) + e = W(\beta)a + e$$

(3)

where the noise vector $e$ is constructed similarly to $y$, and

$$W(\beta) = [ w_1(\beta_1) \ldots w_P(\beta_P) ]$$

(4)

$$w_{f_k}(\beta_k) = [ e^{2\pi f_k \beta_k} \ldots e^{2\pi f_k \beta_k N} ]^T$$

(5)

$$a = [ \alpha_1 \ldots \alpha_P ]^T$$

(6)

$$\beta = [ \beta_1 \ldots \beta_P ]^T$$

(7)

with $P$ denoting the total number of dictionary elements, typically formed such that $\{f_k\}_{k=1}^P$ are chosen uniformly over the frequency range of interest, e.g., $[0, 1]$, for a given set of damping factors, $\{\beta_1\}$, for $\ell = 1, \ldots, P$. To allow for high-resolution estimates without imposing any prior knowledge of the frequency locality of the modes, we here assumed that $P \gg N$, although to reduce the computational complexity in our implementation, we actually form the dictionary using only about $N$ elements, which combined with an iterative zooming, as detailed below, is more than sufficient to yield the desired resolution. As the damping factors are unknown, the dictionary elements need to be initiated to some reasonable estimate, for instance, setting all damping constants equal to zero. The unknown amplitudes and damping constants may then be found as the solution to the optimization problem

$$\min_{a, \beta} ||y - W(\beta)a||_2^2 + \rho(a)$$

(8)

where $\rho(a)$ is a sparsity promoting penalty function, on $a$, often selected as the $\ell_1$ norm [6, 16], leading to a convex optimization problem, for a fixed $\beta$, which will thus reliably converge to a global minimum without the need of a good initialization (see, e.g., [17]). We note that in many cases, even further sparsity is desirable, which may then be obtained by using other heuristics for approximating the function that counts the number of non-zero elements, e.g., by using a concave penalty function [18], so that larger non-zero elements are penalized (relatively) less than small non-zero elements, or by iteratively forming improved $\ell_0$ approximations [19], although it should be stressed that such approaches lead to more difficult, non-convex, optimization problems, commonly suffering from numerous local minima. Without loss of generality, we are in this work limiting our attention to using only an $\ell_1$ penalty, as the found solutions are deemed to be sufficiently sparse, but the algorithm could easily be modified to using other penalties as well. After forming the initial estimates of $a$ and $\beta$ using (8), we proceed to the dictionary learning step wherein both the frequency and damping for each candidate mode in the dictionary are updated appropriately, and where, if desired, the number of dictionary elements may be reduced or expanded as deemed appropriate. The Fourier transform of an exponentially damped sinusoid reveals that the power spectral density decays as

$$\frac{2\beta}{(2\pi \Delta f)^2 + \beta^2}$$

(9)

where $\Delta f$ is the distance from the frequency from frequency of the actual spectral peak and $\beta$ the damping factor. The initial estimates will thus not be sparse, although the found spectral peak will still be at the correct frequency. Similarly, for a coarse initial grid spacing, closely spaced components or off-grid frequencies will still result in peaks in neighboring grid-points. As a result, one would obtain a sparser solution if one estimated a suitable damping factor for each of the peaks found in the initial estimate, as well as refine the used frequency grid for the peak location to allow for off-grid frequencies fitting the data better. This is done by estimating the frequency and damping factor for each of the modes found in (8), restricting the frequencies to move only within the range of the neighboring frequency grid points, such that a dictionary element may not move away further in frequency than a few grid points, updating the dictionary element accordingly. The dictionary is then updated using these refined dictionary elements, whereafter the entire process is iterated. To clarify, in order to form the refined dictionary elements, consider the residual from the solution of (8), i.e.,

$$\varepsilon = y - W(\beta)a$$

(10)

Algorithm 1 Seperating frequency and damping algorithm

1: Initialize $\beta^0 = 0$, $f_k^0 = k/N$ for $k = 1, \ldots, N$
2: for $m = 1, \ldots$ do
3: $a^m = \arg\min_a ||y - W(\beta^{m-1})a||_2^2 + \rho(a)$
4: $\varepsilon^m = y - W(\beta^{m-1})a^m$
5: $P^m = \text{number of peaks of } a^m$
6: for $j = 1, \ldots, P^m$ do
7: $I_j = \text{index of peak } j$
8: $\Gamma_j = \text{indexes of peak } j$ and the adjacent elements.
9: $\{\hat{\beta}_j, \hat{f}_j\} = \arg\max_{\beta, f} \{ \varepsilon^*_T(\beta) \Pi_{\varepsilon}(\beta) \}$
10: end for
11: Update $W$ with $\beta^m$ and a finer grid around $f^m$.
12: end for
Fig. 1. The RMSE of the frequency estimation as a function of SNR. The signal contains three modes.

where \( \hat{a} \) denotes the vector minimizing (8) with \( \beta = 0 \). Let \( I_j \) denote the index of the \( j \)th peak in \( \hat{a} \), and form \( \Gamma_j \) from the adjacent indices on each side of \( I_j \), i.e.,

\[
\Gamma_j = \begin{bmatrix} I_{j-1} & I_j & I_{j+1} \end{bmatrix}^T
\] (11)

The estimates of the corresponding dictionary elements can thus be refined using the minimization

\[
C = \min_{f,\beta,a} \| \epsilon + W_\Gamma(\beta)\hat{a}_\Gamma - w_f(\beta)a \|_2
\]

\[
= \min_{f,\beta,a} \| \epsilon_\Gamma(\beta) - w_f(\beta)a \|_2
\] (13)

with

\[
\epsilon_\Gamma(\beta) \triangleq \epsilon + W_\Gamma(\beta)\hat{a}_\Gamma
\] (14)

where by subscribing a set of indices the corresponding elements or columns of a vector or matrix are inferred, respectively. Solving for the unknown amplitude using least squares, for a fixed frequency and damping constant, yields

\[
\hat{a} = \frac{w_f^T(\beta)\epsilon_\Gamma}{w_f^T(\beta)w_f(\beta)}
\] (15)

with \((\cdot)^*\) denoting the conjugate transpose, which, if inserted in (13), implies

\[
\left\{ \beta_j, f_j \right\} = \arg \max_{f,\beta} \left\{ \epsilon^*_\Gamma(\beta) \Pi_w \epsilon_\Gamma(\beta) \right\}
\] (16)

where

\[
\Pi_w = \frac{w_f(\beta)w_f^T(\beta)}{w_f^T(\beta)w_f(\beta)} = \psi_w w_f(\beta)w_f^T(\beta)
\] (17)

3. NUMERICAL EXAMPLES

We proceed to examine the performance of the proposed method using simulated data. We initially consider \( N = 128 \) samples of a signal containing three modes, where the frequency and damping constants are chosen uniformly over
be the frequency and damping grid for dCapon was selected to 175 Monte-Carlo simulations. In this example, the size of the frequency and damping estimates, here defined as noise. The figures show the root mean squared error (RMSE) where \( \theta \) is the total number of Monte-Carlo simulations, and the damping of mode \( m \) denotes the estimate of either the frequency or the damping of mode \( k \) and Monte-Carlo simulation \( m \), \( M \) is the total number of Monte-Carlo simulations, and \( K \) is the number of modes. These results have been obtained using 175 Monte-Carlo simulations. In this example, the size of the frequency and damping grid for dCapon was selected to be 6000 \( \times \) 6000, uniformly covering frequencies and damping factors in [0, 1] and [0, 0.025], respectively, and using the recommended filter length of \( N/4 \). The SEMA algorithm on the other hand use a dictionary containing only 128 elements, with 40 elements being inserted in the dictionary refinement step in the iteration. As is clear from the figures, the proposed SEMA algorithm yields notably better estimates, without requiring a large dictionary grid over both dimensions, thereby allow for a substantially faster implementation. It is also worth noting that the dCapon estimation errors are here larger than the smallest possible error that is attainable given the current grid size, implying that the grid size does not in itself limit the quality of the estimates.

Next, we examine the ability of the methods to resolve two closely spaced spectral lines. In this case, we consider a signal containing two sinusoidal components with frequencies, \( f_1 = 0.6417 \) and \( f_2 = 0.6456 \), with random damping constants being drawn uniformly from [0, 0.025]. Figure 3 illustrates the resulting frequency estimates as obtained from 5 Monte-Carlo simulations, using \( \text{SNR} = 20 \text{ dB} \), and \( \text{SNR} = 20 \). The figure shows the estimates obtained using SEMA, dCapon, dIAA, and for a Lasso method with a dictionary containing both frequencies and damping factors and exploiting a zooming similar to the one used in SEMA. Here, the frequency grid for dCapon and dIAA has been selected to only be formed on [0.63, 0.67], allowing the methods notable \textit{a priori} information on the frequency region of interest. The damping grid ranges over [0, 0.025] and has size 500 for all methods, except for the used Lasso method, where, due to complexity reasons, it is set to 10. As seen in the figure, both the proposed method and the Lasso method clearly manage to resolve the two peaks, whereas dCapon and dIAA, in most cases, are unable to find the correct peaks. In the figure, the red square indicates the region 1/(2N) around the true frequencies. Throughout these simulations, we have used \( \rho(\lambda) = \lambda ||\alpha||_1 \), there \( \lambda \) has been selected such as being the average of the dominant amplitudes and that of the remaining spectrum, where the former is found as the mean of the amplitude of the \( k_0 \) largest peaks of the periodogram, whereas the latter is computed as the mean of the remaining periodogram estimate. Here, \( k_0 \) denotes the number of considered peaks, being equivalent to the model order given to the competitive algorithms. It should be stressed that selecting \( \lambda \) in this way does not imply using \( k_0 \) as the assumed model order, only acting as a guideline for how this weighting might be selected.

4. REFERENCES


